



**BROWNFIELD SITE INVESTIGATION REPORT  
LACKAWANNA FOUNDRY  
LACKAWANNA, ERIE COUNTY, NEW YORK**

**CERCLIS ID No.: NYR000073064**

EPA Contract No.: 68-W-00-121  
W.A. No.: 002  
W.O. No.: 20103.001.001.1002.00  
Document Control No.: 20103.1002.52

15 June 2001

Prepared for:

**UNITED STATES ENVIRONMENTAL PROTECTION AGENCY**

Prepared by:

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Date 6/14/01

## **1.0 INTRODUCTION**

The U.S. Environmental Protection Agency (EPA) tasked the Roy F. Weston, Inc. Site Assessment Team (SAT) to conduct a Brownfield Site Investigation (BSI) at the Lackawanna Foundry site located in Lackawanna, Erie County, New York. The purpose of the BSI was to characterize any residual contamination in soil under a concrete pad on-site, investigate the potential for groundwater contamination, and assess risks posed by the contamination. The sample locations were based on a request/recommendation by the New York State Department of Environmental Conservation. The Comprehensive Environmental Response Compensation and Liability Information System Identification Number (CERCLIS ID No.) for the Lackawanna Foundry site is NYR000073064 (Ref. No. 1, p.1). The Lackawanna Foundry property is being considered for future use as a park to be operated by the City of Lackawanna (Ref. Nos. 1, p.3; 2, p. 5).

### **1.1 Site Location/Description**

The Lackawanna Foundry site is a 2-acre property located at 2 Elm Street in Lackawanna, Erie County, New York (Ref. Nos. 1, p. 1; 3, p. 1; 4, p.17). This site was the location of a foundry that produced specialty cast iron molds for over 60 years. The current owner of the site inherited the business from his father, who operated the foundry at this location for more than 30 years. The previous owner of the property is unknown at this time. Local residents have stated that foundry operations have steadily decreased over the past 20 years. The current owner ceased operations completely in 1997 (Ref. No. 3, p.4).

The site is located in a mixed residential and heavily industrialized area of Lackawanna (Ref. Nos. 1, pp. 1, 2; 3, p. 4). The site is bordered to the north by Smoke Creek; to the south by a multi-use city park, to the east by an active rail yard, and to the west by residential properties (Ref. Nos. 1, p. 2; 3, p.4; 4, p.19). Twelve residential properties are immediately adjacent to the western border of the site property, and approximately 100 residences are located within one-eighth mile of the site. The population within 0.5 mile is approximately 1,000 (Ref. No. 3, p.4). Figure 1 presents a Site Location Map.

Prior to the EPA Removal Action, which began in March 1999, the site consisted of several foundry buildings, approximately 100 drums in varying stages of deterioration located outside of the buildings, and an additional 200 drums, laboratory containers, compressed gas cylinders, and transformers located within the foundry buildings. Wastes on site included slag, foundry sand, binders, oils, resins, and other wastes (Ref. Nos. 1, p. 2; 3, p. 2). The EPA Removal Action was completed in September 2000, and the site now consists of an open grassy plot with a vegetated wetland area located in the north central portion of the property. (Ref. Nos. 1, p.3; 2, p.5; 4, p. 19). The site is currently well vegetated and is surrounded by a fence with a locked gate to control access (Ref. No. 2, p. 5; 4, p.19).

### **1.2 Previous Investigations/Regulatory History**

In 1997, the New York State Department of Environmental Conservation (NYSDEC) performed an

inspection of the site and required the property owner to conduct an inventory of all containers on site, and to collect samples from transformers and suspect areas of polychlorinated biphenyl (PCB) soil contamination (Ref. Nos. 1, p. 2; 3, p. 7). Analytical results for the sampling event indicated that the transformer fluid contained up to 86 percent PCBs and nearby soil contained up to 210 parts per million (ppm) of PCBs (Ref. No. 1, p. 2; 3, pp.6-8). The inventory provided by the property owner listed more than 100 drums located at the site, which were in various stages of deterioration.

NYSDEC requested the property owner to take actions to address the threats at the Lackawanna Foundry site; however, the property owner was financially unable to perform the cleanup (Ref. Nos. 1, p. 2; 3, pp. 7, 8).

At the request of NYSDEC, EPA performed a Removal Assessment of the Lackawanna Foundry property that was completed in March 1999 (Ref. Nos. 1, p. 2; 3, p. 2; 5, p.1). The entire site was unfenced and accessible to the public, and evidence of trespassing was observed. Approximately 100 drums were located outside the foundry buildings and another 200 drums, laboratory chemicals, compressed gas cylinders, and PCB transformers were located within the foundry buildings. In addition, the buildings containing the drums and transformers were in extremely poor condition and several access points to the buildings were observed (Ref. Nos. 1, p. 2; 3, p. 2).

On 19 March 1999, EPA was granted verbal authorization to conduct emergency stabilization actions at the Lackawanna Foundry site (Ref. No. 3, pp. 2, 3). EPA mobilized to the site on 22 March 1999 to begin stabilization procedures. EPA installed a fence along the perimeter of the site, inventoried and moved all containers of hazardous substances into the foundry buildings, and overpacked the leaking drums and containers as necessary. EPA then secured the foundry buildings (Ref. No. 1, p. 2; 3, p.3).

During the emergency stabilization operations, field characterization of drum contents revealed the presence of flammable, oxidizing, water reactive, and corrosive substances. Label information indicated that containers of paint-related materials had various concentrations of xylene, acetone, toluene, and naphtha (Ref. No. 3, p.3). Over 1,000 pounds of laboratory chemicals were inventoried and stabilized during the emergency action. Label information from these containers indicated that hazardous substances and wastes such as sulfuric acid, phosphoric acid, flammable alcohols and acetone, ammonium hydroxide, calcium carbide, and many unknown materials were present in these containers (Ref. No. 3, p.3). Eight compressed gas cylinders were collected from throughout the property and staged according to compatibility. Four cylinders contained liquified petroleum gas, one contained propane, one contained oxygen, one contained polymeric isocyanate, and one cylinder contained unknown substances (Ref. No. 3, p.3).

On 14 April 1999, the EPA Environmental Response Team (ERT) and the EPA Response Engineering and Analytical Contract team (Weston) conducted surface and subsurface soil sampling to determine the extent of contamination outside and inside the on-site foundry buildings (Ref. No. 1, p. 2; 3, p. 3; 6, p. 1). Metals detected in the soil at levels exceeding NYSDEC cleanup levels included arsenic, beryllium, cadmium, chromium, copper, iron, mercury, nickel, and zinc (Ref. No. 6, p. 3, Table 3). Arsenic contamination in soil adjacent to the site in the recreational area appeared to be associated with the former use of a baghouse dust collection system. PCBs were detected in

soil at levels up to 39 mg/kg (Ref. No. 6, p.3, Table 1). Semivolatile organic compounds (SVOCs) detected in the soil above NYSDEC cleanup levels included phenol, 2-methylphenol, benzo(a)anthracene, chrysene, benzo(b)fluoranthene, benzo(k)fluoranthene, indeno(1,2,3-cd)pyrene, and dibenz(a,h)anthracene. Phenol was the most frequently detected SVOC, and was detected at the highest concentrations of all SVOCs (Ref. No. 6, p. 3, Table 2).

An Action Memorandum, signed by EPA on 13 August 1999, authorized a Removal Action at the Lackawanna Foundry site to complete the removal actions started in April 1999. On 19 October 1999, the EPA Superfund Technical Assessment and Response Team (Weston) collected 34 soil samples south of the Lackawanna Foundry building, near the recreational area owned by the City of Lackawanna. The surface and subsurface soil samples were collected from predetermined grid locations and analyzed for arsenic to determine its extent (Ref. Nos. 7, p. 2; 8, p. 2, Table 1). The on-site buildings were demolished in November 1999 and excavation of contaminated soil was completed in April 2000 (Ref. No. 1, p. 3).

Hazardous wastes transported off site for disposal included approximately 680 tons of contaminated soils with less than 500 mg/kg PCBs, 900 tons of arsenic-contaminated soils, 160 cubic yards of PCB-contaminated debris, 360 cubic yards of contaminated soil with greater than 500 mg/kg PCBs, 8 PCB-contaminated transformers, 153 PCB capacitors, 2 drums of PCB oils, and 68 drums containing various hazardous substances (Ref. No. 1, p. 4).

Site restoration actions, including the placement of approximately two feet of clean topsoil over the excavated areas, were completed in June 2000. The site was seeded with grass, and the wetland in the center portion of the site was restored by the use of clean marsh sediment from a New York State Parks dredging project (Ref. Nos. 1, p 3; 2, p.4; 4, p.18.).

## **2.0 BROWNFIELD INVESTIGATION**

The Brownfield Site Investigation at the Lackawanna Foundry site consisted of a file review of available background and analytical information, a site reconnaissance to determine soil sample and monitoring well locations, subsurface soil sampling, monitoring well installation, and collection of groundwater samples from the monitoring wells (Ref. No. 2, pp. 1, 4; 4, pp.17-21).

### **2.1 Site Reconnaissance**

On 4 October 2000, EPA and EPA Site Assessment Team (Weston) personnel conducted a site reconnaissance at the Lackawanna Foundry site. The purpose of this visit was to determine the locations for monitoring well installation and subsurface soil sample collection (Ref. No. 2, p. 5; 4, pp.17-21). During the site reconnaissance, no buildings or other structures were observed on the Lackawanna Foundry property. The relatively flat site was well vegetated and surrounded by a fence with a locked gate to control access. All known waste sources were removed from the property during the EPA Removal Action conducted earlier in the year. There were no signs of stressed vegetation or contaminated soil (Ref. No. 2, p. 5; 4, pp.17-21).

## 2.2 Subsurface Soil Sampling Analytical Results

On 7 November 2000, the EPA Site Assessment Team (Weston) collected 14 subsurface soil samples from six soil boring locations at the Lackawanna Foundry property. Soil samples SS-01, SS-02, SS-03, and SS-07 (duplicate of SS-03) were collected from beneath a buried concrete pad located on the eastern portion of the site. Soil samples SS-04 (A through D), SS-05 (A through C), and SS-06 (A through C) were collected from boring locations at depths of 2 to 10 feet or until groundwater was encountered. Figure 2, Sample Location Map, depicts the soil sampling locations. The soil samples were analyzed for Target Compound List (TCL) organic and Target Analyte List (TAL) inorganic parameters according to EPA Contract Laboratory Program (CLP) protocol (Ref. No. 10, pp. 1, 2, 5, 6; 11, pp.1-4).

Table 1A provides the organic analytical results for the subsurface soil samples collected at the Lackawanna Foundry site. No volatile organic compounds (VOCs) were detected in the soil samples at concentrations exceeding the NYSDEC Technical and Administrative Guidance Memorandum (TAGM) Recommended Soil Cleanup Objectives (RSCO) (Ref. Nos. 12, pp. 1, 2; 13, pp. 83-127).

Five semivolatile organic compounds (SVOCs) were detected at levels exceeding their respective RSCOs. Phenol was detected above its RSCO (30 parts per billion [ppb]) in two samples collected from one borehole location (SS-04). Sample SS-04A (2-4 feet), contained 7,900 ppb of phenol, and sample SS-04B (4-6 feet) contained 3,500 ppb of phenol. Chrysene was detected above its RSCO (400 ppb) in one sample, SS-04B, at 2,800 ppb. Benzo(b)fluoranthene was detected above its RSCO (1,100 ppb) in one sample, SS-04B, at 2,100 ppb. Benzo(a)pyrene was detected in three samples above its RSCO (61 ppb): sample SS-01 (7-9 feet) at an estimated concentration of 170 J ppb, sample SS-02 (6-8 feet) at 230 J ppb, and sample SS-06A (2-4 feet) at 190 J ppb. Dibenz(a,h)anthracene was detected above its RSCO (14 ppb) in one sample, SS-04B, at 530 ppb (Ref. Nos. 12, pp. 3-5; 13, pp. 128-172). No pesticides or PCBs were detected in the soil samples at concentrations exceeding their respective RSCOs (Ref. Nos. 12, pp. 1-9; 13, pp. 173-187).

Table 1B provides the inorganic analytical results for the subsurface soil samples collected at the Lackawanna Foundry site. RSCOs for inorganic hazardous substances are based on site background concentrations, or the higher of site background concentrations and a given numerical value. Since background concentration data are not available, sampling results were compared to the given numerical value (where applicable). Eight inorganic hazardous substances (metals) were detected at concentrations levels exceeding their respective RSCO. Arsenic was detected above its RSCO (7.5 ppm) in two samples, SS-04D (8-10 feet) at 16.9 ppm and SS-06C (6-8 feet) at 8.0 ppm. Chromium was detected above its RSCO (10 ppm) in 12 samples, ranging from 11.5 ppm in sample SS-01 to 27.2 ppm in sample SS-07 (duplicate of sample SS-03). Copper was detected above its RSCO (25 ppm) in three samples, SS-04D at 41.9J ppm, SS-05B (4-6 feet) at 25.1 ppm, and SS-06A (2-4 feet) at 34.4J ppm. Iron was detected above its RSCO (2,000 ppm) in 14 samples, ranging from 7,390 ppm in sample SS-04A to 52,800 ppm in sample SS-04D. Mercury was detected above its RSCO (0.10 ppm) in one sample, SS-04B, at 0.14J ppm. Nickel was detected above its RSCO (13 ppm) in 12 samples, ranging from 27.0 ppm in sample SS-01 to 48.9 ppm in sample SS-04D. Selenium was detected above its RSCO (2 ppm) in two samples, SS-06B at 2.3 ppm and SS-06C at 2.2 ppm. Zinc was detected above its RSCO (20 ppm) in 14 samples, ranging from 23.8 ppm in

sample SS-04A to 139 ppm in sample SS-06A. Beryllium was detected at concentrations exceeding its RSCO (0.16 ppm) in 12 samples, ranging in concentration from 0.45B ppm in sample SS-01 to 1.0B ppm in sample SS-04D (Ref. Nos. 12, pp. 1-9; 14, pp. 38-51).

### **2.3 Monitoring Well Installation/Groundwater Sampling**

One bedrock groundwater monitoring well and four overburden monitoring wells were installed on the Lackawanna Foundry property during 8-13 November. The bedrock well, GW-01, was installed to a depth of 67 feet (Ref. No. 11, p.10). The four overburden wells, GW-02 to GW-05, were drilled to a depth of 20 feet each (Ref. No. 11, pp.4-13).

WESTON SAT collected five groundwater samples (GW-01, GW-03, GW-04, GW-05, and GW-06[duplicate of GW-01]) from four monitoring wells on 5 December 2000 (Ref. No. 11, pp.14-16).

Figure 2, Sample Location Map, depicts the monitoring well/groundwater sample locations. The samples were analyzed for TCL organic and TAL inorganic parameters according to EPA CLP protocols. Monitoring well GW-02 was not sampled due to its inability to recharge (Ref. No. 11, p.16)

Table 2 provides the organic and inorganic analytical results for the groundwater sampling event conducted at the Lackawanna Foundry site (Ref. No.15, pp. 1, 6, 8). Results were compared to NYSDEC Ambient Water Quality Standards and Guidance Value (AWQS/GV). No organic compounds were detected at concentrations exceeding their respective AWQS/GVs (Ref. No. 16, Table 1; 17, pp. 2-28). Four inorganic concentrations exceeded their respective AWQS/GVs. Iron exceeded its AWQS/GV (300 ppb) in all five groundwater samples, ranging from 6,200 ppb in sample GW-03 to 42,000 ppb in sample GW-04. Magnesium exceeded its AWQS/GV (35,000 ppb) in one groundwater sample and its duplicate, sample GW-01 at 130,000 ppb and sample GW-06 at 120,000 ppb. Manganese exceeded its AWQS/GV (300 ppb) in three samples: GW-03 at 1,700 ppb, GW-04 at 1,700 ppb, and GW-05 at 2,200 ppb. Sodium exceeded its AWQS/GV (20,000 ppb) in four samples: GW-01 at 540,000 ppb, GW-03 at 22,000 ppb, GW-04 at 25,000 ppb, and GW-06 (duplicate of GW-01) at 550,000 ppb (Ref. No. 16, Table 1; 17, pp. 5, 9, 10, 14, 18, 22, 23, 26-30).

### **3.0 CONCLUSIONS AND RECOMMENDATIONS**

Analytical results for subsurface soil samples collected during the EPA BSI indicate the presence of five SVOCs (phenol and four polynuclear aromatic hydrocarbons, or PAHs) at concentrations exceeding their corresponding NYSDEC RSCOs. Phenols are associated with foundry sands and therefore are attributed to site operations. Eight metals were also found at concentrations exceeding RSCOs. No site background concentration data exist for metals to establish site-specific RSCOs for comparison to the analytical results. However, with the possible exception of iron, metals concentrations were generally found to be comparable across the site, and within expected background ranges. Exceedences of RSCOs were observed in soil samples collected under the concrete pad and from monitoring well pilot borings.

Iron, magnesium, manganese, and sodium exceeded their AWQS/GVs in groundwater. Magnesium

is a common component of subsurface materials in the region, and iron and manganese concentrations were lower in the downgradient monitoring well than in upgradient monitoring wells.

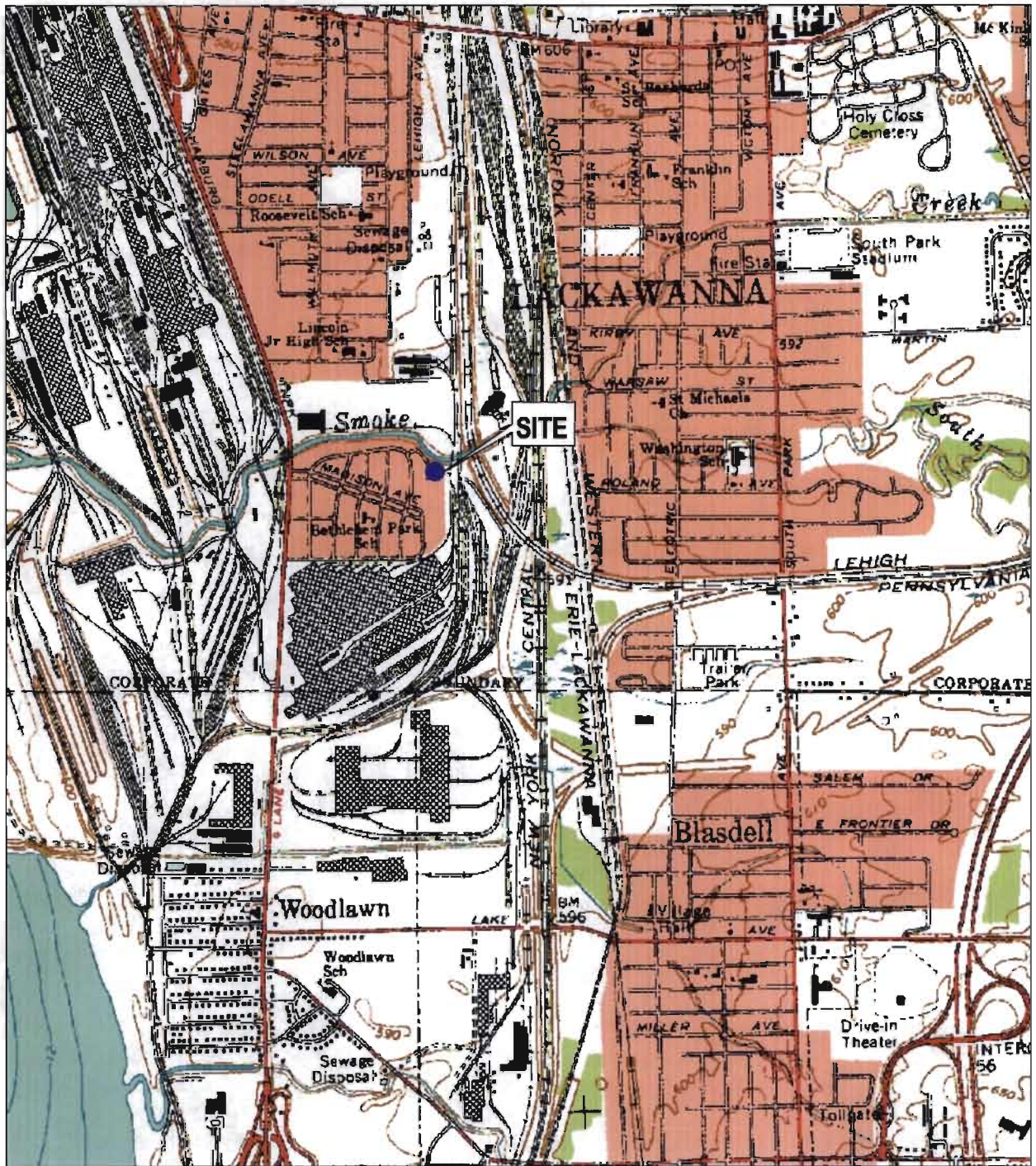
Based on the above, contamination attributable to former site operations exists in soil at depths below that which had been previously removed. However, a two-foot-thick cap of clean fill was emplaced at the site. Previous site operations do not appear to have impacted groundwater. Based on the response/removal actions conducted to date and the existing analytical data, the site appears to be suitable for nonresidential use. The current soil cap must be maintained to prevent exposure to subsurface contamination, and a deed restriction should be executed to control future disturbances of subsurface soil. If the concrete pad is removed from the site enough soil should be removed from beneath it to achieve a total excavation depth of two feet. The excavation should be backfilled with two feet of certified clean fill. Concurrence regarding the above should be obtained from the NYSDEC by the entity ultimately responsible for control, redevelopment, and/or future use of the property.



## REFERENCES

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# Figures

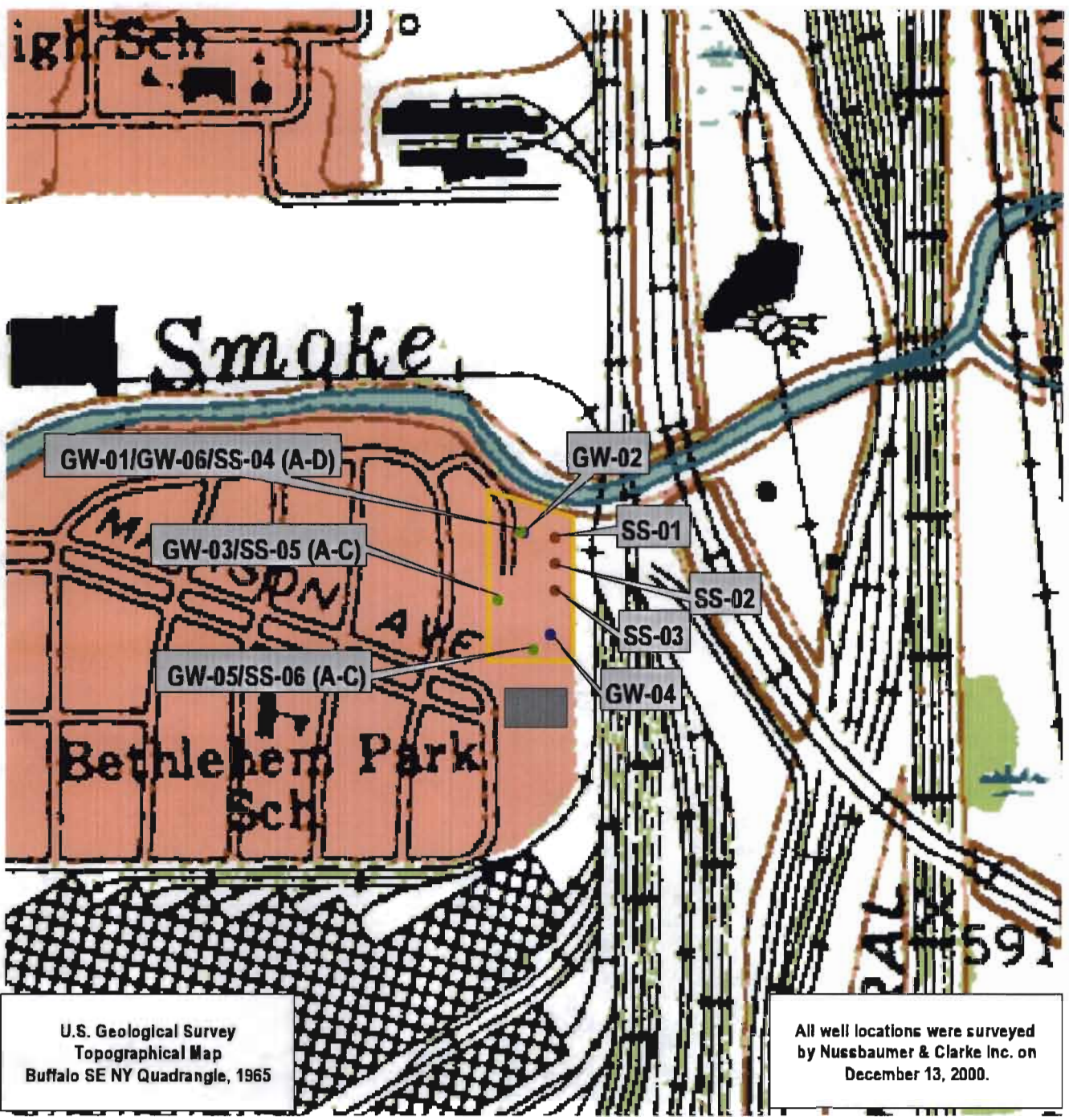


U.S. Geological Survey  
 7.5 Minute Series Topographical Map  
 Buffalo Southeast, NY Quadrangle  
 1948, Revised 1965



Lackawanna Foundry Site  
 Lackawanna, Erie Co., NY  
 Figure 1  
 Site Location Map

● Site Location



U.S. Geological Survey  
Topographical Map  
Buffalo SE NY Quadrangle, 1965

All well locations were surveyed  
by Nussbaumer & Clarke Inc. on  
December 13, 2000.



- Groundwater and Soil Sample Location
- Groundwater Sample Location
- Soil Sample Location
- City of Lackawanna Building
- Approximate Site Boundary

- A = Depth of 2-4 feet
- B = Depth of 4-6 feet
- C = Depth of 6-8 feet
- D = Depth of 8-10 feet



Lackawanna Foundry Site  
Figure 2  
Sample Location Map



# Tables

**TABLE 1A**  
**SOIL SAMPLE ORGANIC ANALYSIS RESULTS FOR**  
**EPA SAT BROWNFIELD SITE INVESTIGATION (NOVEMBER 2000)**  
**LACKAWANNA FOUNDRY, LACKAWANNA, NEW YORK**

| Field Sample No.<br>CLP Organic Sample No.<br>Time Sampled<br>Hazardous Substance | SS-01<br>BNP05<br>10:05 | SS-02<br>BNP06<br>10:55 | SS-03<br>BNP07<br>11:25 | SS-07<br>BNQ91<br>11:35<br>Dup. Of SS-03 | SS-04A<br>BNP08<br>14:40 | SS-04B<br>BNP09<br>14:50 | SS-04C<br>BNP10<br>14:55 | NYSDEC<br>Recommended Soil<br>Cleanup Objective |
|---|-------------------------|-------------------------|-------------------------|--|--------------------------|--------------------------|--------------------------|---|
| <b>VOCs</b>   |                         |                         |                         |  |                          |                          |                          |   |
| Carbon disulfide  | 2 J                     | ---                     | ---                     | ---                                      | ---                      | ---                      | ---                      | 2700  |
| 2-Butanone  | ---                     | 29                      | ---                     | ---                                      | ---                      | ---                      | ---                      | 300   |
| <b>SVOCs</b>  |                         |                         |                         |  |                          |                          |                          |   |
| Phenol  | ---                     | ---                     | ---                     | ---                                      | 7,900*                   | 3,500*                   | ---                      | 30  |
| Acetophenone  | ---                     | ---                     | ---                     | ---                                      | 830                      | ---                      | ---                      | NA  |
| Naphthalene   | ---                     | ---                     | ---                     | ---                                      | 1,100                    | 380                      | ---                      | 13000   |
| 2-Methylnaphthalene   | ---                     | ---                     | ---                     | ---                                      | 130 J                    | 150 J                    | ---                      | 36400   |
| Acenaphthene  | ---                     | ---                     | ---                     | ---                                      | ---                      | 300 J                    | ---                      | 50000   |
| Phenanthrene  | ---                     | ---                     | ---                     | ---                                      | ---                      | 3,700*                   | ---                      | 50000   |
| Anthracene  | ---                     | ---                     | ---                     | ---                                      | ---                      | 600                      | ---                      | 50000   |
| Di-n-butylphthalate   | ---                     | ---                     | ---                     | ---                                      | 58 J                     | 61 J                     | ---                      | 8100  |
| Fluoranthene  | ---                     | ---                     | ---                     | ---                                      | 67 J                     | ---                      | ---                      | 50000   |
| Pyrene  | ---                     | ---                     | ---                     | ---                                      | 59 J                     | ---                      | ---                      | 50000   |
| Benzo(a)anthracene  | ---                     | ---                     | ---                     | ---                                      | ---                      | ---                      | ---                      | 224   |
| Chrysene  | ---                     | ---                     | ---                     | ---                                      | 41 J                     | 2,800                    | ---                      | 400   |

Notes:  
All results are recorded in units of ug/kg.  
Dashed (--) lines indicate hazardous substances not detected.  
\*-Value reported from dilution run.  
NA-None available.

J-Estimated Value  
JN-Presumptive evidence of the compound at an estimated value.  
Shading indicates exceedence of NYSDEC Recommended Soil  
Cleanup Objective.

TABLE 1A (CONTINUED)

SOIL SAMPLE ORGANIC ANALYSIS RESULTS FOR  
EPA SAT BROWNFIELD SITE INVESTIGATION (NOVEMBER 2000)  
LACKAWANNA FOUNDRY, LACKAWANNA, NEW YORK

| Field Sample No.<br>CLP Organic Sample No.<br>Time Sampled<br>Hazardous Substance | SS-01<br>BNP05<br>10:05 | SS-02<br>BNP06<br>10:55 | SS-03<br>BNP07<br>11:25 | SS-07<br>BNQ91<br>11:35<br>Dup. Of SS-03 | SS-04A<br>BNP08<br>14:40 | SS-04B<br>BNP09<br>14:50 | SS-04C<br>BNP10<br>14:55 | NYSDEC<br>Recommended Soil<br>Cleanup Objective |
|---|-------------------------|-------------------------|-------------------------|--|--------------------------|--------------------------|--------------------------|---|
| <b>SVOCs (cont'd)</b>   |                         |                         |                         |  |                          |                          |                          |   |
| bis(2-Ethylhexyl)phthalate  | ---                     | 60 J                    | 380 J                   | 83 J                                     | 520                      | 870                      | 94 J                     | 50000   |
| Benzo(b)fluoranthene  | ---                     | ---                     | ---                     | ---                                      | ---                      | 2,100*                   | ---                      | 1100  |
| Benzo(k)fluoranthene  | ---                     | ---                     | ---                     | ---                                      | ---                      | 760                      | ---                      | 1100  |
| Benzo(a)pyrene  | 170 J                   | 230 J                   | ---                     | ---                                      | ---                      | ---                      | ---                      | 61  |
| Indeno(1,2,3-cd)pyrene  | ---                     | ---                     | ---                     | ---                                      | ---                      | 1,300                    | ---                      | 3200  |
| Dibenz(a,h)anthracene   | ---                     | ---                     | ---                     | ---                                      | ---                      | 530                      | ---                      | 14  |
| Benzo(g,h,i)perylene  | ---                     | ---                     | ---                     | ---                                      | ---                      | 1,300                    | ---                      | 50000   |
| <b>PESTICIDES</b>   |                         |                         |                         |  |                          |                          |                          |   |
| gamma-Chlordane   | 0.27 J                  | ---                     | ---                     | ---                                      | ---                      | ---                      | ---                      | 540   |
| alpha-BHC   | ---                     | ---                     | 0.38 J                  | ---                                      | ---                      | ---                      | 0.34 J                   | 111   |
| Heptachlor epoxide  | ---                     | 4.6                     | 0.77 J                  | ---                                      | ---                      | ---                      | ---                      | 20  |
| Dieldrin  | 0.68 J                  | ---                     | ---                     | ---                                      | ---                      | ---                      | ---                      | 44  |
| Endosulfan II   | ---                     | 2.5 J                   | ---                     | ---                                      | ---                      | ---                      | ---                      | 900   |
| <b>PCBs</b>   |                         |                         |                         |  |                          |                          |                          |   |
| Aroclor-1248  | ---                     | ---                     | ---                     | ---                                      | 170                      | 210                      | ---                      | 1000(surface), 10000(subsurface)                |

Notes:

All results are recorded in units of ug/kg.

Dashed (--) lines indicate hazardous substances not detected.

\*-Value reported from dilution run.

NA-None available.

J-Estimated Value

JN-Presumptive evidence of the compound at an estimated value.

Shading indicates exceedence of NYSDEC Recommended Soil

Cleanup Objective.

TABLE 1B

SOIL SAMPLE INORGANIC ANALYSIS RESULTS  
FOR EPA SAT BROWNFIELD SITE INVESTIGATION (NOVEMBER 2000)  
LACKAWANNA FOUNDRY, LACKAWANNA NEW YORK

| Field Sample No.<br>CLP Sample No.<br>Time Sampled<br>Hazardous Substance | SS-01<br>MBNJ51<br>10:05 | SS-02<br>MBNJ52<br>10:55 | SS-03<br>MBNJ53<br>11:25 | SS-07<br>MBNJ66<br>11:35<br>Dup. Of SS-03 | SS-04A<br>MBNJ54<br>14:40 | SS-04B<br>MBNJ55<br>14:50 | SS-04C<br>MBNJ56<br>14:55 | NYSDEC<br>Recommended<br>Soil Cleanup<br>Objective |
|---|--------------------------|--------------------------|--------------------------|---|---------------------------|---------------------------|---------------------------|--|
| Aluminum  | 6,770                    | 8,590                    | 15,900                   | 18,500                                    | 686                       | 4,170                     | 12,300                    | SB   |
| Antimony  | ---                      | ---                      | ---                      | ---                                       | ---                       | ---                       | ---                       | SB   |
| Arsenic   | 2.9 B                    | ---                      | 4.1                      | 2.2 B                                     | ---                       | 1.8 B                     | 4.2                       | 7.5 or SB  |
| Barium  | 67.3                     | 97.3                     | 187                      | 200                                       | 9.4 B                     | 35.1 B                    | 110                       | 300 or SB  |
| Beryllium   | 0.45 B                   | 0.72 B                   | 0.73 B                   | 0.76 B                                    | ---                       | ---                       | 0.66 B                    | 0.16 or SB   |
| Cadmium   | ---                      | ---                      | ---                      | ---                                       | ---                       | ---                       | ---                       | 1 or SB  |
| Calcium   | 2,140                    | 3,220                    | 2,640                    | 2,940                                     | 936 B                     | 60,300                    | 1,900                     | SB   |
| Chromium  | 11.5                     | 14.7                     | 26.2                     | 27.2                                      | 6.2                       | 8                         | 18.9                      | 10 or SB   |
| Cobalt  | 9.1 B                    | 9.7 B                    | 20.3                     | 14.6                                      | 0.99 B                    | 2.4 B                     | 10.1 B                    | 30 or SB   |
| Copper  | 19.7 J                   | 20.0 J                   | 21.0 J                   | 22.7 J                                    | 12.5 J                    | 14.3 J                    | 21.3 J                    | 25 or SB   |
| Iron  | 12,100                   | 16,200                   | 30,100                   | 27,700                                    | 7,390                     | 9,500                     | 24,800                    | 2000 or SB   |
| Lead  | 12.6                     | 15.3                     | 21.6                     | 18.1                                      | 15.2                      | 24.6                      | 18.4                      | SB   |
| Magnesium   | 2,630                    | 3,390                    | 5,510                    | 5,910                                     | 244 B                     | 2,490                     | 4,430                     | SB   |
| Manganese   | 117                      | 138                      | 1,050                    | 754                                       | 56.1                      | 145                       | 217                       | SB   |
| Mercury   | ---                      | ---                      | ---                      | ---                                       | 0.09 J                    | 0.14 J                    | ---                       | 0.1  |
| Nickel  | 27                       | 27.3                     | 48.7                     | 47.4                                      | 3.6 B                     | 7.7 B                     | 34.5                      | 13 or SB   |
| Potassium   | 639 B                    | 938 B                    | 1,080 B                  | 1,630 J                                   | 90.1 B                    | 425 B                     | 847 B                     | SB   |
| Selenium  | 1.3 B                    | ---                      | ---                      | ---                                       | ---                       | ---                       | ---                       | 2 or SB  |
| Silver  | ---                      | ---                      | 0.88 B                   | 0.63 B                                    | ---                       | ---                       | 0.53 B                    | SB   |
| Sodium  | 161 B                    | ---                      | ---                      | ---                                       | ---                       | 88.4 B                    | 228 B                     | SB   |
| Thallium  | ---                      | ---                      | ---                      | 0.94 B                                    | ---                       | ---                       | ---                       | SB   |
| Vanadium  | 12.8 B                   | 14.4 B                   | 24                       | 23.6                                      | 3.0 B                     | 6.0 B                     | 19.7                      | 150 or SB  |
| Zinc  | 71.4                     | 72.1                     | 128                      | 136                                       | 23.8                      | 46.4                      | 91                        | 20 or SB   |
| Cyanide   | ---                      | ---                      | ---                      | ---                                       | ---                       | ---                       | ---                       | NA   |

Notes:

All results are reported in units of mg/kg.

Dashed lines (--) denote hazardous substance not detected.

Shading denotes concentration exceeding NYSDEC Recommended Soil Cleanup Objective.

J-Estimated value.

B-Indicates concentration between the instrument detection limit and the contract required detection limit

SB-site background

NA-None available



TABLE 1A (CONTINUED)

SOIL SAMPLE ORGANIC ANALYSIS RESULTS FOR  
 EPA SAT BROWNFIELD SITE INVESTIGATION (NOVEMBER 2000)  
 LACKAWANNA FOUNDRY, LACKAWANNA, NEW YORK

| Field Sample No.<br>CLP Organic Sample No.<br>Time Sampled<br>Hazardous Substance | SS-04D<br>BNP11<br>15:15 | SS-05A<br>BNP12<br>13:25 | 22-05B<br>BNP13<br>13:40 | 22-05C<br>BNQ85<br>13:45 | SS-06A<br>BNQ87<br>13:00 | SS-06B<br>BNQ88<br>13:05 | SS06C<br>BNQ89<br>13:10 | NYSDEC<br>Recommended Soil<br>Cleanup Objective |
|---|--------------------------|--------------------------|--------------------------|--------------------------|--------------------------|--------------------------|-------------------------|---|
| VOCs  |                          |                          |                          |                          |                          |                          |                         |   |
| 2-Butanone  | ---                      | 9 J                      | 13                       | 8 J                      | 5 J                      | ---                      | ---                     | 300   |
| SVOCs   |                          |                          |                          |                          |                          |                          |                         |   |
| Phenanthrene  | ---                      | ---                      | 150 J                    | ---                      | ---                      | ---                      | ---                     | 50000   |
| Pyrene  | ---                      | 130 J                    | ---                      | ---                      | ---                      | ---                      | ---                     | 50000   |
| bis(2-<br>ethylhexyl)phthalate  | 240 J                    | 59 J                     | ---                      | 78 J                     | ---                      | 100 J                    | 81 J                    | 50000   |
| Di-n-octylphthalate   | ---                      | 120 J                    | ---                      | ---                      | ---                      | ---                      | ---                     | 50000   |
| Benzo(a)pyrene  | ---                      | ---                      | ---                      | ---                      | 190 J                    | ---                      | ---                     | 61  |
| PESTICIDES  |                          |                          |                          |                          |                          |                          |                         |   |
| Heptachlor  | ---                      | ---                      | 0.32 J                   | ---                      | ---                      | ---                      | ---                     | 100   |
| Heptachlor epoxide  | ---                      | 0.79 J                   | ---                      | ---                      | ---                      | ---                      | ---                     | 20  |
| Endosulfan I  | ---                      | 0.64 J                   | 0.26 J                   | ---                      | ---                      | ---                      | ---                     | 900   |
| Endrin  | ---                      | 2.3 J                    | ---                      | ---                      | ---                      | ---                      | ---                     | 100   |
| Endosulfan II   | ---                      | ---                      | ---                      | 0.14 J                   | ---                      | ---                      | ---                     | 900   |
| 4,4'-DDD  | ---                      | 1.8 J                    | 0.97 J                   | ---                      | ---                      | ---                      | ---                     | 2900  |
| Endosulfan sulfate  | ---                      | 6.1                      | ---                      | ---                      | ---                      | 0.13 J                   | ---                     | 1000  |
| Methoxychlor  | ---                      | 9.1 J                    | 7.0 J                    | ---                      | ---                      | ---                      | ---                     | Total VOCs < 10000                              |
| Endrin ketone   | ---                      | 4.1 J                    | ---                      | ---                      | ---                      | ---                      | ---                     | NA  |
| Endrin aldehyde   | ---                      | 7.7 JN                   | 5.0 J                    | ---                      | ---                      | ---                      | ---                     | NA  |
| gamma-Chlordane   | ---                      | 0.36 J                   | 0.35 J                   | ---                      | ---                      | ---                      | ---                     | 540   |

Notes:

All results are recorded in units of ug/kg.

Dashed (---) lines indicate hazardous substances not detected.

\*-Value reported from dilution run.

NA-None available.

J-Estimated Value

JN-Presumptive evidence of the compound at an estimated value.

Shading indicates exceedence of NYSDEC Recommended Soil

Cleanup Objective.

TABLE 1B (Continued)

SOIL SAMPLE INORGANIC ANALYSIS RESULTS  
FOR EPA SAT BROWNFIELD SITE INVESTIGATION (NOVEMBER 2000)  
LACKAWANNA FOUNDRY, LACKAWANNA NEW YORK

| Field Sample No.<br>CLP Sample No.<br>Time Sampled<br>Hazardous Substance | SS-04D<br>MBNJ57<br>15:15 | SS-05A<br>MBNJ58<br>13:35 | SS-05B<br>MBNJ59<br>13:40 | SS-05C<br>MBNJ60<br>13:45<br>Dup. Of SS-03 | SS-06A<br>MBNJ62<br>13:00 | SS-06B<br>MBNJ63<br>13:05 | SS-06C<br>MBNJ64<br>13:10 | NYSDEC<br>Recommended<br>Soil Cleanup<br>Objective |
|---|---------------------------|---------------------------|---------------------------|--|---------------------------|---------------------------|---------------------------|--|
| Aluminum  | 15,100                    | 13,800                    | 16,200                    | 18,300                                     | 15,400                    | 13,500                    | 12,200                    | SB   |
| Antimony  | ---                       | ---                       | ---                       | ---  | ---                       | ---                       | ---                       | SB   |
| Arsenic   | 16.9                      | 5.4                       | 4.8                       | 1.9 B                                      | 4.3                       | 6.9                       | 8.0                       | 7.5 or SB  |
| Barium  | 166                       | 129                       | 157                       | 179  | 182                       | 114                       | 89.8                      | 300 or SB  |
| Beryllium   | 1.0 B                     | 0.72 B                    | 0.84 B                    | 0.89 B                                     | 0.85 B                    | 0.76 B                    | 0.58 B                    | 0.16 or SB   |
| Cadmium   | ---                       | ---                       | ---                       | ---  | ---                       | ---                       | ---                       | 1 or SB  |
| Calcium   | 2,630                     | 2,210                     | 2,260                     | 2,250                                      | 5,220                     | 2,450                     | 2,000                     | SB   |
| Chromium  | 23.7                      | 20.4                      | 24                        | 26.1                                       | 21.8                      | 19.2                      | 17.6                      | 10 or SB   |
| Cobalt  | 12.7                      | 12.6                      | 14.9                      | 11.0 B                                     | 13.2                      | 15.2                      | 13                        | 30 or SB   |
| Copper  | 41.9 J                    | 23.4 J                    | 25.1 J                    | 19.9 J                                     | 34.4 J                    | 23.9 J                    | 22.0 J                    | 25 or SB   |
| Iron  | 52,800                    | 26,700                    | 31,000                    | 27,300                                     | 26,000                    | 29,800                    | 28,300                    | 2000 or SB   |
| Lead  | 25.2                      | 19.5                      | 28.8                      | 17.8                                       | 24                        | 18.8                      | 14.1                      | SB   |
| Magnesium   | 5,450                     | 4,730                     | 5,190                     | 5,830                                      | 4,810                     | 4,250                     | 4,000                     | SB   |
| Manganese   | 197                       | 364                       | 391                       | 261  | 484                       | 427                       | 457                       | SB   |
| Mercury   | ---                       | ---                       | ---                       | ---  | ---                       | ---                       | ---                       | 0.1  |
| Nickel  | 48.9                      | 38.8                      | 42.2                      | 45.9                                       | 44.8                      | 36.5                      | 32.5                      | 13 or SB   |
| Potassium   | 1,030 B                   | 1,200 B                   | 1,420 J                   | 1,940 J                                    | 1,870 J                   | 1,420 J                   | 1,140 B                   | SB   |
| Selenium  | ---                       | 1.8                       | 1.5                       | 1.8  | 1.4                       | 2.3                       | 2.2                       | 2 or SB  |
| Silver  | 1.3 B                     | 0.56 B                    | 0.73 B                    | 0.43 B                                     | 0.42 B                    | 0.58 B                    | 0.50 B                    | SB   |
| Sodium  | ---                       | ---                       | 138 B                     | 161 B                                      | ---                       | 108 B                     | 92. B                     | SB   |
| Thallium  | ---                       | 2.3 B                     | ---                       | 1.2 B                                      | ---                       | ---                       | 2.2 B                     | SB   |
| Vanadium  | 24                        | 20.6                      | 24                        | 24.7                                       | 23.8                      | 21.8                      | 18.5                      | 150 or SB  |
| Zinc  | 127                       | 108                       | 134                       | 134  | 139                       | 96.9                      | 88.5                      | 20 or SB   |
| Cyanide   | ---                       | ---                       | ---                       | ---  | ---                       | ---                       | ---                       | NA   |

Notes:

All results are reported in units of mg/kg.

Dashed lines (--) denote hazardous substance not detected.

Shading denotes concentration exceeding NYSDEC Recommended Soil Cleanup Objective.

J-Estimated value.

B-Indicates concentration between the instrument detection limit and the contract required detection limit.

SB-site background

NA-None available

**TABLE 2**

**GROUNDWATER SAMPLE ANALYSIS RESULTS  
FOR EPA SAT BROWNFIELD SITE INVESTIGATION (NOVEMBER 2000)  
LACKAWANNA FOUNDRY, LACKAWANNA, NEW YORK**

| Sample No.<br>Hazardous Substance | GW-01<br>(Bedrock/<br>Downgradient) | GW-03<br>(Overburden/<br>Sidegradient) | GW-04<br>(Overburden/<br>Upgradient) | GW-05<br>(Overburden/<br>Upgradient) | GW-06 (dup.<br>of GW-01) | NYSDEC Ambient<br>Water Quality<br>Standard or<br>Guidance Values |
|-----------------------------------|-------------------------------------|--|--------------------------------------|--------------------------------------|--------------------------|---|
| <b>VOCs</b>                       |                                     |  |                                      |                                      |                          |   |
| Acetone                           | ---                                 | 11                                     | 9.0 J                                | 6.0 J                                | ---                      | 50  |
| <b>METALS</b>                     |                                     |  |                                      |                                      |                          |   |
| Aluminum                          | ---                                 | ---                                    | 3,500                                | 2,500                                | ---                      | NA  |
| Arsenic                           | ---                                 | ---                                    | 13                                   | 12                                   | ---                      | 25  |
| Barium                            | 380                                 | 240                                    | 220                                  | 300                                  | 380                      | 1000  |
| Calcium                           | 250,000                             | 100,000                                | 60,000                               | 87,000                               | 250,000                  | NA  |
| Chromium                          | ---                                 | ---                                    | ---                                  | 23                                   | ---                      | 50  |
| Iron                              | 13,000                              | 6,200                                  | 42,000                               | 34,000                               | 13,000                   | 300   |
| Lead                              | ---                                 | ---                                    | ---                                  | 3.4                                  | ---                      | 25  |
| Magnesium                         | 130,000                             | 22,000                                 | 14,000                               | 22,000                               | 120,000                  | 35000   |
| Manganese                         | 300                                 | 1,700                                  | 1,700                                | 2,200                                | 300                      | 300   |
| Potassium                         | 8,100                               | ---                                    | ---                                  | ---                                  | 8,000                    | NA  |
| Sodium                            | 540,000                             | 22,000                                 | 25,000                               | 20,000                               | 550,000                  | 20000   |
| Zinc                              | ---                                 | 33                                     | 76                                   | ---                                  | ---                      | 2000  |
| Cyanide                           | ---                                 | ---                                    | 12                                   | ---                                  | ---                      | 200   |

Notes:

All units in ug/l.

J - Estimated value.

- Not Detected

Shading Denotes concentration exceeding New York State Department of Environmental Conservation Ambient Water Quality Standards and Guidance Values.

Analytes not listed were not detected in any groundwater sample.

Rinsate blank and trip blank were non-detect for all analytes.

NA - None available

REFERENCE NO. 1

**U.S. ENVIRONMENTAL PROTECTION AGENCY  
FINAL POLLUTION REPORT**

**I. HEADING**

**Date:** September 15, 2000  
**From:** Kevin M. Matheis, On-Scene-Coordinator  
U.S. EPA Region II  
**To:** R. Salkie, EPA  
T. Johnson, 5202G  
R. Cahill, 2CD-PAT  
P. Simon, 2ORC-NYCSFB  
J. LaPadula, 2ERRD-NYRB  
T. Rivero, 2OPM-GCMB  
M. O'Toole, NYSDEC  
ERD-Washington (E-Mail)  
START  
V. Pitruzzello, 2ERRD-PSB  
J. Rotola, EPA  
B. Bellow, 2CD  
M. Basile, 2CD-POB  
B. Carr, 2ORC-NYCSFB  
R. Byrnes, 2OIG  
B. Deese, 2ERRD-RAB  
M. Doster, NYSDEC IX  
J. Guyer, City of Lackawanna  
P. Brandt, EPA

**Subject:** Lackawanna Foundry Site, Lackawanna, New York - PCB transformers and capacitors, PCB-contaminated soil, arsenic contaminated soil, drums, lab chemicals and compressed gas cylinders

**POLREP NO:** Fifteen (15) and **Final**

**II. BACKGROUND**

**SITE/SPILL NO.:** 02MY  
**RESPONSE AUTHORITY:** CERCLA/SARA  
**NPL STATUS:** Non-NPL  
**CERCLIS ID:** NYR000073064  
**START DATE:** March 22, 1999  
**COMPLETION DATE:** **September 15, 2000**  
**FUNDING DATE:** Verbal authorization of funds from Deputy Division Director on March 19, 1999. Action Memo Signed on August 13, 1999

**VERBAL AUTHORIZATION**

**MEMO STATUS:** Sent to Deputy Director on March 25  
**ACTION MEMO STATUS:** Signed August 13, 1999

**III. RESPONSE INFORMATION**

**A. Situation**

The Lackawanna Foundry Site is located at 2 Elm Street, Lackawanna, New York. This site was the location of a foundry that produced speciality cast iron molds for over 60 years. The foundry owner passed away approximately seven years ago and operations decreased on-site until discontinuing prior to EPA's removal action. The Site is located in a mixed residential and

heavily industrialized area of Lackawanna, New York, adjacent to Smokes Creek, a direct tributary to Lake Erie. Twelve residential properties are immediately adjacent to the site to the west, railroad tracks border the east, Smokes Creek to the north, and a multi-use city park to the South.

Prior to EPA actions at the site, the New York State Department of Environmental Conservation (NYSDEC) ordered the property owner to collect samples from transformers and soil on-site and provide NYSDEC with an inventory of chemicals on-site. The owner complied with this request and analysis of transformers and soils indicated that two transformers contain 86% PCBs. Soil contamination in proximity of the transformers indicated the presence of PCB contaminated soils up to 210 ppm. The inventory also showed more than 100 drums located at the site, stored in various stages of deteriorated condition. NYSDEC requested that the owner undertake a cleanup at the site, but he did not have the financial ability to comply with this request.

As a result, the NYSDEC requested that EPA evaluate the site for a removal action. EPA performed a removal assessment which was completed on March 15, 1999. During this assessment, EPA became aware of several problem areas at the site. The site was completely unfenced and contained over 100 unknown drums outside the foundry buildings, and another 200 drums, laboratory chemicals, compressed gas cylinders, and PCB transformers within the foundry buildings. In addition, the buildings that contain the drums and transformers are in extremely poor condition and several access points to the building were identified. Evidence of trespassing also existed throughout the property. As recently as February, 1999, the attorney for the owner has written to the City of Lackawanna and told them to demolish the buildings on-site because the owner does not have the financial ability to undertake a removal action or comply with building code violations.

## **B. Actions Taken**

EPA began its removal action in March 22, 1999. Initial actions began with the fencing of the perimeter of the site, including fence installation along the residential properties adjacent to the site. EPA then inventoried and moved all containers of hazardous substances into the foundry buildings and over-packed the leaking drums and containers as necessary. EPA then secured the foundry buildings. EPA's Environmental Response Team (ERT) was mobilized to the site and they collected soil samples from the site property and the recreational area adjacent to the site. Results of the sampling showed the presence of PCB contaminated soils inside the foundry buildings and arsenic contaminated soils in the recreational areas.

The Action Memorandum to address the cleanup at the site was signed on August 13, 1999 and authorized the following actions at the site: the removal of all PCB transformers, capacitors and contaminated soil and debris; removal of drums, laboratory chemicals, and compressed gas cylinders; removal of arsenic contaminated soils from recreational area adjacent to site; possible demolition of on-site buildings; and restoration of site upon completion of removal actions.

Due to the location of the soil contamination in proximity to the site buildings and overall

instability of the site buildings, EPA determined that demolition of the buildings would be necessary to complete its removal action. The site buildings were demolished in November 1999. Soil excavations of contaminated soils were completed in April 2000. Site restoration actions were completed in June 2000. The site was seeded with grass and the wetlands in the center portion of the site was restored by the use of clean marsh sediment from a New York State Parks dredging project.

The site restoration has been satisfactorily completed. The City of Lackawanna will install barriers to prevent unauthorized vehicular access.

**C. Future Actions**

EPA has designated this site as eligible for an EPA-funded targeted site assessment. This site assessment, when implemented and completed, will provide analytical data to the NYSDEC for site classification purposes and will hopefully lead to acquisition of the site by the City of Lackawanna.

**D. Key Issues**

None

**IV. COST INFORMATION: As of 09/15/00**

| <b>Site Ceiling as of 09/15/00</b> |                      |                      |                         |
|------------------------------------|----------------------|----------------------|-------------------------|
|                                    | <b>Total Ceiling</b> | <b>Costs to Date</b> | <b>Amount Remaining</b> |
| Earth Tech                         | \$ 1,070,000         | \$ 1,060,000         | \$ 10,000               |
| START/EPA                          | \$ 200,000           | \$ 198,000           | \$ 2,000                |
| <b>Totals</b>                      | <b>\$ 1,270,000</b>  | <b>\$ 1,258,000</b>  | <b>\$ 12,000</b>        |

**The above accounting of expenditures is an estimate based on figures known to the OSC at the time this report was written. The OSC does not necessarily receive specific figures on final payments made to any contractor(s). Other financial data which the OSC must rely upon may not be entirely up-to-date. The cost accounting provided in this report does not necessarily represent an exact monetary figure which the government may include in any claim for cost recovery.**

**V. DISPOSITION OF WASTES**

**LACKAWANNA FOUNDRY SITE  
WASTE DISPOSAL TABLE**

| Waste Material Description                         | Volume                                    | Disposal Facility  | Disposal Method                     |
|--|---|--|-------------------------------------|
| PCB Contaminated Debris<br>>500ppm                 | 160 y <sup>3</sup>                        | CWM - Model City,<br>NY  | Landfill                            |
| PCB Contaminated Soil<br>>500ppm                   | 360 y <sup>3</sup>                        | CWM - Model City,<br>NY  | Landfill                            |
| PCB Capacitors                                     | 49 drums<br>containing 153<br>capacitors. | CWM - Port Arthur,<br>TX   | Incineration                        |
| PCB Oils   | 2 drums                                   | Safety Kleen -<br>Twinsburg,OH   | Treatment/<br>Incineration          |
| PCB Transformers >500ppm                           | 6 transformers<br>(1,916 lbs.)            | Safety Kleen -<br>Twinsburg,OH   | Treatment/<br>Incineration          |
| PCB Transformers <500ppm                           | 2 transformers<br>(2,467 lbs.)            | Safety Kleen -<br>Twinsburg,OH   | Treatment/<br>Incineration          |
| Drums containing various<br>hazardous substances   | 68 drums                                  | Varies   | Treatment/Incin<br>eration/Landfill |
| Foundry slag                                       | 20 tons                                   | CID Refuse,<br>Chaffee,NY  | Landfill                            |
| Hazardous substance<br>containing debris, non-RCRA | 550 y <sup>3</sup>                        | BFI Landfill,<br>Niagara Falls, NY   | Landfill                            |
| Concrete from building<br>demolition               | 1,600 tons                                | CTS Recycling  | Recycling                           |
| Structural and non-structural<br>steel             | 220 tons                                  | Gateway Recycling,<br>Cheektowaga, NY and<br>Lake Erie Recycling,<br>Buffalo, NY | Recycling                           |
| Wood and miscellaneous<br>debris from demolition   | 650 y <sup>3</sup>                        | Modern Disposal,<br>Model City, NY   | Landfill                            |
| Arsenic contaminated soils                         | 900 tons                                  | Chaffee Landfill,<br>Chaffee, NY   | Landfill                            |
| PCB contaminated soils<br><50ppm                   | 400 tons                                  | Modern Disposal,<br>Model City, NY   | Landfill                            |
| PCB contaminated soils (50 -<br>500ppm)            | 280 tons                                  | CWM, Model City,<br>NY   | Landfill                            |



REFERENCE NO. 2



**FIELD SAMPLING PLAN  
FOR  
LACKAWANNA FOUNDRY SITE  
LACKAWANNA, ERIE COUNTY, NEW YORK**

**Prepared for:**  
**U.S. Environmental Protection Agency**  
Region II  
New York, New York 10007

**Prepared by:**  
**Region II Site Assessment Team**  
Roy F. Weston, Inc.  
Edison, New Jersey 08837

U. S. EPA Contract No. 68-W-00-121

October 2000



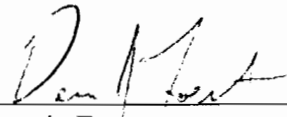
**FIELD SAMPLING PLAN  
FOR  
LACKAWANNA FOUNDRY SITE  
LACKAWANNA, ERIE COUNTY, NEW YORK**

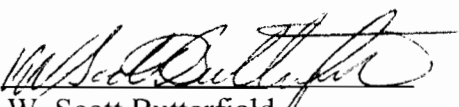
**Prepared by:**  
Region II Site Assessment Team  
Roy F. Weston, Inc.  
U.S. EPA CONTRACT NO. 68-W-00-121

DCN: SAT - 1002.013

October 2000

Prepared by:  Date: 10/31/00  
Donna Janda  
Project Manager

Approved by:  Date: 10/31/00  
Dennis Foerter  
Quality Assurance Officer

Approved by:  Date: 10/31/00  
W. Scott Butterfield  
SAT Program Manager

Approved by: \_\_\_\_\_ Date: \_\_\_\_\_  
Dennis Munhall  
U.S. EPA Work Assignment Manager

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## 1.0 INTRODUCTION

Presented herein is the Field Sampling Plan for the Lackawanna Foundry Site for sampling to be conducted by the Roy F. Weston, Inc. (WESTON®) Region II Site Assessment Team (SAT). This field sampling plan has been developed at the request of the United States Environmental Protection Agency (U.S. EPA) in accordance with the U.S. EPA Region II CERCLA Quality Assurance Manual (October 1989) and the SAT Quality Assurance Project Plan (September 2000).

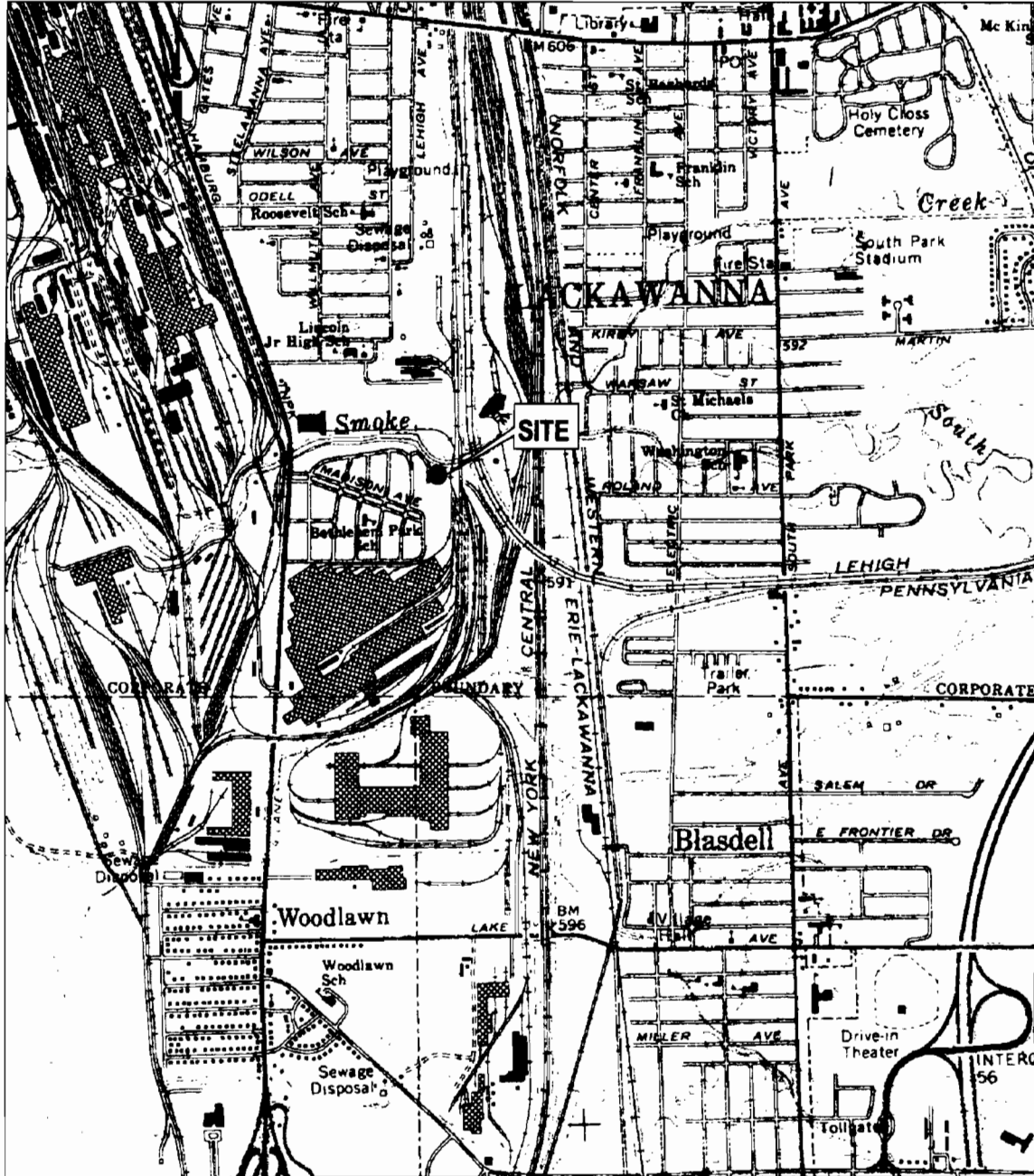
The sampling strategy listed within the sampling plan will emphasize the collection of samples required to evaluate certain pathways of concern to the CERCLA Program. The sampling plan includes the following sections: Site Reconnaissance (2.0), Sampling Visit and Sampling Procedures (3.0), Quality Assurance/Quality Control (4.0), and Field Changes and Corrective Actions (5.0).

### 1.1 Site Description

The Lackawanna Foundry Site is located at 3 Elm Street in Lackawanna, Erie County, New York. The two-acre property is located in a mixed residential and heavily industrialized area adjacent to Smokes Creek, which is a direct tributary to Lake Erie. The Lackawanna Foundry property is bordered by Smoke Creek to the north, by an open field to the south, by an active rail yard to the east, and by private residences to the west. A vegetated wetland area is located in the north central portion of the property. The residential neighborhood immediately adjacent to the Site consists of 12 residences with over 100 residences located within 1/8 mile of the Site. Approximately 1,000 people reside within 0.5 mile of the Site. The location and layout of the Lackawanna Foundry Site can be found in Figure 1, Site Location Map and Figure 2, Site Map.

A cast-iron foundry has operated on this property since the early 1900s. The current property owner inherited the business from his father who operated the foundry at this location for more than 30 years. The current owner ceased operations in approximately 1997.

On 22 December 1998, the New York State Department of Environmental Conservation (NYSDEC) requested that the U.S. Environmental Protection Agency (EPA) evaluate the site and perform a Removal Action, as appropriate, to address the threats to human health, welfare, and the environment from drums, polychlorinated biphenyls (PCBs) in transformers, and PCB-contaminated soil. A drum inventory completed by the property owner listed over 233 drums of various foundry raw materials and waste.



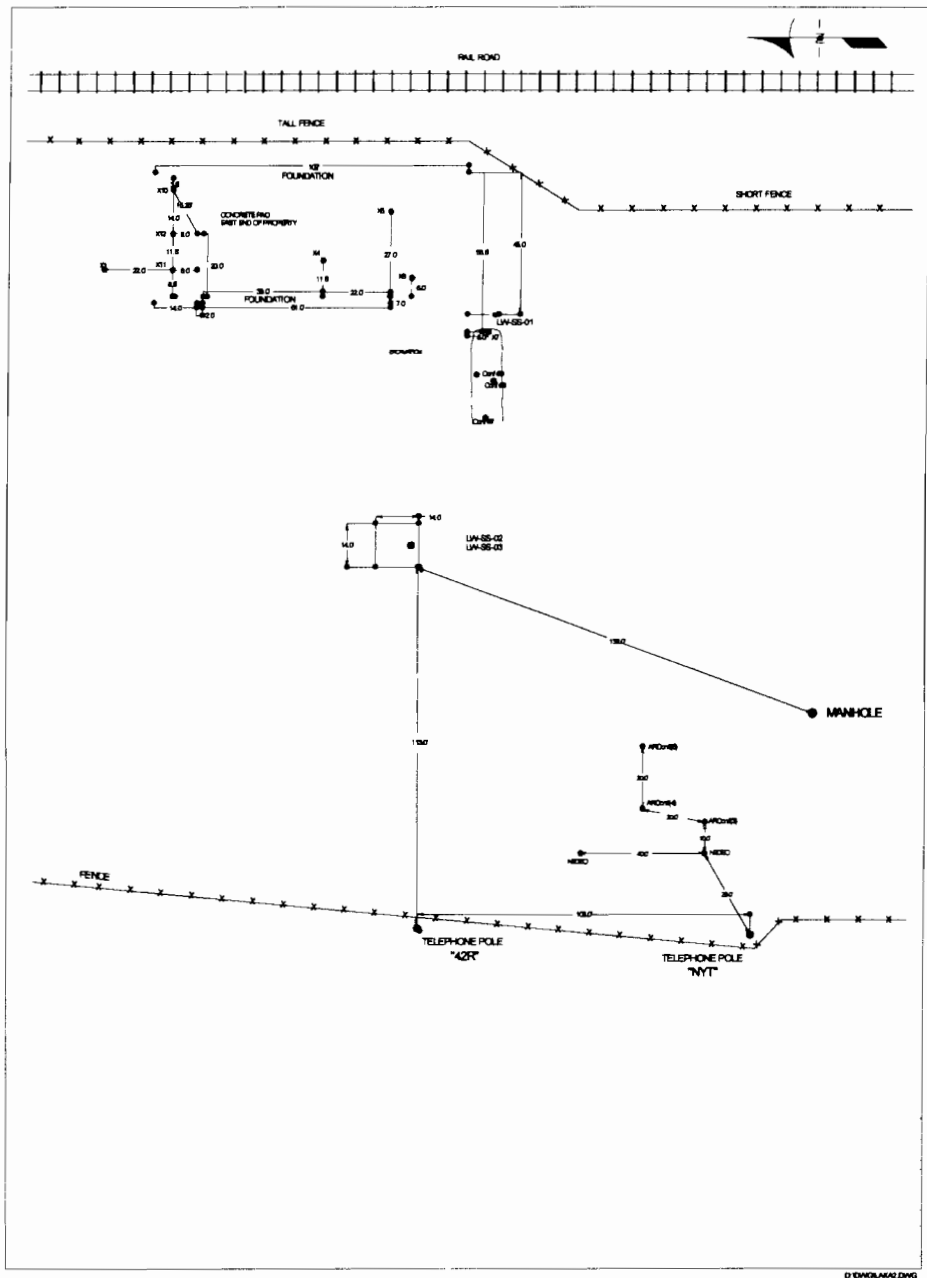
U.S. Geological Survey  
 7.5 Minute Series Topographical Map  
 Buffalo Southeast, NY Quadrangle  
 1948, Revised 1965

1000 0 1000 2000 Feet



Lackawanna Foudry Site  
 Lackawanna, Erie Co., NY  
 Figure 1  
 Site Location Map





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**Figure 2  
Site Map**

**Lackawanna Foundry Site  
Lackawanna, Erie Co., NY**



## **1.2 Previous Work at the Site**

On 22 March 1999, the EPA conducted emergency stabilization activities at the Site, which included fencing the perimeter of the property, securing the buildings, moving drums into a secure location, and stabilizing leaking containers. Two primary areas of contamination were identified at the Lackawanna Foundry Site. The first area included the on-site buildings, which were contaminated with PCBs and contained laboratory chemicals, compressed gas cylinders, paint cans, and several hundred drums. The second area is a recreational area, which is owned by the City of Lackawanna and is located adjacent to the Lackawanna Foundry property. The recreational area is a grassy area that the City maintains by frequent grass and weed cuttings.

Analytical data from an EPA Environmental Response Team soil sampling event conducted in April 1999 indicated arsenic levels as high as 77 parts per million (ppm) in the recreational area. The arsenic contamination in this area appears to be in the location of a former baghouse dust collection system used during foundry operations. The contamination was caused by either airborne deposition or runoff from the baghouse collection area.

EPA conducted a Removal Action at the Lackawanna Foundry property during the spring of 2000. The Removal Action included excavation of contaminated soil and buried PCB-contaminated capacitors, dismantlement of on-site buildings, extent of contamination sampling, confirmation sampling, transportation and disposal of drums, contaminated soil and demolition material, restoration of on-site wetlands area, and backfilling of excavated areas with certified clean topsoil.

## **1.3 Schedule**

EPA and SAT conducted an on-site reconnaissance on 4 October 2000.

The anticipated date for subsurface soil sampling and monitoring well installation is the week of 6 November 2000. The anticipated date for monitoring well sampling is the week of 11 December 2000.

Sample analysis is expected to take 10 to 12 weeks. Receipt of analytical data is expected during March 2001.

The draft Brownfields Site Assessment Report is anticipated to be completed by April 2001. The final report is anticipated to be completed within 30 days of receipt of EPA comments.

## 2.0 SITE RECONNAISSANCE

Region II SAT accompanied the EPA Work Assignment Manager and On-Scene Coordinator on a site reconnaissance of the Lackawanna Foundry property on 4 October 2000. There are currently no buildings or other structures on the Lackawanna Foundry property. The relatively flat site is well vegetated and is surrounded by a fence with a locked gate to control access. All known waste sources were removed from the property during the Removal Action conducted earlier this year. There were no signs of stressed vegetation or contaminated soil.

Sample locations were selected based on the recommendations of the NYSDEC to assist in a Brownfields site investigation. The Lackawanna Foundry property is being considered for future use as a recreational park. The nearest residences are located adjacent to the west side of the property. The Bethlehem Steel property is located approximately 0.25 mile south of the Lackawanna Foundry Site.

### 3.0 SAMPLING VISIT AND SAMPLING PROCEDURES

This section outlines overall sample management and control procedures to be implemented by WESTON personnel during field activities. Standard analytical methods, preservation, holding times, and sample containers are summarized in Table 3-1.

#### 3.1 Sample Tracking System

##### 3.1.1 Sample Identification System

Each sample will be designated by a numeric code which will identify the site. The numeric code will be the site-specific task number (1002). The media type will follow the site code. A hyphen will separate the site code and media type. Specific media types are as follows:

GW - Groundwater  
RB - Rinsate Blank  
SS - Subsurface Soil  
TB - Trip Blank

After the media type, the sequential sample numbers will be listed; sample numbers will begin with 01 and increase accordingly. For example, two groundwater samples collected from a site may be designated as Sample Nos. 1002-GW-01 and 1002-GW-02. A duplicate sample will be identified in the same manner as other samples and will be distinguished in the field logbook.

##### 3.1.2 Sample Bottles

Sample bottles will be obtained from qualified vendors via a competitive bid process and will meet all guidelines specified in OSWER Directive 9240.0-05A, Specifications and Guidance for Obtaining Contaminant-Free Sample Containers (December, 1992).

##### 3.1.3 Sample Packaging and Shipping

Samples will be packaged and shipped according to the U.S. EPA User's Guide to the Contract Laboratory Program (January 1991). Chain of custody forms, sample labels, custody seals, and other sample documents will be completed as specified in the above reference manual. All entries will be made in permanent ink. If errors are made when completing any of these forms, the error will be crossed out with a single line, initialed, and dated by the sampler. Each environmental sample will be properly identified and sealed in a polyethylene bag. The bag shall then be placed in a plastic cooler which has also been lined with a large polyethylene bag. When required, samples shall be packed with sufficient ice (sealed in polyethylene bags) to cool the samples to 4°C. Sufficient non-combustible, adsorbent cushioning material shall be placed in the cooler so as to minimize the possibility of container breakage. The large plastic bag shall then be sealed and the container closed. Custody seals and strapping tape shall then be affixed to the outer packaging. All samples will either be hand-delivered or shipped via common carrier to the laboratory within 24 hours of collection. Sample shipment will conform to Roy F. Weston, Inc., Guidelines For

Classifying Field Sample Shipments and the most current International Air Transport Association (IATA) Dangerous Goods Regulations. Information relating to the shipment of samples, including the airbill number, sample quantity, and sample types, will be reported to the U.S. EPA Sample Management Office (SMO) on the day of or morning after shipment.

#### **3.1.4 Sample Documentation**

The sampling team or individual performing a particular sampling activity is required to maintain a field logbook. The bound, numbered, and paginated logbook shall be filled out at the location of sample collection immediately after sampling. The logbook shall contain sampling information, including: sample number, sample collection time, sample location, sample descriptions, sampling methods, weather conditions, field measurements, name of sampler, site-specific observations, and any deviations from protocol. All entries will be entered legibly in permanent ink. If errors are made when completing this logbook, the error will be crossed out with a single line, initialed, and dated by the sampler.

#### **3.2 Sampling Program**

A description of all site-specific samples, including the rationale for the collection of each sample, are presented in Table 3-2. All proposed sample locations are presented in Figure 3, Sample Location Map. Site-specific sample analyses, bottle types, and sample preservatives are presented in Table 3-3. Samples will be analyzed for Target Compound List (TCL) and Target Analyte List (TAL) parameters through the U.S. EPA Contract Laboratory Program (CLP).

Listed below are standard operating procedures which will be adhered to during field sampling activities conducted by Region II SAT.

**TABLE 3-1  
CLP ROUTINE ANALYTICAL SERVICES**

| Sample Location | Number of Samples  | Matrix                             | Sampling Device               | Sample Container (1)   | Sample Preservation                     | Holding Time*  | CLP Laboratory Analyses (2) |
|-----------------|--|------------------------------------|-------------------------------|--|---|--|-----------------------------|
| GW-01 to GW-06  | 6 (1 at each location & 1 duplicate)                                     | Aqueous (Low/Medium Concentration) | Low flow pump                 | Two 40-ml glass vials w/ teflon septum caps                      | HCl to pH<2<br>Cool to 4°C              | 10 days (preserved); (7 days [unpreserved] to analyze aromatic hydrocarbons) | TCL VOAs<br>CLP SOW         |
|                 |  |                                    |                               | Four 1-L amber glass bottles                                     | Cool to 4°C                             | 5 days to extract;<br>40 days to analyze                                     | TCL Extractables<br>CLP SOW |
|                 |  |                                    |                               | One 1-L polyethylene bottle                                      | HNO <sub>3</sub> to pH<2<br>Cool to 4°C | 6 months to analyze (except Hg - 28 days)                                    | TAL Inorganics<br>CLP SOW   |
|                 |  |                                    |                               | One 1-L polyethylene bottle                                      | NaOH to pH>12<br>Cool to 4°C            | 12 days to analyze   | Cyanide<br>CLP SOW          |
| SS-01 to SS-07  | 16 (1 each at SS-01 to SS-03; 4 each at SS-04 to SS-06; & one duplicate) | Soil (Low/Medium Concentration)    | Split spoon<br>Bowl<br>Trowel | Two 40-ml wide-mouth jars w/ teflon septum caps (3)              | Cool to 4°C                             | 10 days to analyze   | TCL VOAs<br>CLP SOW         |
|                 |  |                                    |                               | One 8-oz. wide-mouth glass jar                                   | Cool to 4°C                             | 5 days to extract;<br>40 days to analyze                                     | TCL Extractables<br>CLP SOW |
|                 |  |                                    |                               | One 8-oz. wide-mouth glass jar                                   | Cool to 4°C                             | 6 months to analyze (except Hg - 28 days)                                    | TAL Inorganics<br>CLP SOW   |
|                 |  |                                    |                               | One 8-oz. wide-mouth glass jar (included in TAL sample fraction) | Cool to 4°C                             | 12 days to analyze   | Cyanide<br>CLP SOW          |

\* = All holding times are calculated from Verified Time of Sample Receipt (VTSR).

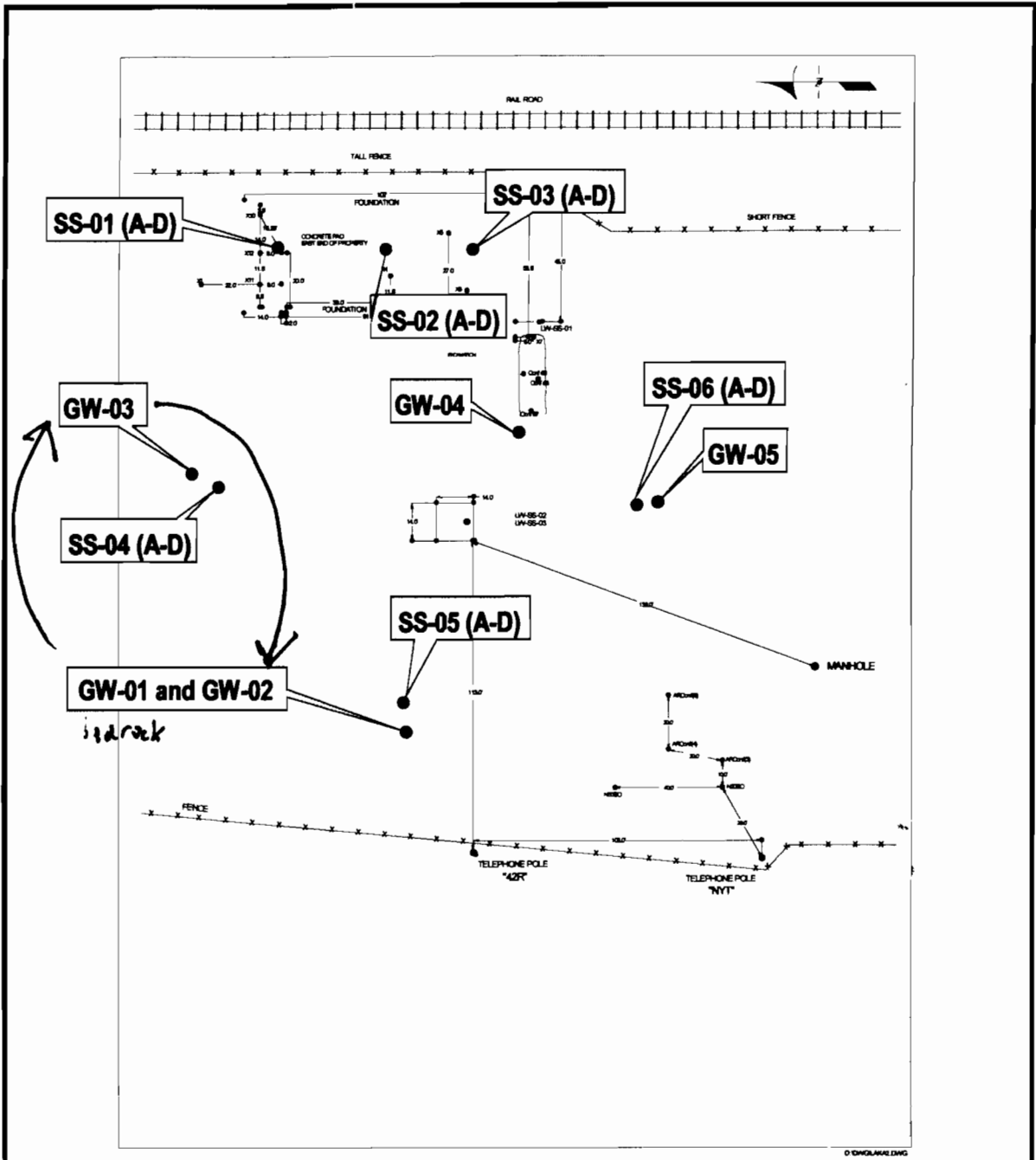
(1) = Sample containers are certified clean by the manufacturer.

(2) = Contract Laboratory Program (CLP) Statement of Work (SOW) for Organic Analysis OLM04-2, or most current revision; and, CLP SOW for Inorganic Analysis ILM04-1, or most current revision.

(3) = Two 120-ml wide-mouth glass jars may be used if large soil particles preclude 40-ml vials from being completely filled.

p. 8

*Mon. drum cuttings, develop curb boxes*



- Proposed ground-water sample location
  - Proposed subsurface soil sample location
  - Historical sample location
- A = Depth of 2-4 feet  
 B = Depth of 4-6 feet  
 C = Depth of 6-8 feet  
 D = Depth of 8-10 feet



**Figure 3**  
**Sample Location Map**

**Lackawanna Foundry Site**  
**Lackawanna, Erie Co., NY**

### 3.2.1 Subsurface Soil Sampling

The following procedures apply to the collection of subsurface soil, from a depth greater than 6 inches, using a stainless steel split spoon:

- 1) Wear protective gear as specified in the Health and Safety Plan. Samplers shall don new outer gloves prior to sampling at each location.
- 2) Use a decontaminated, 2" or 3" diameter stainless steel split spoon.
- 3) Remove unnecessary rocks, twigs, and other non-soil materials from the selected sampling location.
- 4) Drive split spoon utilizing the direct push method (i.e., Geoprobe). Do not drive past the bottom of the head piece as this will result in the compression of the sample.
- 5) Withdraw the spoon and open by unscrewing bit and head and splitting barrel. Transfer the samples using clean stainless steel utensils into a clean stainless steel bowl. Collect PID and/or FID readings; record results in field logbook. Repeat step number 4 at an adjacent location until sufficient soil is collected to satisfy analytical requirements.
- 6) Fill the volatile sample bottles immediately so as to not compromise sample integrity.
- 7) Homogenize remaining soil for the non-volatile organic analysis fractions in the bowl using a decontaminated, stainless steel utensil. Transfer the samples into required sample containers. Homogenization shall be completed as per the following procedure:

After collection of the volatile sample fraction, the soil in the stainless steel bowl will be scraped from the sides, corners and bottom of the bowl, rolled to the middle of the bowl, and mixed. The soil will then be quartered and moved to the four corners of the bowl. Each quarter will then be mixed individually, and when completed be rolled to the center of the bowl and mixed once again.
- 8) Place samples in cooler and chill with ice. Samples will be hand-delivered or shipped within 24 hours of collection to the designated CLP laboratory(ies).
- 9) Fill out field logbook, custody seals, sample labels, and chain of custody forms.

**TABLE 3-2**  
**SAMPLE DESCRIPTIONS/RATIONALE**

| SAMPLE NUMBER  | DESCRIPTION/RATIONALE   |
|----------------|---|
| 1002-GW-01     | Sample to be collected from on-site monitoring well MW-01 to determine if hazardous substances are migrating from on-site soils into groundwater.   |
| 1002-GW-02     | Sample to be collected from on-site monitoring well MW-02 to determine if hazardous substances are migrating from on-site soils into groundwater.   |
| 1002-GW-03     | Sample to be collected from on-site monitoring well MW-03 to determine if hazardous substances are migrating from on-site soils into groundwater.   |
| 1002-GW-04     | Sample to be collected from on-site monitoring well MW-04 to determine if hazardous substances are migrating from on-site soils into groundwater.   |
| 1002-GW-05     | Sample to be collected from on-site monitoring well MW-05 to determine if hazardous substances are migrating from on-site soils into groundwater.   |
| 1002-GW-06     | Duplicate sample of 1002-GW-05.   |
| 1002-SS-01     | Sample to be collected from under a concrete pad, which is located 3 to 4 feet below the ground surface, to determine the presence of hazardous substances in on-site subsurface soil.  |
| 1002-SS-02     | Sample to be collected from under a concrete pad, which is located 3 to 4 feet below the ground surface, to determine the presence of hazardous substances in on-site subsurface soil.  |
| 1002-SS-03     | Sample to be collected from under a concrete pad, which is located 3 to 4 feet below the ground surface, to determine the presence of hazardous substances in on-site subsurface soil.  |
| 1002-SS-04 A-D | Sample to be collected from an on-site soil boring location at two-foot intervals (A: 2-4 feet; B: 4-6 feet; C: 6-8 feet; and D: 8-10 feet, if necessary) to determine the presence of hazardous substances in the on-site subsurface soil. |
| 1002-SS-05 A-D | Sample to be collected from an on-site soil boring location at two-foot intervals (A: 2-4 feet; B: 4-6 feet; C: 6-8 feet; and D: 8-10 feet, if necessary) to determine the presence of hazardous substances in the on-site subsurface soil. |
| 1002-SS-06 A-D | Sample to be collected from an on-site soil boring location at two-foot intervals (A: 2-4 feet; B: 4-6 feet; C: 6-8 feet; and D: 8-10 feet, if necessary) to determine the presence of hazardous substances in the on-site subsurface soil. |
| SS-07          | Duplicate of 1002-SS-03.  |



**TABLE 3-3  
SAMPLE ANALYSES, BOTTLE TYPES, AND PRESERVATIVES**

| SAMPLE NUMBER                               | SAMPLE BOTTLES                                | ANALYSIS                    | PRESERVATION                            |
|---|---|-----------------------------|---|
| 1002-GW-01                                  | Two 40-mL glass vials<br>w/teflon septum caps | TCL VOA<br>CLP SOW          | HCl to pH<2<br>Cool to 4°C              |
|   | Four 1-L amber glass bottles                  | TCL Extractables<br>CLP SOW | Cool to 4°C                             |
|   | One 1-L polyethylene bottle                   | TAL Inorganics<br>CLP SOW   | HNO <sub>3</sub> to pH<2<br>Cool to 4°C |
|   | One 1-L polyethylene bottle                   | Cyanide<br>CLP SOW          | NaOH to pH>12<br>Cool to 4°C            |
| 1002-GW-02<br>(Triple volume for<br>MS/MSD) | Six 40-mL glass vials<br>w/teflon septum caps | TCL VOA<br>CLP SOW          | HCl to pH<2<br>Cool to 4°C              |
|   | Twelve 1-L amber glass<br>bottles             | TCL Extractables<br>CLP SOW | Cool to 4°C                             |
|   | Three 1-L polyethylene<br>bottle              | TAL Inorganics<br>CLP SOW   | HNO <sub>3</sub> to pH<2<br>Cool to 4°C |
|   | Three 1-L polyethylene<br>bottle              | Cyanide<br>CLP SOW          | NaOH to pH>12<br>Cool to 4°C            |
| 1002-GW-03                                  | Two 40-mL glass vials<br>w/teflon septum caps | TCL VOA<br>CLP SOW          | HCl to pH<2<br>Cool to 4°C              |
|   | Four 1-L amber glass bottles                  | TCL Extractables<br>CLP SOW | Cool to 4°C                             |
|   | One 1-L polyethylene bottle                   | TAL Inorganics<br>CLP SOW   | HNO <sub>3</sub> to pH<2<br>Cool to 4°C |
|   | One 1-L polyethylene bottle                   | Cyanide<br>CLP SOW          | NaOH to pH>12<br>Cool to 4°C            |
| 1002-GW-04                                  | Two 40-mL glass vials<br>w/teflon septum caps | TCL VOA<br>CLP SOW          | HCl to pH<2<br>Cool to 4°C              |
|   | Four 1-L amber glass bottles                  | TCL Extractables<br>CLP SOW | Cool to 4°C                             |
|   | One 1-L polyethylene bottle                   | TAL Inorganics<br>CLP SOW   | HNO <sub>3</sub> to pH<2<br>Cool to 4°C |
|   | One 1-L polyethylene bottle                   | Cyanide<br>CLP SOW          | NaOH to pH>12<br>Cool to 4°C            |
| 1002-GW-05                                  | Two 40-mL glass vials<br>w/teflon septum caps | TCL VOA<br>CLP SOW          | HCl to pH<2<br>Cool to 4°C              |
|   | Four 1-L amber glass bottles                  | TCL Extractables<br>CLP SOW | Cool to 4°C                             |

| SAMPLE NUMBER                      | SAMPLE BOTTLES                                | ANALYSIS                    | PRESERVATION                            |
|------------------------------------|---|-----------------------------|---|
| 1002-GW-05                         | One 1-L polyethylene bottle                   | TAL Inorganics<br>CLP SOW   | HNO <sub>3</sub> to pH<2<br>Cool to 4°C |
|                                    | One 1-L polyethylene bottle                   | Cyanide<br>CLP SOW          | NaOH to pH>12<br>Cool to 4°C            |
| 1002-GW-06<br>(Duplicate of GW-05) | Two 40-mL glass vials<br>w/teflon septum caps | TCL VOA<br>CLP SOW          | HCl to pH<2<br>Cool to 4°C              |
|                                    | Four 1-L amber glass bottles                  | TCL Extractables<br>CLP SOW | Cool to 4°C                             |
|                                    | One 1-L polyethylene bottle                   | TAL Inorganics<br>CLP SOW   | HNO <sub>3</sub> to pH<2<br>Cool to 4°C |
|                                    | One 1-L polyethylene bottle                   | Cyanide<br>CLP SOW          | NaOH to pH>12<br>Cool to 4°C            |
| 1002-TB-01                         | Two 40-mL glass vials<br>w/teflon septum caps | TCL VOA<br>CLP SOW          | HCl to pH<2<br>Cool to 4°C              |
| 1002-RB-01                         | Two 40-mL glass vials<br>w/teflon septum caps | TCL VOA<br>CLP SOW          | HCl to pH<2<br>Cool to 4°C              |
|                                    | Four 1-L amber glass bottles                  | TCL Extractables<br>CLP SOW | Cool to 4°C                             |
|                                    | One 1-L polyethylene bottle                   | TAL Inorganics<br>CLP SOW   | HNO <sub>3</sub> to pH<2<br>Cool to 4°C |
|                                    | One 1-L polyethylene bottle                   | Cyanide<br>CLP SOW          | NaOH to pH>12<br>Cool to 4°C            |
| 1002-RB-02                         | Two 40-mL glass vials<br>w/teflon septum caps | TCL VOA<br>CLP SOW          | HCl to pH<2<br>Cool to 4°C              |
|                                    | Four 1-L amber glass bottles                  | TCL Extractables<br>CLP SOW | Cool to 4°C                             |
|                                    | One 1-L polyethylene bottle                   | TAL Inorganics<br>CLP SOW   | HNO <sub>3</sub> to pH<2<br>Cool to 4°C |
|                                    | One 1-L polyethylene bottle                   | Cyanide<br>CLP SOW          | NaOH to pH>12<br>Cool to 4°C            |
| 1002-RB-03                         | Two 40-mL glass vials<br>w/teflon septum caps | TCL VOA<br>CLP SOW          | HCl to pH<2<br>Cool to 4°C              |
|                                    | Four 1-L amber glass bottles                  | TCL Extractables<br>CLP SOW | Cool to 4°C                             |
|                                    | One 1-L polyethylene bottle                   | TAL Inorganics<br>CLP SOW   | HNO <sub>3</sub> to pH<2<br>Cool to 4°C |
|                                    | One 1-L polyethylene bottle                   | Cyanide<br>CLP SOW          | NaOH to pH>12<br>Cool to 4°C            |

| SAMPLE NUMBER   | SAMPLE BOTTLES                                    | ANALYSIS                          | PRESERVATION |
|---|---|-----------------------------------|--------------|
| 1002-SS-01  | Two 40-mL wide-mouth jars<br>w/teflon septum caps | TCL VOA                           | Cool to 4°C  |
|   | One 8-oz. wide-mouth glass<br>jar                 | TCL Extractables<br>CLP SOW       | Cool to 4°C  |
|   | One 8-oz. wide-mouth glass<br>jar                 | TAL Inorganics/Cyanide<br>CLP SOW | Cool to 4°C  |
| 1002-SS-02  | Two 40-mL wide-mouth jars<br>w/teflon septum caps | TCL VOA                           | Cool to 4°C  |
|   | One 8-oz. wide-mouth glass<br>jar                 | TCL Extractables<br>CLP SOW       | Cool to 4°C  |
|   | One 8-oz. wide-mouth glass<br>jar                 | TAL Inorganics/Cyanide<br>CLP SOW | Cool to 4°C  |
| 1002-SS-03  | Two 40-mL wide-mouth jars<br>w/teflon septum caps | TCL VOA                           | Cool to 4°C  |
|   | One 8-oz. wide-mouth glass<br>jar                 | TCL Extractables<br>CLP SOW       | Cool to 4°C  |
|   | One 8-oz. wide-mouth glass<br>jar                 | TAL Inorganics/Cyanide<br>CLP SOW | Cool to 4°C  |
| 1002-SS-04<br>A (2-4 ft)<br>B (4-6 ft)<br>C (6-8 ft)<br>D (8-10 ft) | Two 40-mL wide-mouth jars<br>w/teflon septum caps | TCL VOA                           | Cool to 4°C  |
|   | One 8-oz. wide-mouth glass<br>jar                 | TCL Extractables<br>CLP SOW       | Cool to 4°C  |
|   | One 8-oz. wide-mouth glass<br>jar                 | TAL Inorganics/Cyanide<br>CLP SOW | Cool to 4°C  |
| 1002-SS-05<br>A (2-4 ft)<br>B (4-6 ft)<br>C (6-8 ft)<br>D (8-10 ft) | Two 40-mL wide-mouth jars<br>w/teflon septum caps | TCL VOA                           | Cool to 4°C  |
|   | One 8-oz. wide-mouth glass<br>jar                 | TCL Extractables<br>CLP SOW       | Cool to 4°C  |
|   | One 8-oz. wide-mouth glass<br>jar                 | TAL Inorganics/Cyanide<br>CLP SOW | Cool to 4°C  |
| 1002-SS-06<br>A (2-4 ft)<br>B (4-6 ft)<br>C (6-8 ft)<br>D (8-10 ft) | Two 40-mL wide-mouth jars<br>w/teflon septum caps | TCL VOA                           | Cool to 4°C  |
|   | One 8-oz. wide-mouth glass<br>jar                 | TCL Extractables<br>CLP SOW       | Cool to 4°C  |
|   | One 8-oz. wide-mouth glass<br>jar                 | TAL Inorganics/Cyanide<br>CLP SOW | Cool to 4°C  |

| SAMPLE NUMBER                      | SAMPLE BOTTLES                                    | ANALYSIS                          | PRESERVATION |
|------------------------------------|---|-----------------------------------|--------------|
| 1002-SS-07<br>(Duplicate of SS-03) | Two 40-mL wide-mouth jars<br>w/teflon septum caps | TCL VOA                           | Cool to 4°C  |
|                                    | One 8-oz. wide-mouth glass<br>jar                 | TCL Extractables<br>CLP SOW       | Cool to 4°C  |
|                                    | One 8-oz. wide-mouth glass<br>jar                 | TAL Inorganics/Cyanide<br>CLP SOW | Cool to 4°C  |

### 3.2.2 Monitoring Well Sampling

Groundwater sampling will be conducted in a manner consistent with current U.S. EPA Region II guidelines. Aeration of the groundwater and disturbance of the sediments/particulate matter within the well bore during sampling will be minimized by matching the pump intake velocity with the natural groundwater flow velocity, thereby providing samples which are more representative of the aquifer conditions. This sampling method generally yields low turbidity samples, eliminating the need for the collection and analysis of filtered groundwater samples.

Six groundwater samples will be collected from five monitoring wells to be installed on site during the week of 6 November 2000. All sampling locations are presented in Figure 3.

Wells will generally be purged using a stainless steel, electric, adjustable rate submersible pump. Pumps will be decontaminated before and after sampling events, as well as between well sampling locations. Pumps shall use ASTM 2239 Teflon or Teflon-lined polyethylene tubing dedicated to each well; the same tubing will be used for purging and sampling. The Teflon or Teflon-lined tubing is suitable for the collection of both organic and inorganic sample fractions. When sampling for inorganic analysis, PVC, Tygon, or polyethylene tubing may be used. If immiscible layers, such as Light and Dense Non-Aqueous Phase Liquids (LNAPLs, DNAPLs), are detected using the interface probe prior to sampling, the LNAPL/DNAPL samples may be collected using a bottom-filling bailer of stainless steel and/or Teflon construction.

The following procedures apply to the collection of monitoring well groundwater samples:

- 1) Wear protective gear as specified in the Health and Safety Plan. Samplers shall don new outer sampling gloves prior to sampling at each location.
- 2) Visually examine the exterior of the monitoring well for signs of damage or tampering; record findings in the field logbook.
- 3) Collect PID and/or FID readings at the well rim; record results in field logbook.
- 4) Unlock well cap, if present.

- 5) Collect PID and/or FID readings; record results in field logbook.
- 6) Measure the static water level in the well with a steel tape or electronic water level indicator. Caution will be exercised in order to minimize disturbance of the water column and any associated sediment/particulate matter. The depths to water of all wells to be sampled will be measured and recorded in the field logbook prior to purging of the wells. The tape or water level indicator will be rinsed with deionized water in between individual wells to prevent cross contamination.
- 7) Measure the depth of any immiscible phases (e.g., DNAPLs/LNAPLs) using an interface probe; record observations in field logbook. During measurement, care will be taken in order to minimize disturbance of any accumulated sediment. The probe will be rinsed with deionized water between individual wells to prevent cross contamination.
- 8) Place clean, polyethylene sheeting around well casing to prevent contamination of sampling equipment in the event equipment is dropped.
- 9) Slowly lower the pump apparatus into the well until it is situated in the middle or slightly above the well screen, ensuring that the pump intake is kept at least 2 feet above the well bottom; this pump position will minimize the potential of pumping stagnant water and the disturbance of bore sediments and DNAPLs.
- 10) Measure the water level again with the pump in the well; record observations in the field logbook.
- 11) Begin purging the well at a low rate (e.g., 500 ml/min). At a minimum, the water level should be monitored every three to five minutes during pumping. A flow rate resulting in a stabilized water table should be established and maintained; drawdown should be less than 0.3 foot. Pump suction will also be maintained in order to avoid introduction of air into the tubing during pumping. In the field logbook, record each pumping rate adjustment and subsequent water level measurement. In the event that the well recharge rate is very low, caution will be used in order to prevent pressure loss in the pump tubing, groundwater recharge cascading through the sand pack, and complete evacuation of the well.

During well purging, measure the following field indicator parameters every three to five minutes and record in the field logbook: temperature, pH, turbidity, dissolved oxygen (DO), specific conductance, and redox potential (Eh). When possible, in-line analyzers and continuous readout displays will be utilized. Measurements for DO and Eh will be obtained from groundwater prior to exposure to air. The remainder of parameters may be either in-line or from a clean glass container holding the groundwater sample. Purging will be continued until well is stabilized; the well is considered stabilized when all

field indicator parameters are  $\pm 10\%$  for three consecutive readings *and* turbidity is less than 50 NTU. If the field indicators have stabilized and the turbidity is greater than 50 NTU, the pump rate will be decreased to  $\leq 250$  ml/minute and the field indicator parameters will be continued to be monitored. When the turbidity decreases to less than 50 NTU, the groundwater well samples can be collected. The recommended sample collection flow rate is 100 to 250 ml/min.

The maximum time between purging and sampling will be three hours. If the well is allowed to sit longer than three hours, the well will be re-evacuated since the water in the casing may no longer be representative of the aquifer.

- 12) Upon U.S. EPA approval, purge water will be discharged to the ground surface in the immediate vicinity of the well. Should the approval for purge water discharge to ground surface be denied, the purge water will be containerized and disposed of as per OERR Directive 9345.3-02, Management of Investigation-Derived Wastes During Site Inspections (May 1991), or most current directive, and Region II SAT procedures for disposal of investigation-derived waste.
- 13) Collect the VOA sample fraction first directly from the tubing into pre-preserved glass vials to prevent degassing of the water. (See Item No. 14, below, for preservation procedures.) During sample collection, the containers will be held so the discharge causes minimal internal disturbance. Collect the remaining organic and inorganic sample fractions.
- 14) Samples shall be preserved as follows:
  - a) VOAs - Determine the amount of 1:1 HCl preservative required to adjust the pH of the sample to less than 2 in an extra 40 ml glass vial. If effervescence occurs when the bottle is tapped, volatile samples will be submitted without preservative and noted as such on the respective Traffic Report. Add the determined volume of acid to the empty 40 ml glass vials prior to sampling. Fill each container with sample to just overflowing so that no air bubbles are trapped inside. This procedure will be repeated at each new sampling station.
  - b) Other Parameters - Fill each container and preserve immediately as required in Table 2-1. When adjusting the pH for sample preservation, pour a minimal portion of sample onto broad range pH paper to verify if the appropriate pH level has been obtained.
- 15) Place samples in cooler and chill with ice. Samples will be hand-delivered or shipped within 24 hours of collection to the designated CLP laboratory(ies).
- 16) Remove the pump apparatus. The used tubing will be disposed of properly. Gross surface decontamination of the pump will be accomplished by placing

the pump in a large container filled with a detergent/water solution. The pump will then be submerged in another large container filled with potable water and re-started. Water will be allowed to circulate through the pump until it is thoroughly rinsed. The pump will be removed from the container and rinsed with deionized water.

- 17) Measure the well depth; record observation in field logbook.
- 18) Re-lock well cap, if present.
- 19) Fill out field logbook, sample labels, custody seals, and chain of custody forms.

### **3.3 Decontamination**

As detailed in the previous sections, all stainless steel equipment involved in field sampling activities will be decontaminated prior to and subsequent to sampling. Decontamination of sampling equipment will be kept to a minimum in the field and whenever possible dedicated sampling equipment will be used. Decontamination of sampling equipment including stainless steel scoops, trowels, and bowls will be conducted as follows:

- 1) Alconox detergent and potable water scrub,
- 2) Potable water rinse,
- 3) A 10% nitric acid rinse (ultra pure grade) when sampling for inorganic parameters,
- 4) Distilled or potable water rinse,
- 5) An acetone *only* rinse or a methanol rinse followed by a hexane rinse (pesticide grade or better),
- 6) Deionized water rinse,
- 7) Air dry (sufficient time will be allowed for the equipment to completely dry), and
- 8) Wrap or cover exposed ends of sampling equipment with aluminum foil (shiny side out) for transport and handling.

## **4.0 QUALITY ASSURANCE/QUALITY CONTROL**

This section details the Quality Assurance/Quality Control (QA/QC) requirements for field activities performed during the sampling effort.

### **4.1 Field Instrument Calibration and Preventive Maintenance**

The sampling team is responsible for assuring that a master calibration/maintenance log will be brought into the field and maintained for each measuring device. Each log will include at a minimum, where applicable:

- name of device and/or instrument calibrated
- device/instrument serial and/or ID number
- frequency of calibration
- date of calibration
- results of calibration
- name of person performing the calibration
- identification of the calibrant (PID, FID, pH meter)

Equipment to be used each day shall be calibrated prior to the commencement of daily activities.

### **4.2 QA/QC Sample Collection**

#### **4.2.1 Trip Blanks**

A trip blank will be collected for events involving aqueous sampling for volatile organic parameters. A trip blank is an aliquot of deionized (DI), demonstrated analyte-free water which is prepared in the field prior to the initiation of field work and sealed in 40-ml glass vials with teflon-lined septum caps. Analytical results of the trip blank sample are utilized during sample data validation to determine if any cross contamination has occurred between samples during shipment/storage, or if on-site atmospheric contaminants are seeping into the sample vials. These sealed bottles will be placed in a plastic cooler and will accompany field personnel to the sampling locations.

Trip blanks will be collected in accordance with the procedures listed below:

- 1) Pour DI water into an extra vial. Determine the amount of 1:1 HCl required to lower sample pH to less than 2, verifying with broad range pH paper. Add the predetermined volume of HCL to the two sample vials.
- 2) Proceed to fill the two vials with DI water just to overflowing and seal so that no air bubbles are trapped inside. Place in sample cooler. After the sampling event is completed, the trip blank sample will be packaged and hand-delivered or shipped



within 24 hours of collection, along with the environmental samples, to the designated laboratory.

- 3) Complete sample labels, custody seals, and chain of custody forms. Record in field logbook.

#### **4.2.2 Field Rinsate Blanks**

A field rinsate blank will consist of DI, demonstrated analyte-free water that has been poured over decontaminated sampling equipment. The field rinsate blank analytical results will be utilized in evaluation of potential cross contamination resulting from inadequate decontamination. The frequency of field rinsate blank collection is one blank per decontamination event per type of equipment, not to exceed more than one per day. For the purposes of sampling associated with pre-remedial field activities, field rinsate blank collection will not exceed a total of three samples. Blanks will be collected for all parameters of interest (excluding physical parameters) and shipped with the samples collected the same day.

Field rinsate blanks will be collected in accordance with the procedure listed below:

- 1) Decontaminate sampling equipment using the procedure specified in Section 3.4 of this plan.
- 2) Pour DI water over the sampling device and collect the rinsate in the appropriate sample containers. Collect the VOA vials first and preserve in the same manner as trip blank vials.
- 3) Preserve remaining samples as specified in Table 2-1 of this plan. Test pH by pouring a small portion of sample on broad range pH paper over a collection bowl. Place samples in cooler.
- 4) Complete sample labels, custody seals, and chain of custody forms. Record in field logbook.

#### **4.2.3 Deionized Water Blanks**

The distilled deionized (DI) water utilized for the trip and field blanks will be certified as such. A copy of this certificate will be kept on site and another in the site-specific project file. The criteria to be demonstrated as analyte-free will be consistent with that specified in the U.S. EPA Region II CERCLA Quality Assurance Manual (October 1989), and is as follows:

Purgeable organics < 10 ppb  
Semi-volatile organics < CRQL  
Pesticides/PCBs < CRQL

## Inorganics < CRDL

where the CRQL is represented by the Contract Required Quantitation Limit and the CRDL is represented by the Contract Required Detection Limit in the most recent CLP Statement of Work. For specific common laboratory contaminants such as methylene chloride, acetone, toluene, 2-butanone and phthalates, the allowable limits are three times the respective CRQLs.

### **4.2.4 Duplicate Samples**

Duplicate samples will be sent for laboratory analysis to evaluate the ability of reproducing the sampling methods. At a minimum a rate of one duplicate sample per 20 samples, or one duplicate sample per batch of less than 20 samples, will be obtained for each matrix. For the purpose of pre-remedial site assessment projects, soil and sediment matrices will be considered as same matrix.

In addition, a minimum of one matrix spike/matrix spike duplicate (MS/MSD) sample per matrix will be collected per 20 samples, or one MS/MSD sample per matrix per batch of less than 20 samples. The analysis of CLP TCL parameters for the MS/MSD sample will involve the collection of triple sample volume for aqueous samples only. Extra sample volumes are not required for TCL solid matrix (i.e., soil/sediment matrix) and TAL parameters MS/MSD analyses.

### **4.2.5 Data Validation**

Personnel, trained and approved by U.S. EPA Region II Monitoring Management Branch, will perform all data validation utilizing the most current U.S. EPA Region II data validation guidelines.

## **5.0 FIELD CHANGES AND CORRECTIVE ACTIONS**

The SAT Project Manager (PM) or his/her designee may be required to modify generic site procedures to accommodate site-specific needs or unforeseeable events. In the event it becomes necessary to modify a procedure, the PM will notify the U.S. EPA Region II WAM. Deviations from the field sampling plan are to be documented in the field logbook and signed by the initiator and the PM.

REFERENCE NO. 3



UNITED STATES ENVIRONMENTAL PROTECTION AGENCY

REGION 2

290 BROADWAY

NEW YORK, NY 10007-1866

AUG 13 1999

**ACTION MEMORANDUM**

**DATE:**

**SUBJECT:** Confirmation of Verbal Authorization of Funding and Request for Ceiling Increase at the Lackawanna Foundry Site, Lackawanna, New York

**FROM:** Kevin M. Matheis, On-Scene Coordinator  
Removal Action Branch

**TO:** Richard L. Caspe, Director  
Emergency and Remedial Response Division

**THRU:** Richard C. Salkie, Chief  
Removal Action Branch

**Site ID#:** MY

**I. PURPOSE**

The purpose of this Action Memorandum is to request and document approval of the proposed removal action described herein for the Lackawanna Foundry Site (Site), located at 2 Elm Street, Lackawanna, Erie County, New York 14218. On March 19, 1999 the Acting Division Director of the Emergency and Remedial Response Division granted verbal authorization for the U.S. Environmental Protection Agency (EPA) to undertake emergency stabilization actions at the Site, authorizing a total project ceiling of \$100,000 (\$90,000 in contractor mitigation costs and \$10,000 in EPA costs) to initiate removal activities to secure and stabilize the Site. This stabilization action was initiated on March 22, 1999.

The actions described herein seek to complete the removal actions at the Site by removing drums, cylinders, PCB contaminated soils and equipment and arsenic contaminated soils at the Site. The proposed project ceiling for this action will be \$1,500,000, of which \$930,000 is for mitigation contracting.

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**CONCURRENCES**

Name: Morgan Chem.

Init: sb

Date: 08/04/99

Filename: A.M.#0166

| Symbol  | ERRD-RAB | ERRD-RAB | ERRD-RAB | ORC-NYCSUP | ERRD-DD | ERRD-D  |
|---------|----------|----------|----------|------------|---------|---------|
| Surname | Matheis  | Rotolo   | Salkie   | Simon      | McCabe  | Caspe   |
| Date    | 8/11/99  | 8/11/99  | 8/11/99  | 8/13/99    | 8/13/99 | 8/13/99 |

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The actions described herein seek to complete the removal actions at the Site by removing drums, cylinders, PCB contaminated soils and equipment and arsenic contaminated soils at the Site. The proposed project ceiling for this action will be \$1,500,000, of which \$930,000 is for mitigation contracting.

The Site is not on the National Priorities List (NPL). There are no nationally significant or precedent-setting issues associated with the proposed removal action.

## **II. SITE CONDITIONS AND BACKGROUND**

The Comprehensive Environmental Response, Compensation and Liability Information System ID Number for the Site is pending assignment through the application submitted by EPA. Once assigned, the ID will accompany all appropriate paperwork regarding the Site.

### **A. Site Description**

#### **1. Removal site evaluation (RSE)**

On December 22, 1998, the New York State Department of Environmental Conservation (NYSDEC) requested that EPA evaluate the Site and perform a removal action, as appropriate, to address the threats at the Site from drums, PCB transformers, and PCB contaminated soils. Included in the referral was an drum inventory performed by the site owner that showed the presence of over 223 drums of various foundry raw materials and waste. This inventory showed 25 drums or binders, oils and resins that were suspected flammable or combustible materials, and 190 drums of slag waste, unknown materials, and foundry sands. One drum of reactive material, calcium carbide, was also listed. Also included in the referral, analysis of transformers indicated that two transformers contain oil with PCBs at concentrations as high as 86 percent and adjacent soils with PCBs as high as 210 parts per million (ppm). On March 16, 1999, EPA performed a RSE. During the RSE, it was observed that the Site had no fencing or security, the buildings were open and evidence of trespassing were quite evident. Graffiti was present at the outer and interior buildings on-site and evidence of splat-ball games were prevalent throughout the building.

It was apparent that the inventory provided by NYSDEC underestimated the amount of drums discovered at the Site. EPA identified over 100 unknown drums outside the foundry buildings and approximately 100 drums, laboratory chemicals, compressed gas cylinders and PCB transformers were noticed within the foundry buildings.

Based upon observations during the RSE and analysis provided by the NYSDEC, verbal authorization of funding was granted by the acting Division Director on March 19, 1999. This funding authorized stabilization actions to commence at the Site. These actions involved fencing the perimeter of the Site and securing the buildings, moving drums into a secure location and stabilizing leaking containers.

During the emergency stabilization operations, EPA's contractor sampled drums and collected label information from containers to identify hazardous substances and to characterize the waste for segregation purposes. During the field characterization, four drums were found to be D001 flammable, 18 drums were D001 oxidizers, one drum was D003 water reactive, five drums were D002 corrosive and the remaining 116 drums contain various solids and chemicals that have yet to be characterized.

In addition, 53 five-gallon and 63 one-gallon containers of paint-related materials were discovered and inventoried. From label information collected, the containers contain various concentrations of xylene, acetone, toluene and naphtha. Many of the paint-related materials indicate they are flammable.

Over 1,000 pounds of laboratory chemicals were inventoried and stabilized during the initial action. Label information from these containers indicate that hazardous substances and wastes such as sulfuric acid, phosphoric acid, flammable alcohols and acetone, ammonium hydroxide, calcium carbide and many unknown materials are contained within these laboratory chemicals. These containers are in poor condition and many have leaked their contents into the site buildings.

Compressed gas cylinders were collected from throughout the property and staged according to compatibility. The following were inventoried: four cylinders contain liquified petroleum gas, a flammable gas; one cylinder contains propane, a flammable gas; one cylinder contains oxygen, an oxidizer in compressed state; one cylinder contains, polymeric isocyanate, a poisonous gas; and one cylinder contains unknown materials.

Key problem areas that exist on-site include leaking drums of unknown solids and liquids, soil contaminated with high levels of PCBs in areas of PCB capacitor dumping, unknown and known cylinders in poor condition, laboratory chemicals stored in poor containers, drums of liquids in poor condition that contain corrosive and flammable liquids, arsenic contaminated soils in an adjacent recreational area and a building that is in an advanced stage of deterioration.

In April 1999, EPA's Environmental Response Team (ERT) performed a preliminary assessment (PA) at the Site to ascertain the extent of surface soil and subsurface soil contamination outside and inside the buildings on-site. The site soil sampling results showed elevated levels of arsenic contamination in a recreational area (66ppm, 48ppm, and 77ppm). In addition, in the areas where the PCB capacitors were dumped, soil contamination for PCBs was found at 256,000 ppm.



Due to the above mentioned conditions, the Site continues to pose a threat to public health and the environment, and thereby qualifies for a removal action.

The site referral from NYSDEC is included as Attachment 1.

## **2. Physical location**

The Site is located in a mixed residential and heavily industrialized area of Lackawanna, New York, adjacent to Smokes Creek, a direct tributary to Lake Erie. Twelve residential properties are immediately adjacent to the Site to the west, railroad tracks border the east, Smokes Creek to the north and a multi-use city park to the south.

The residential neighborhood immediately adjacent to the Site consists of twelve residences with over 100 residences being located within 1/8 mile of the Site. The population of the area within a 1/2 mile of the Site consists of nearly 1,000 people.

## **3. Site characteristics**

The Site has been the location of a cast-iron foundry since the early 1900's. The current owner of the Site inherited the business from his father who operated the foundry at this location for more than 30 years. Residences adjacent to the Site have told EPA that foundry operations have been small-scale and have steadily decreased over the past 20 years. Over the past five years, the current owner has operated a marginal business and has ceased operations completely since approximately 1997. To date, two primary areas of contamination have been identified and are described below:

On-Site Buildings - PCBs located at the Site are suspected to have originated from electrical equipment and machinery operated at the foundry. Items containing PCBs have deteriorated over time and leaked, which resulted in the contamination of soils within the buildings. Sampling from PCB capacitors dumped inside of buildings showed the presence of 100 percent PCBs. Several of these capacitors have leaked inside the on-site buildings and samples showed the soil contaminated with 25 percent PCBs. The removal of these capacitors and associated soil contamination will be part of this proposed removal action. In addition to the high concentrations of PCBs found in the areas of PCB capacitor storage and dumping, several areas within the on-site buildings are suspected to contain hazardous substances and will need to be sampled. A basement area containing runoff water and sediment from the Site will be sampled and evaluated as part of this removal. The drums, laboratory chemicals and cylinders appear to have been utilized as part of the foundry operations and have been abandoned since the operations have ceased.

Recreational Area - The recreational area adjacent to the Site is approximately two acres in size and is owned by the City of Lackawanna. A bocce court and baseball field is on this parcel and is frequented by citizens who live adjacent to the Site and in the surrounding Bethlehem Park community. The recreational area is a grassy area that Lackawanna maintains by frequent grass and weed cutting and trimming. A wide group of citizens of various age groups use the recreational area on a daily basis. The Site is topographically higher than the recreational area. The arsenic contamination adjacent to the Site in the recreational area appears to be in the location of a bag-house dust collection system used during foundry operations. The contamination was caused by either airborne deposition or runoff from the baghouse collection area. Of the arsenic data collected on-site, the highest concentrations of arsenic contaminated soils have been identified in this area.

The actions proposed in this memoranda will be a continuation of site stabilization and cleanup activities. Presently, EPA inspects the Site on a routine basis and maintains a command post and response capabilities in the event of an emergency.

**4. Release or threatened release into the environment of a hazardous substance, or pollutant, or contaminant.**

The inventory obtained by EPA as part of the RSE included the following hazardous substances as defined by Section 101(33) of the Comprehensive Environmental Response, Compensation and Liability Act (CERCLA), as amended, 42 U.S.C. 9601 et seq. and listed in 40 CFR Table 302.1:

| Material                          | Quantity             | Storage Method              | Primary Hazard  |
|-----------------------------------|----------------------|-----------------------------|---|
| PCB contaminated soils and debris | 1,000 y <sup>3</sup> | Soils inside building       | toxic, carcinogen<br>25.6% PCBs   |
| PCB equipment                     | 120 units            | Transformers and capacitors | toxic, carcinogen<br>100% Pure PCBs   |
| Arsenic contaminated soils        | 800 y <sup>3</sup>   | Soils outside buildings     | toxic, carcinogen   |
| Laboratory Chemicals              | 1,000 lbs            | Various sized containers    | D001 - flammables,<br>D002 - corrosives,<br>D001 - oxidizers,<br>D003 - reactives |
| Compressed gas cylinders          | 9                    | Cylinders                   | D001 - flammables,<br>poisons, D001 -<br>oxidizers,<br>uncharacterized            |

|  |                                |   |   |
|--|--------------------------------|---|---|
| Drums, known and yet to be characterized | 144 x 55 gallon drums          | Interior storage buildings, poor condition  | D001 - flammables, D002 - corrosives, D003 - reactives, and uncharacterized |
| Five gallon pails and paints             | 53 x 5 gallon<br>63 x 1 gallon | Interior storage buildings, poor conditions | D001 - flammables, D002 - corrosives and uncharacterized                    |

Sampling results and a map depicting sample locations are included in Attachment 2.

Several release mechanisms exist at the Site for the various hazardous substances found. These release mechanisms include the following:

- There is a fire and explosion threat from potential releases of the drums containing flammable, reactive or oxidizing materials. Resulting emissions from a release in conjunction with the amount of uncharacterized materials contained within drums on-site would cause a significant threat to the adjacent residential area. Since the Site has been historically the subject of frequent vandalism and break-ins, the threat of release from intentional causes is high. The drums known to be reactive and/or flammables and oxidizers were identified through thorough on-site field characterization.
- The PCB equipment and highly contaminated PCB soils on-site present a threat of release from poor building conditions that may cause contaminants to migrate. In the area where the PCB capacitors were found dumped, water exists on top of the soil contamination areas. Due to the poor condition of the roof, rainfall events may cause the contamination to migrate further. In addition, the PCB transformers and capacitors are in poor condition and several items have already leaked. Actions are required to remove this equipment off-site to disposal facilities. The concentrations of the PCB capacitors dumped on-site consist of PCBs in the 100 percent pure category. Samples of transformers in the referral by NYSDEC indicated that several transformers contain 86 percent PCBs. Soil samples taken from the location of PCB capacitors leaking on-site indicate the presence of PCBs at 25 percent. Summary of PCB results follows:

| Capacitor Sample Results  | Transformer Sample Results | Soil Contamination Results |
|---------------------------|----------------------------|----------------------------|
| 1,000,000 ppm (100%) PCBs | 866,000 ppm (86%) PCBs     | 250,000 ppm (25%) PCBs     |

- In the area of arsenic soil contamination, the contaminated soil is unsecured and adjacent to a recreational area frequented by citizens of all ages utilizing the property for multi-purposes. The runoff area from the bag house collection system into the recreational area appear to have areas of soil exposed at the surface where vegetation is non-existent. This area would be a prime pathway for exposure to the contaminated soil and for arsenic

contaminated soils to further migrate. The recreational area contains a bocce field and baseball area and is fully grassed, being maintained by the City of Lackawanna. The area is frequented on a daily basis by citizens utilizing this open green space in the Bethlehem Park community. The arsenic contaminated soils, in concentrations ranging from 21-77 ppm arsenic in surface and subsurface soils abuts the foundry buildings in bulldozed piles and weather conditions such as rainfall may cause the contamination to spread further. Actions are needed to remove the arsenic contaminated soils from this area. Sample results from the site adjacent to the residential area follows:

| Sample 22 (0 - 12") Arsenic Results | Sample 26 (0 - 12") Arsenic Results | Sample #26D (1' -3' depth) Arsenic Results |
|-------------------------------------|-------------------------------------|--|
| 77 ppm                              | 66 ppm                              | 48 ppm                                     |

All three samples taken in this area exhibited elevated arsenic concentration levels. Other arsenic samples taken away from the bag house dust collection area decreased in arsenic concentrations. These results show a definable area of arsenic concentration that can be remediated as part of this removal action.

This removal action addresses the disposal of these hazardous materials from the Site.

**5. NPL status**

The Site is not listed or proposed to be included on the NPL. The site assessment program will be contacted to provide an evaluation of NPL inclusion as further site information is collected and evaluated.

**6. Maps, pictures, and other graphic representations**

Figures 1, 2 and 3 are included as Attachment 3 and provide the location and configuration of the Site.

**B. Other Actions to Date**

**1. Previous actions**

In 1997, the facility was marginally operational and the NYSDEC had performed an inspection at the Site which identified drums and transformers. At that time, they required the owner of the Site to undertake an inventory of all containers on-site and to take samples and analysis of transformers and suspected areas of PCB soil contamination. Results of the sampling indicated that the transformers contained up to 86 percent PCBs and soil contamination from the

transformers were up to 210 ppm of PCBs. As a result, the NYSDEC demanded that the owner take actions to address the threats at his property. The owner has been unable financially to perform a cleanup and has abandoned the property.

Since the site buildings are in such poor condition, the City of Lackawanna has also ordered the owner of the property to clean the debris from the property and to demolish several of the structures on-site due to poor structural integrity. The owner has not complied with these requests and has communicated with city officials that they can do what is necessary to bring the building into compliance with city ordinances.

## **2. Current actions**

EPA mobilized to the Site on March 22, 1999 to begin site stabilization operations. During these operations, EPA fenced the perimeter of the buildings and the adjacent residential area. In addition, all drums stored outside the buildings were moved inside, leaking liquid drums were over packed and all drums were sampled and field characterized. PCB transformers were secured and all laboratory chemicals, five-gallon pails and paint cans were organized and segregated. Cylinders of compressed gas were collected from throughout the Site and consolidated into a secure storage area. Over 100 PCB capacitors were discovered in a hidden trench on-site and were excavated and placed into steel drums. The trench was sampled and found to contain over 26 percent PCBs in the soil. This area was roped with caution tape pending excavation and removal to off-site disposal facilities.

ERT mobilized to the Site in April 1999 and collected surface and subsurface soil samples from the interior of the building and in areas adjacent to the building. Sampling results indicated that arsenic contamination had migrated to the recreation area located to the south of the Site.

Erie County has identified the recreational area and Site property that borders Smokes Creek as a target for environmental restoration under the 1996 New York Clean Water/Clean Air Bond Act. Upon completion of EPA's actions at the Site, the area may be turned into a recreational area with waterfront access.

EPA has informed the community about the status and the future of the removal action at the Site. Fact sheets have been delivered and personal interviews have been held with residents immediately adjacent to the Site. EPA intends to expand community relations actions when additional removal actions commence.

## **C. State and Local Authorities' Role**

### **1. State and local actions to date**

On December 22, 1998, the NYSDEC requested that EPA evaluate the Site and perform a removal action, as appropriate, to address the threats at the Site from drums, PCB transformers and PCB contaminated soils.

The City of Lackawanna and EPA continue to monitor the Site for breaches in security and vandalism.

### **2. Potential for continued State/local response**

There are no actions being taken by State or local government agencies to address the hazardous substances located at the Site.

Erie County has identified the recreational area and Site property that borders Smokes Creek as a target for environmental restoration under the 1996 New York Clean Water/Clean Air Bond Act. Upon completion of EPA's actions at the Site, the area may be turned into a recreational area with waterfront access.

## **III. THREATS TO PUBLIC HEALTH, OR WELFARE, OR THE ENVIRONMENT AND STATUTORY AND REGULATORY AUTHORITIES**

The following factors described in 40 CFR Part 300.415(b)(2) of the National Contingency Plan (NCP) were applied in determining the appropriateness of a removal action at the Site.

- (i) Actual or potential exposure to nearby human populations, animals, or the food chain from hazardous substances, or pollutants, or contaminants;
- (iii) Hazardous substances, or pollutants, or contaminants in drums, barrels, tanks, or other bulk storage containers, that may pose a threat of release;
- (iv) High levels of hazardous substances, or pollutants, or contaminants in soils largely at or near the surface, that may migrate;
- (v) Weather conditions that may cause hazardous substances, pollutants, or contaminants to migrate or be released;
- (vi) Threat of fire or explosion; and
- (vii) The availability of other appropriate federal or State response mechanisms to respond to the release.

**A. Threats to Public Health or Welfare**

- (i) Actual or potential exposure to nearby human populations, animals, or the food chain from hazardous substances, or pollutants, or contaminants;**

The arsenic contaminated soils in the recreational area is unsecured and presents a threat to persons that comes in contact with the contaminated surface soils. The site buildings contain drums of hazardous substances such as PCBs, xylene, toluene, flammable liquids, corrosive liquids and poisons stored in containers that are in poor condition and not intended to be stored indefinitely. Many of these containers have leaked and present a threat of exposure. The PCB contaminated soils contained within the buildings may migrate outside the buildings and cause a direct contact threat.

- (iii) Hazardous substances, or pollutants, or contaminants in drums, barrels, tanks, or other bulk storage containers, that may pose a threat of release;**

The drums found at the Site contain the hazardous substances as listed in Section II.A.4 of this Action Memorandum. The hazardous substances contained within these containers present a threat of continuing release.

- (iv) High levels of hazardous substances, or pollutants, or contaminants in soils largely at or near the surface, that may migrate;**

The site perimeter contains arsenic contaminated soils that is unsecured and impacting a recreational area next to the Site. The location of the arsenic soils is topographically higher than the surrounding area and the hazardous substances will continue to migrate off-site if not remediated. In addition, high levels of PCB contaminated soils exist in the warehouse on-site and are covered by rainwater infiltration. If the PCB contaminated soil is not removed, the contamination will continue to migrate, perhaps beyond the building perimeter walls.

- (v) Weather conditions that may cause hazardous substances, pollutants, or contaminants to migrate or be released;**

The building is in poor structural condition with leaking roof and walls and partially collapsed areas. The Site has no utilities and temperature controls and drums and other containers are subject to freeze/thaw cycles. Rainfall causes flooding in several areas of the buildings and contamination from the PCB soils may continue to migrate.

- (vi) Threat of fire or explosion; and**

The Site contains flammable liquids contained in drums, five-gallon pails, paint cans and laboratory chemicals. In addition, several drums contain reactive materials. Due to the poor conditions of the buildings and the threat of vandalism and chemical incompatibility, a threat of

fire and explosion exists on-site. The Site is immediately adjacent to a residential area and a resulting fire would require the evacuation of the adjacent residential area.

- (vii) The availability of other appropriate federal or State response mechanisms to respond to the release.**

No other federal or State response mechanism is available to respond to the significant threat to the environment that the Site presents.

**B. Threats to the Environment**

- (i) Actual or potential exposure to nearby human populations, animals, or the food chain from hazardous substances, or pollutants, or contaminants;**

The arsenic contaminated soils in the recreational area is unsecured and presents a threat to animals that comes in contact with the contaminated surface soils. The site buildings contain drums of hazardous substances such as PCBs, xylene, toluene, flammable liquids, corrosive liquids and poisons stored in containers that are in poor condition and not intended to be stored indefinitely. Many of these containers have leaked and present a threat of exposure. The PCB contaminated soils contained within the buildings may migrate outside the buildings and cause a direct contact threat. The Site is immediately adjacent to Smokes Creek, a direct tributary to Lake Erie. Fishing is very prevalent in the creek and fish can bioaccumulate PCBs from the Site. This bioaccumulation would cause damage to the food chain as a result.

- (iii) Hazardous substances, or pollutants, or contaminants in drums, barrels, tanks, or other bulk storage containers, that may pose a threat of release;**

The drums found at the Site contain the hazardous substances as listed in Section II.A.4 of this Action Memorandum. The hazardous substances contained within these containers present a threat of continuing release to the environment.

- (iv) High levels of hazardous substances, or pollutants, or contaminants in soils largely at or near the surface, that may migrate;**

The Site perimeter contains arsenic contaminated soils that is unsecured and impacting a recreational area next to the Site. The location of the arsenic soils is topographically higher than the surrounding area and the hazardous substances will continue to migrate off-site if not remediated. In addition, high levels of PCB contaminated soils exist in the warehouse on-site and are covered by rainwater infiltration. If the PCB contaminated soil is not removed, the contamination will continue to migrate, perhaps beyond the building perimeter walls. The Site is immediately adjacent to Smokes Creek, a direct tributary of Lake Erie. Run-off from the Site would directly enter the creek, causing contamination to migrate off-site.



- (v) **Weather conditions that may cause hazardous substances, pollutants, or contaminants to migrate or be released;**

The building is in poor structural condition with leaking roof and walls and partially collapsed areas. The Site has no utilities and temperature controls and drums and other containers are subject to freeze/thaw cycles. Rainfall causes flooding in several areas of the buildings and contamination from the PCB soils may continue to migrate.

- (vi) **Threat of fire or explosion; and**

The Site contains flammable liquids contained in drums, five-gallon pails, paint cans and laboratory chemicals. In addition, several drums contain reactive materials. Due to the poor conditions of the buildings and the threat of vandalism and chemical incompatibility, a threat of fire and explosion exists on-site. Firefighting water from the resulting fire may enter the adjacent creek causing contamination and hazardous substances to further migrate.

- (vii) **The availability of other appropriate federal or State response mechanisms to respond to the release.**

No other federal or State response mechanism is available to respond to the significant threat to the environment that the Site presents.

#### **IV. ENDANGERMENT DETERMINATION**

Actual or threatened releases of hazardous substances from the Site, if not addressed by implementing the response action selected in this Action Memorandum, may present an imminent and substantial endangerment to public health, or welfare, or the environment.

#### **V. PROPOSED ACTIONS AND ESTIMATED COST**

##### **A. Proposed Actions**

##### **1. Proposed action description**

The following activities are proposed to address the immediate threats to human health and the environment posed by hazardous substances, or pollutants, or contaminants present at the Site.

- a. Stabilization - All containers that are open or of questionable integrity will be over packed or transferred into a new container.
- b. Sampling - All hazardous substances, pollutants and contaminants in containers, soil and debris will be sampled prior to disposal. Where possible, composite samples will be taken to reduce the total amount of samples analyzed. In order to quantify and delineate the

extent of soil contamination, additional samples will be taken in areas of arsenic and PCB soil contamination prior to excavation.

- c. Analysis - All samples will be evaluated for compatibility. The samples will be analyzed for disposal parameters, anticipated to be full toxicity characteristic leaching procedure (TCLP) analysis. The analysis will be sent to disposal facilities in compliance with EPA's off-site disposal rule to ensure proper disposal.
- d. Excavation - Upon receipt of approvals at disposal facilities, the areas of arsenic and PCB soil contamination will be excavated and sent to off-site disposal facilities. For the arsenic contaminated soils, a cleanup level of 20 ppm will be utilized. For PCB contaminated soils, a cleanup level of ten ppm will be implemented. Delineation of the contamination will be performed prior to excavation. Based upon existing data, the volume estimate for soil to be removed in this Action Memorandum should be similar to the area identified in the proposed delineation. Excavations will occur in areas that are adjacent to support structures of the steel framed building and partial demolition of the building may be required. The building is constructed of structural steel and sheet metal so the fabric and structure of the building will be recycled as scrap, if demolition is necessary.
- e. Disposal - Upon receipt of disposal analysis, waste profiles will be completed and sent to disposal facilities for acceptance. Hazardous substance, pollutants and contaminants will be sent to off-site disposal facilities in compliance with EPA's Off-Site Disposal Rule.
- f. Restoration - Areas of the recreational area impacted by the excavation of arsenic contaminated soils will be returned to original condition by replacing soil with certified clean fill and re-seeding the grassy areas. In addition, any access road necessary to excavate this area will be removed.

## 2. Contribution to remedial performance

The removal action at the Site is consistent with the requirement of Section 104(a)(2) of CERCLA, which states, "any removal action undertaken . . . should . . . to the extent practicable, contribute to the efficient performance of any long-term remedial action with respect to the release or the threatened release concerned." Since any remedial action undertaken would encompass the elements in this response, this removal action is consistent with any future remedial work.

## 3. Description of alternative technologies

Because of the quantities and types of the hazardous substances and/or wastes at the Site, on-site treatment and/or incineration is not appropriate. The selected removal action includes the characterization of hazardous substances found at the Site and the transportation of all hazardous substances off-site for treatment and/or disposal. The selected removal action has been

determined to be the appropriate response action for the Site based upon the criteria of effectiveness, implement ability and cost.

**4. EE/CA**

Due to the time-critical nature of this removal action, an EE/CA will not be prepared.

**5. Applicable or Relevant and Appropriate Requirements (ARARS)**

ARARS that are within the scope of this removal action will be met to the extent practicable. Federal ARARS determined to be applicable for the proposed scope of work include the Resource Conservation and Recovery Act, Occupational Safety and Health Act and Hazardous Materials Transportation Uniform Safety Act.

**6. Project schedule**

It is anticipated that the project will be completed within three months. Upon mobilization to the Site, all samples will be collected for disposal approvals. Upon receipt of disposal approvals, the removal of hazardous substances will commence.

**Estimated Costs:**

The estimated costs for the completion of this project are summarized below. A summary of costs is included as Attachment 4.

**Extramural Costs:**

|  | <u>Previous<br/>Ceiling</u> | <u>Increase<br/>Requested</u> | <u>Total<br/>Ceiling</u> |
|--|-----------------------------|-------------------------------|--------------------------|
| <b><u>Regional Allowance Costs:</u></b>  |                             |                               |                          |
| Total ERCS Cost  | \$ 90,000                   | \$ 840,000                    | \$ 930,000               |
| 20% contingency  |                             | \$ 168,000                    | \$ 168,000               |
| <b><u>Other Extramural Costs Not Funded From<br/>the Regional Allowance:</u></b> |                             |                               |                          |
| Total START costs  | \$ 2,000                    | \$ 128,000                    | \$ 130,000               |
| <b>TOTAL, EXTRAMURAL COSTS</b>   | <b>\$ 92,000</b>            | <b>\$ 1,136,000</b>           | <b>\$ 1,228,000</b>      |

**Intramural Costs:**

|                                       |                  |                     |                     |
|---------------------------------------|------------------|---------------------|---------------------|
| Intramural Direct Costs               | \$ 3,000         | \$ 88,000           | \$ 91,000           |
| Intramural Indirect Costs             | <u>\$ 5,000</u>  | <u>\$ 176,000</u>   | <u>\$ 181,000</u>   |
| <b>TOTAL, INTRAMURAL COSTS</b>        | <b>\$ 8,000</b>  | <b>\$ 264,000</b>   | <b>\$ 272,000</b>   |
| <b>TOTAL, REMOVAL PROJECT CEILING</b> | <b>\$100,000</b> | <b>\$ 1,400,000</b> | <b>\$ 1,500,000</b> |

**VI. EXPECTED CHANGE IN THE SITUATION SHOULD NO ACTION BE TAKEN OR ACTION DELAYED**

Should no action be taken or the planned action be delayed, hazardous substances in drums or contaminated soils could be released. A release of hazardous substances from these containers and soil could result in the exposure of the neighboring population and/or contamination of the environment to the adjacent creek. Releases of contaminants to the ground and/or groundwater would increase the cost of the required removal action.

**VII. OUTSTANDING POLICY ISSUES**

No known outstanding policy issues are associated with the Site.

**VIII. ENFORCEMENT**

EPA has a signed access agreement from the property owner and has received information on his financial ability to undertake a removal action. At this time it appears that the owner does not have the ability to undertake an emergency removal action at the Site. EPA will continue to evaluate the ability of the owner to undertake a removal action at various times during the removal, but expects to pursue cost-recovery as a means to reimburse actions undertaken by EPA.

**IX. RECOMMENDATION**

This decision document represents the selected removal action for the Lackawanna Foundry Site located in Lackawanna, New York, developed in accordance with CERCLA, as amended, and not inconsistent with the NCP. This decision is based on the administrative record for the Site.

Conditions at the Site meet the NCP Section 300.415(b)(2) criteria for a removal and I recommend your approval of the proposed removal action. The total project ceiling, including actions from the verbal authorization of funding if approved, will be \$1,500,000. Of this, \$90,000

has already been allocated to the project and an estimated \$840,000 additional will come from the Regional removal allowance. Funds for this removal action are currently within the Regional Advice of Allowance.

Please indicate your approval, or disapproval, and authorization of funding as per current Delegation of Authority, by signing below.

**APPROVAL:** \_\_\_\_\_ **DATE:** \_\_\_\_\_

Richard L. Caspe, Director  
Emergency and Remedial Response Division

**DISAPPROVAL:** \_\_\_\_\_ **DATE:** \_\_\_\_\_

Richard L. Caspe, Director  
Emergency and Remedial Response Division

cc: (after approval)  
W. Muszynski, DRA  
R. Caspe, ERRD-D  
W. McCabe, ERRD-DD  
R. Salkie, ERRD-RAB  
J. Rotola, ERRD-RAB  
G. Zachos, ACSMO  
B. Dease, ERRD-RPB  
B. Bellow, EPD  
P. Simon, ORC-NYCSUP  
K. Weaver, OPM-FIN  
R. Gherardi, OPM-FIN  
C. Moyik, ERRD-PS  
T. Johnson, 5202G  
M. O'Toole, NYSDEC  
T. Vickerson, NYSDEC  
P. McKechnie, OIG  
A. Raddant, DOI  
G. Wheaton, NOAA  
O. Douglas, START

# Attachment 1

**New York State Department of Environmental Conservation**  
**Division of Environmental Remediation, Room 260B**  
50 Wolf Road, Albany, New York 12233-7010  
Phone: (518) 457-5861 FAX: (518) 485-8404

FJV:ts  
cc: M. O'Toole (2)  
H. Koelling  
P. Buechi - NYSDEC, Region 9  
T. Vickerson  
Dayfile  
tjv.discl:hightech1.wpd

RECEIVED

DEC 22 1998

NYSDEC - REG. 9  
FOIL  
\*REL UNREL

f  
High Tech Metals  
Co. Co. (unlisted)

Mr. Richard Caspe  
Director  
Emergency & Remedial Response Division  
USEPA, Region II  
290 Broadway  
New York, New York 10007-1866

Dear Mr. Caspe:

Re: High Tech Metals Site  
Lackawanna (C), Erie County  
Request for Emergency Removal

The New York State Department of Environmental Conservation (NYSDEC) hereby requests the United States Environmental Protection Agency (USEPA) to perform an appropriate CERCLA/SARA authorized emergency response action at the former High Tech Metals site (a.k.a. Lackawanna Foundry).

The site consists of a several story block, brick and wood foundry building in a mixed residential/industrial area off Elm and Jackson Streets in the City of Lackawanna. A residential area is located adjacent to the west of this site and surface water drainage is mostly to Smokes Creek less than 500 feet north of the site.

Previous site investigations by NYSDEC staff and a limited sampling program by the owner have reported accumulations of numerous pails, drums and other containers of wastes from past site operations. An inventory of these drums has been completed but contents were not verified by sampling. Additionally, and of most concern are eight transformers located in the building. A total of five samples of transformer oil were collected. All five samples contained PCB contamination from 12 ppm to 860,000 ppm. Surface soils surrounding the transformer areas have been sampled. PCB contamination was noted as high as 210 ppm from a composite sample. Evidence of past releases have been witnessed and documented by NYSDEC staff during site visits. The current property owner has not been responsive to requests for further action from the Department and has since vacated the premises and has not paid taxes on the facility for a number of quarters.

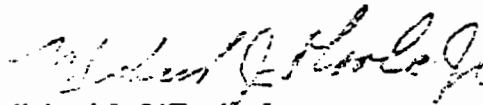
Due to location of the facility adjacent to a residential area and to Smokes Creek (a tributary of Lake Erie) and indications that a release(s) has/have been noted (i.e., visible stains originating from containers, deteriorating/bulging containers) and the potential for vandalism that would result in a release of undetermined wastes and/or PCB contaminated oil into the environment, it is necessary that a timely response action be undertaken to stabilize, identify and dispose of these materials properly.

Messrs. Joe Rotola and Kevin Matheis of your Removal Action Branch have visited the site with DEC staff and are aware of the situation concerning the PRP and the abandoned wastes which require disposal.

A report entitled "Summary of Sampling Results Lackawanna Foundry, 184 Madison Avenue, Buffalo, New York" is available and will be provided, along with other information, as necessary.

If you have any questions, please contact Mr. Peter Buechi at (716) 851-7220 or Tom Vickerson at (518) 457-7878.

Sincerely,



Michael J. O'Toole, Jr.

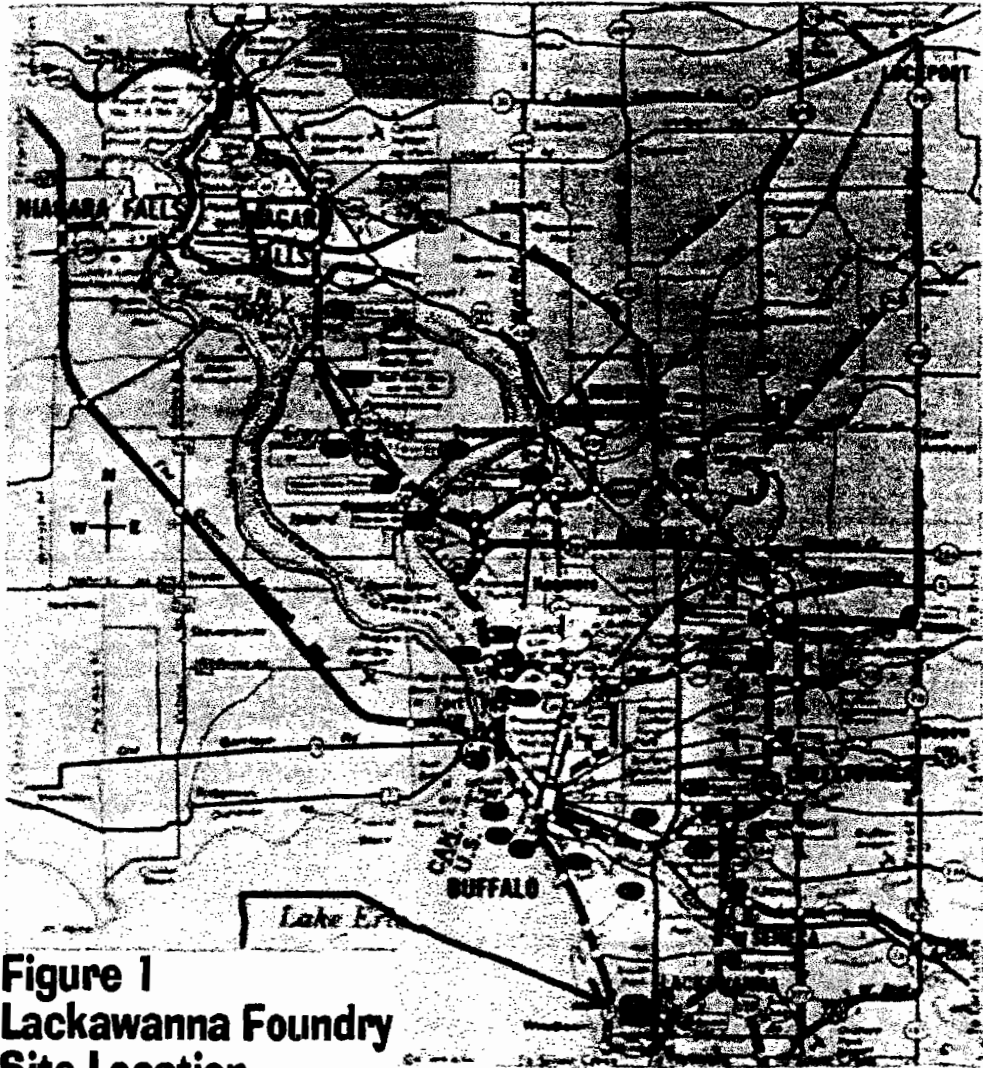
Director

Division of Environmental Remediation

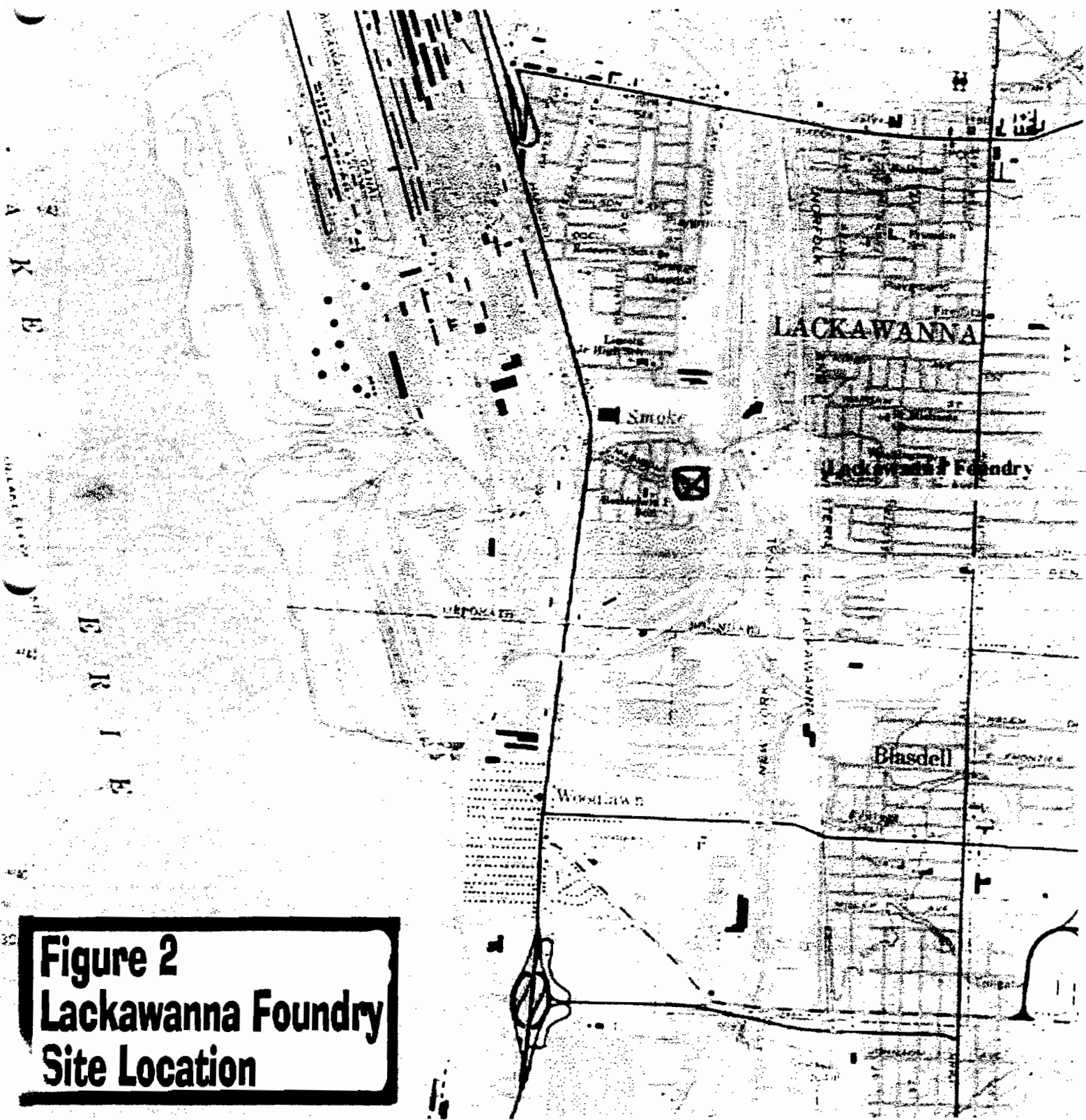
cc: Mr. R. Salkie - USEPA, Region II, Edison, NJ  
Mr. J. Rotola - USEPA, Region II, Edison, NJ  
Mr. B. Sprague - USEPA, Region II, Edison, NJ  
Mr. G. Zachos - USEPA, Region II, Edison, NJ



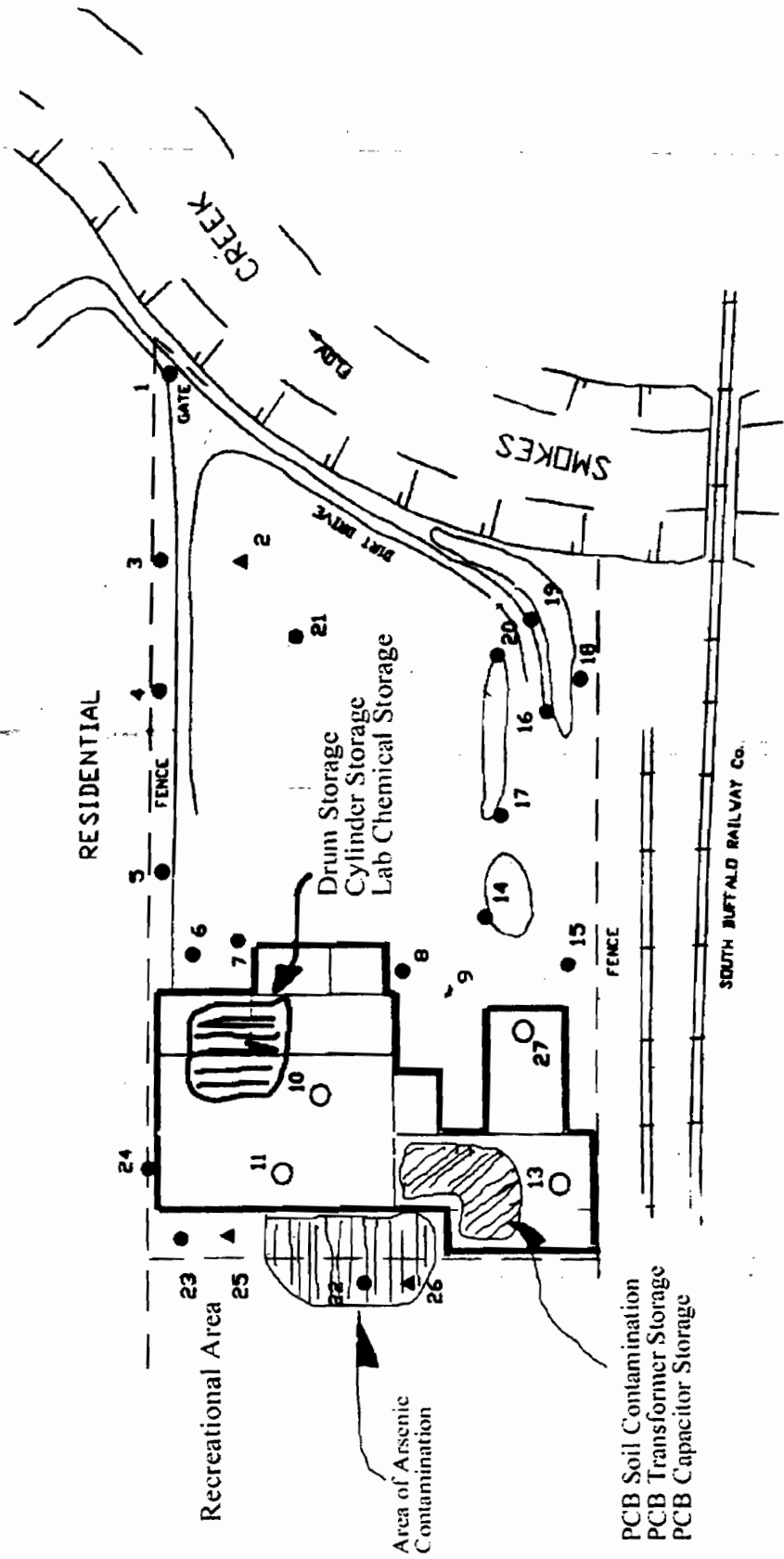
Attachment 2



**Figure 1**  
**Lackawanna Foundry**  
**Site Location**



**Figure 2**  
**Lackawanna Foundry**  
**Site Location**



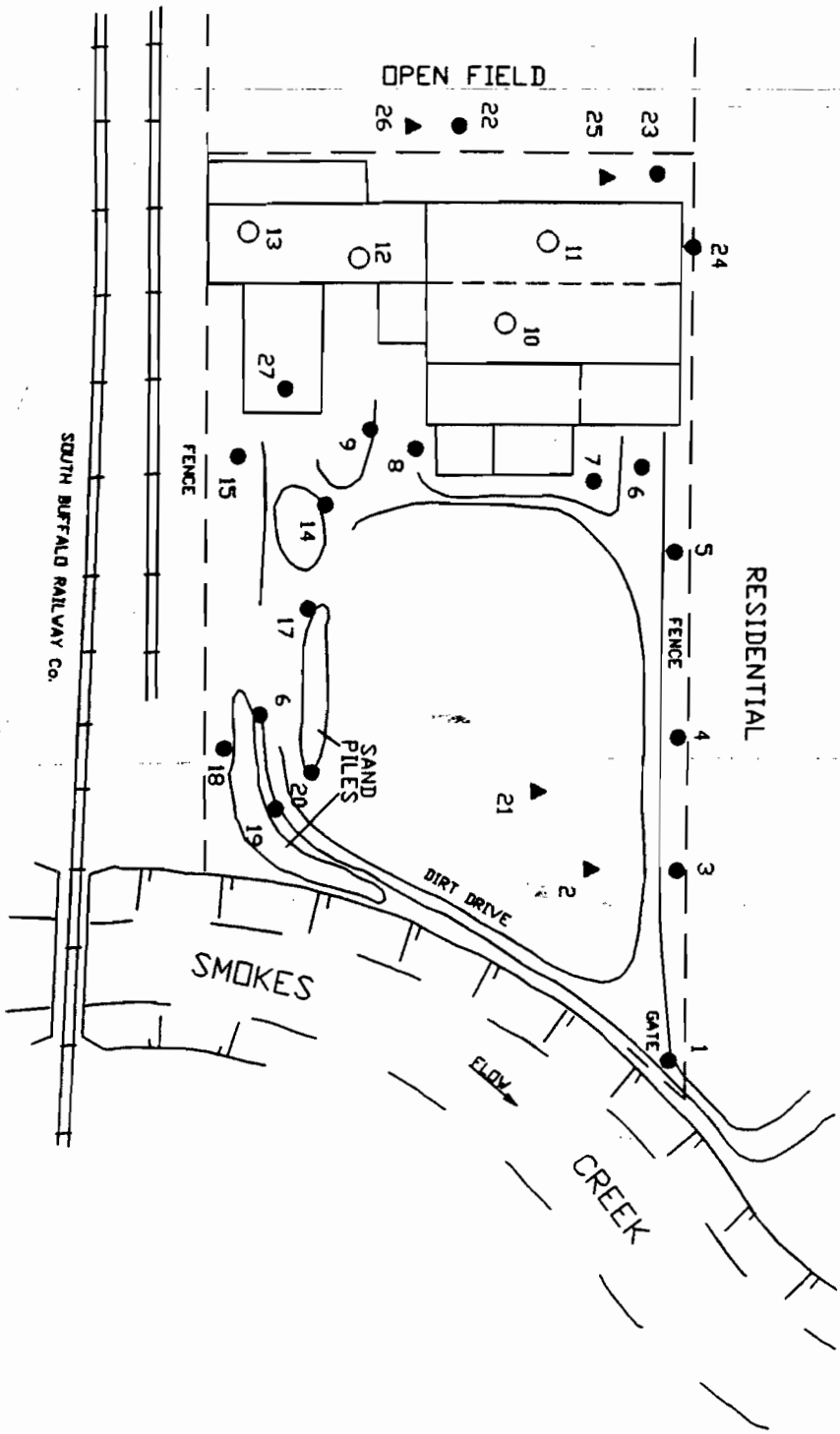
**FIGURE 1**  
**SAMPLING LOCATION MAP**  
**LACKAWANNA FOUNDRY SITE**  
**LACKAWANNA, NEW YORK**  
**MAY 1999**

- SURFACE SAMPLE
- COMPOSITE SAMPLE
- ▲ SURFACE AND SUBSURFACE SAMPLE



U.S. EPA ENVIRONMENTAL RESPONSE TEAM CENTER  
 RESPONSE ENGINEERING AND ANALYTICAL CONTRACT  
 U.S. #123-123456-789-01

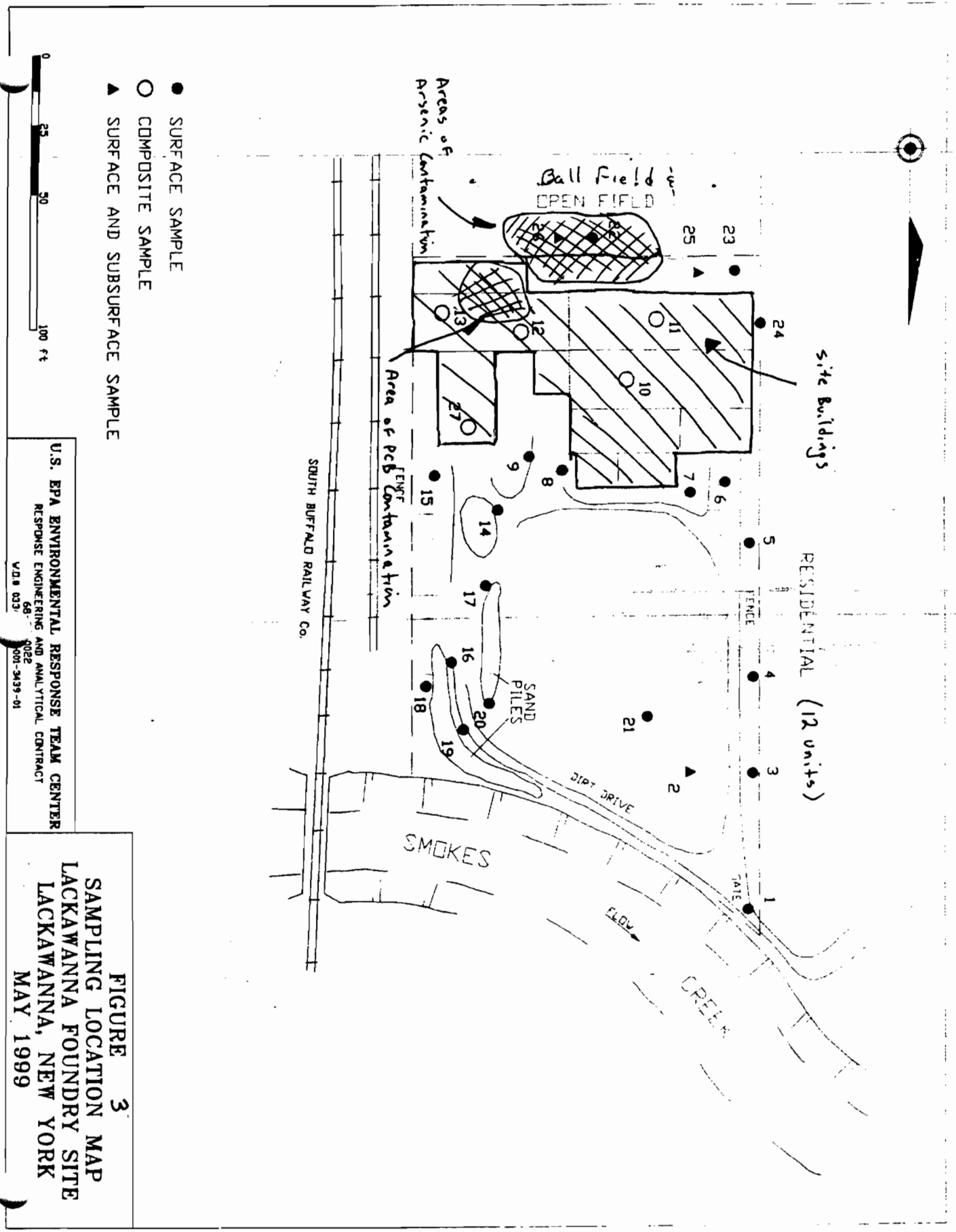
Attachment 3



- SURFACE SAMPLE
- COMPOSITE SAMPLE
- ▲ SURFACE AND SUBSURFACE SAMPLE



LACKAWANNA FOUNDRY SITE  
LACKAWANNA, NEW YORK



U.S. EPA ENVIRONMENTAL RESPONSE TEAM CENTER  
 RESPONSE ENGINEERING AND ANALYTICAL CONTRACT  
 V.O.B. 0337 0022 001-3439-01

**FIGURE 3**  
 SAMPLING LOCATION MAP  
 LACKAWANNA FOUNDRY SITE  
 LACKAWANNA, NEW YORK  
 MAY 1999

Table 1  
Summarized Pesticide/PCB Sampling Results  
Lackawanna Foundry Site  
Lackawanna, New York  
May 1999

| Sample Number<br>Location | TAGM<br>Conc.<br>µg/kg | A16650<br>1    |              | A16651<br>2    |              | A16652<br>2D   |              | A16653<br>3    |              | A16654<br>4    |              | A16655<br>5    |              | A16656<br>6    |              | A16657<br>7    |              |
|---------------------------|------------------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|
|                           |                        | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg |
| Heptachlor                | 100                    | ND             | 5.8          | ND             | 3.5          | ND             | 3.5          | ND             | 4.6          | ND             | 4.1          | ND             | 4.1          | ND             | 3.9          | ND             | 4.4          |
| p,p'-DDE                  | 2100                   | 12             | 5.8          | ND             | 3.5          | ND             | 3.5          | 30             | 4.6          | 19             | 4.1          | 0.9 J          | 4.1          | 12             | 3.9          | 4.5            | 4.4          |
| Dieldrin                  | 44                     | ND             | 5.8          | ND             | 3.5          | ND             | 3.5          | ND             | 4.6          | ND             | 4.1          | ND             | 4.1          | ND             | 3.9          | 2.9 J          | 4.4          |
| p,p'-DDD                  | 2900                   | ND             | 5.8          | ND             | 3.5          | ND             | 3.7          | ND             | 4.6          | 6.9            | 4.1          | ND             | 4.1          | ND             | 3.9          | ND             | 4.4          |
| Endosulfan (II)           | 900                    | ND             | 5.8          | ND             | 3.5          | ND             | 3.5          | ND             | 4.6          | ND             | 4.1          | ND             | 4.1          | ND             | 3.9          | ND             | 4.4          |
| p,p'-DDT                  | 2100                   | 25             | 5.8          | ND             | 3.5          | ND             | 3.5          | 14             | 4.6          | 6.4            | 4.1          | ND             | 4.1          | ND             | 3.9          | 9.6            | 4.4          |
| Aroclor 1242              | Surf.<br>1000          | ND             | 73           | ND             | 3.5          | ND             | 46           | ND             | 57           | ND             | 51           | ND             | 52           | ND             | 3.9          | ND             | 55           |
| Aroclor 1248              |                        | ND             | 73           | 560            | 44           | 2700           | 46           | ND             | 57           | 120            | 51           | 15 J           | 52           | 770            | 49           | 590            | 55           |
| Aroclor 1254              | Sub.<br>10000          | ND             | 73           | ND             | 44           | ND             | 46           | 25 J           | 57           | 81             | 51           | 161W           | 52           | 1200           | 49           | 200            | 55           |
| Aroclor 1260              |                        | ND             | 73           | ND             | 44           | ND             | 46           | 19 J           | 57           | 31 J           | 51           | 13 J           | 52           | 690            | 49           | 67             | 55           |

MDL - Method Detection Limit

Conc. - Concentration

µg/kg - micrograms per kilogram

ND - Not Detected

J - Detected Below the MDL

TAGM - New York State Department of Environmental Conservation-Technical & Administrative Guidance Memorandum

Surf. - Surface

Sub. - Subsurface

NA - Not Applicable

W - Weathering

Shading indicates above NYSDEC-TAGM

p. 27



Table I(Continued)  
 Summarized Pesticide/PCB Sampling Results  
 Lackawanna Foundry Site  
 Lackawanna, New York  
 May 1999

| Sample Number<br>Location | TAGM<br>Conc.<br>µg/kg | A16658<br>8    |              | A16659<br>9    |              | A16660<br>10   |              | A16661<br>11   |              | A16662<br>12   |              | A16663<br>13   |              | A16664<br>14   |              | A16665<br>15   |              |
|---------------------------|------------------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|
|                           |                        | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg |
| Heptachlor                | 100                    | ND             | 3.8          | ND             | 3.7          | ND             | 3.4          | ND             | 3.9          | ND             | 4.2          | ND             | 3.7          | ND             | 4.5          | ND             | 5.3          |
| p,p'-DDE                  | 2100                   | 2.5 J          | 3.8          | 17             | 3.7          | 4.7            | 3.4          | 3.9 J          | 3.9          | 28             | 4.2          | 13             | 3.7          | 93             | 4.5          | ND             | 5.3          |
| Dieldrin                  | 44                     | 2.0 J          | 3.8          | ND             | 3.7          | 2.7 J          | 3.4          | 1.1 J          | 3.9          | 6.3            | 4.2          | 2.6 J          | 3.7          | 2.4 J          | 4.5          | ND             | 5.3          |
| p,p'-DDD                  | 2900                   | ND             | 3.8          | ND             | 3.7          | ND             | 3.4          | ND             | 3.9          | ND             | 4.2          | ND             | 3.7          | ND             | 4.5          | ND             | 5.3          |
| Endosulfan (II)           | 900                    | ND             | 3.8          | ND             | 3.7          | ND             | 3.4          | ND             | 3.9          | ND             | 4.2          | ND             | 3.7          | ND             | 4.5          | ND             | 5.3          |
| p,p'-DDT                  | 2100                   | 5.3            | 3.8          | ND             | 3.7          | ND             | 3.4          | 5.2            | 3.9          | ND             | 4.2          | ND             | 3.7          | ND             | 4.5          | ND             | 5.3          |
| Aroclor 1242              | Surf.<br>1000          | ND             | 48           | ND             | 3.7          | 6500           | 42           | ND             | 49           | ND             | 52           | ND             | 46           | ND             | 4.5          | ND             | 5.3          |
| Aroclor 1248              |                        | 68             | 48           | 3500           | 46           | ND             | 42           | 1000           | 49           | 7400           | 52           | 3100           | 46           | 2800           | 57           | 42 J           | 67           |
| Aroclor 1254              | Sub.<br>10000          | 68             | 48           | 1300           | 46           | ND             | 42           | ND             | 49           | 81             | 52           | 1300           | 46           | ND             | 57           | ND             | 67           |
| Aroclor 1260              |                        | 32 J           | 48           | 320            | 46           | ND             | 42           | ND             | 49           | 31 J           | 52           | 240            | 46           | ND             | 57           | ND             | 67           |

MDL - Method Detection Limit

Conc. - Concentration

µg/kg - micrograms per kilogram

ND - Not Detected

J - Detected Below the MDL

TAGM - New York State Department of Environmental Conservation Technical & Administrative Guidance Memorandum

Surf - Surface

Sub - Subsurface

NA - Not Applicable

Shading indicates above TAGM

0.28

Table 1 (Continued)  
 Summarized Pesticide/PCB Sampling Results  
 Lackawanna Foundry Site  
 Lackawanna, New York  
 May 1999

| Sample Number<br>Location | TAGM<br>Conc.<br>µg/kg | A16666<br>16   |              | A16667<br>17   |              | A16668<br>18   |              | A16669<br>19   |              | A16670<br>20   |              | A16671<br>21   |              | A16672<br>22   |              | A16673<br>23   |              |
|---------------------------|------------------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|
|                           |                        | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg |
| g-BHC                     | 300                    | ND             | 3.7          | ND             | 3.8          | ND             | 3.6          | ND             | 3.6          | ND             | 3.6          | 2.5 J          | 3.7          | ND             | 6.9          | ND             | 4.1          |
| Heptachlor                | 100                    | ND             | 3.7          | ND             | 3.8          | ND             | 3.6          | ND             | 3.6          | ND             | 3.6          | ND             | 3.7          | ND             | 6.9          | ND             | 4.1          |
| p,p'-DDE                  | 2100                   | 0.7 J          | 3.7          | 2.2 J          | 3.8          | ND             | 3.6          | 5.1            | 3.6          | ND             | 3.6          | ND             | 3.7          | 8.4            | 6.9          | 2.3 J          | 4.1          |
| Dieldrin                  | 44                     | ND             | 3.7          | 0.6 J          | 3.8          | ND             | 3.6          | 2.2 J          | 3.6          | ND             | 3.6          | ND             | 3.7          | 21             | 6.9          | ND             | 4.1          |
| p,p'-DDD                  | 2900                   | ND             | 3.7          | ND             | 3.8          | ND             | 3.6          | ND             | 3.6          | ND             | 3.6          | ND             | 3.7          | ND             | 6.9          | 1.6 J          | 4.1          |
| Endosulfan (II)           | 900                    | ND             | 3.7          | ND             | 3.8          | ND             | 3.6          | ND             | 3.6          | ND             | 3.6          | ND             | 3.7          | ND             | 6.9          | ND             | 4.1          |
| p,p'-DDT                  | 2100                   | ND             | 3.7          | ND             | 3.8          | ND             | 3.6          | 4.9            | 3.6          | ND             | 3.6          | ND             | 3.7          | ND             | 6.9          | 1.9 J          | 4.1          |
| Aroclor 1242              | Surf.<br>1000          | 340            | 47           | ND             | 3.8          | ND             | 45           | 2500           | 45           | ND             | 46           | ND             | 46           | ND             | 86           | ND             | 51           |
| Aroclor 1248              |                        | ND             | 47           | 620            | 47           | 77             | 45           | ND             | 45           | ND             | 46           | 320            | 46           | 11JW           | 86           | 110            | 51           |
| Aroclor 1254              | Sub.<br>10000          | ND             | 47           | ND             | 47           | 33 J           | 45           | ND             | 45           | ND             | 46           | 300            | 46           | 16JW           | 86           | ND             | 51           |
| Aroclor 1260              |                        | ND             | 47           | ND             | 47           | ND             | 45           | ND             | 45           | 3900           | 46           | 200            | 46           | ND             | 86           | 14 J           | 51           |

MDL - Method Detection Limit

Conc. - Concentration

µg/kg - micrograms per kilogram

ND - Not Detected

J - Detected Below the MDL

TAGM - New York State Department of Environmental Conservation-Technical & Administrative Guidance Memorandum

Surf - Surface

Sub - Subsurface

NA - Not Applicable

W - Weathering

Shading indicates above TAGM

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Table 1(Continued)  
 Summarized Pesticide/PCB Sampling Results  
 Lackawanna Foundry Site  
 Lackawanna, New York  
 May 1999

| Sample Number<br>Location | TAGM<br>Conc.<br>µg/kg | A16674<br>24   |              | A16675<br>25   |              | A16676<br>25D  |              | A16677<br>26   |              | A16678<br>26-D |              | A16679<br>27   |              |
|---------------------------|------------------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|
|                           |                        | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg |
| g-BHC                     | 300                    | ND             | 4.4          | ND             | 4.0          | ND             | 4.3          | ND             | 4.8          | ND             | 4.4          | ND             | 4.0          |
| Heptachlor                | 100                    | ND             | 4.4          | ND             | 4.0          | ND             | 4.3          | ND             | 4.8          | ND             | 4.4          | ND             | 4.0          |
| p,p'-DDE                  | 2100                   | 1.9 J          | 4.4          | 3.3 J          | 4.0          | 4.3            | 4.3          | 11             | 4.8          | ND             | 4.4          | ND             | 4.0          |
| Dieldrin                  | 44                     | ND             | 4.4          | ND             | 4.0          | 4.3            | 4.3          | ND             | 4.8          | ND             | 4.4          | ND             | 4.0          |
| p,p'-DDD                  | 2900                   | ND             | 4.4          | ND             | 4.0          | 4.3            | 4.3          | ND             | 4.8          | ND             | 4.4          | ND             | 4.0          |
| Endosulfan (II)           | 900                    | ND             | 4.4          | ND             | 4.0          | 4.3            | 4.3          | ND             | 4.8          | ND             | 4.4          | ND             | 4.0          |
| p,p'-DDT                  | 2100                   | ND             | 4.4          | ND             | 4.0          | 4.3            | 4.3          | 18             | 4.8          | ND             | 4.4          | ND             | 4.0          |
| Aroclor 1242              | Surf.<br>1000          | 340            | 55           | ND             | 4.0          | 54             | 54           | ND             | 60           | ND             | 54           | ND             | 50           |
| Aroclor 1248              |                        | 120            | 55           | 270            | 50           | 54             | 54           | 170            | 60           | ND             | 54           | 39000          | 50           |
| Aroclor 1254              | Sub.<br>10000          | ND             | 55           | 83             | 50           | 54             | 54           | ND             | 60           | ND             | 54           | ND             | 50           |
| Aroclor 1260              |                        | 10 J           | 55           | ND             | 50           | 54             | 54           | 42 J           | 60           | 3900           | 54           | ND             | 50           |

MDL - Method Detection Limit

Conc. - Concentration

µg/kg - micrograms per kilogram

ND - Not Detected

J - Detected Below the MDL

TAGM - New York State Department of Environmental Conservation Technical & Administrative Guidance Memorandum

Surf - Surface

Sub - Subsurface

NA - Not Applicable

Shading indicates above TAGM

Table 2  
 Summarized BNAs Sampling Results  
 Lackawanna Foundry Site  
 Lackawanna, New York  
 May 1999

| Sample Number<br>Location  | TAGM           | A166501        | A16651<br>2    | A16652<br>2-D  | A16653<br>3    | A16654<br>4    | A16655<br>5    | A16656<br>6    | A16657         |
|----------------------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|
| Analyte                    | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg |
| Phenol                     | 30             | ND             | <b>16000</b>   | <b>23000</b>   | ND             | ND             | ND             | ND             | 1900 J         |
| 2-Methylphenol             | 100            | ND             | <b>370</b>     | 410 J          | ND             | ND             | ND             | ND             | ND             |
| Naphthalene                | 13000          | ND             | 970            | 960 J          | ND             | ND             | 310 J          | ND             | ND             |
| 2-Methylnaphthalene        | 36400          | ND             | 360            | 640 J          | ND             | ND             | 270 J          | ND             | ND             |
| Acenaphthylene             | 41000          | ND             | ND             | ND             | ND             | ND             | 200 J          | ND             | ND             |
| Acenaphthene               | 50000          | ND             | 95 J           | ND             | ND             | ND             | 740            | ND             | ND             |
| Dibenzofuran               | 6200           | ND             | 100 J          | ND             | ND             | ND             | 490            | ND             | ND             |
| Fluorene                   | 50000          | ND             | 87 J           | ND             | ND             | ND             | 880            | ND             | ND             |
| Phenanthrene               | 50000          | ND             | 1400           | ND             | ND             | ND             | 7900           | ND             | ND             |
| Anthracene                 | 50000          | ND             | 220 J          | ND             | ND             | ND             | 1800           | ND             | ND             |
| Carbazole                  | NA             | ND             | 190 J          | ND             | ND             | ND             | 1200           | ND             | ND             |
| Di-n-butylphthalate        | 8100           | ND             | 150 J          | ND             | ND             | ND             | ND             | ND             | ND             |
| Fluoranthene               | 50000          | ND             | 2000           | 380 J          | 1000 J         | 2300 J         | 9900           | 2000 J         | 2400 J         |
| Pyrene                     | 50000          | ND             | 1500           | ND             | ND             | 1900 J         | 7800           | 1700 J         | 2200 J         |
| Benzo(a)anthracene         | 224            | ND             | 800            | ND             | ND             | 1100 J         | <b>5000</b>    | 1000 J         | 1500 J         |
| Chrysene                   | 400            | ND             | 920            | ND             | ND             | 1400 J         | <b>5300</b>    | 1100 J         | 1900 J         |
| Bis(2-Ethylhexyl)phthalate | 50000          | ND             | 640            | 1300 J         | 1000 J         | 1000 J         | 260 J          | 810 J          | ND             |
| Benzo(b)fluoranthene       | 1100           | ND             | <b>1100</b>    | ND             | ND             | 1700 J         | <b>5800</b>    | 1400 J         | 3100 J         |
| Benzo(k)fluoranthene       | 1100           | ND             | 880            | ND             | ND             | 1400 J         | <b>4500</b>    | 1200 J         | 2400 J         |
| Benzo(a)pyrene             | 61             | ND             | 1.100          | ND             | ND             | 1700 J         | <b>5800</b>    | 1300 J         | 3300 J         |
| Indeno(1,2,3,-cd)pyrene    | 3200           | ND             | 750            | ND             | ND             | 1.300 J        | <b>3400</b>    | 990 J          | 3200 J         |
| Dibenzo(a,h)anthracene     | 14             | ND             | 300 J          | ND             | ND             | ND             | <b>1300</b>    | ND             | 1200 J         |
| Benzo(g,h,i)perylene       | 50000          | ND             | 890            | ND             | ND             | 1600 J         | 3700           | 1100 J         | 3600 J         |

MDL - Method Detection Limit

Conc. - Concentration

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J - Detected Below the MDL

TAGM - New York State Department of Environmental Conservation Technical & Administrative Guidance Memorandum

Surf - Surface

Sub - Subsurface

NA - Not Applicable

Shading indicates above TAGM

## Summarized BNAs Sampling Results

Lackawanna Foundry Site

Lackawanna, New York

May 1999

| Sample Number<br>Location  | TAGM           | A16658<br>8    | A16659<br>9    | A16660<br>10   | A16661<br>11   | A16662<br>12   | A16663<br>13   | A16664<br>14   | A16665<br>15   |
|----------------------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|
| Analyte                    | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg |
| Phenol                     | 30             | 1500           | 4900           | 1800           | 9700           | 4700           | 2200           | 7000           | 1600           |
| 2-Methylphenol             | 100            | ND             | ND             | 570 J          | ND             | ND             | ND             | ND             | ND             |
| Naphthalene                | 13000          | ND             | ND             | 3100           | 2100 J         | 960 J          | 670 J          | 1400 J         | ND             |
| 2-Methylnaphthalene        | 36400          | ND             | ND             | 1500 J         | ND             | 500 J          | 400 J          | ND             | ND             |
| Acenaphthylene             | 41000          | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Acenaphthene               | 50000          | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Dibenzofuran               | 6200           | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Fluorene                   | 50000          | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Phenanthrene               | 50000          | ND             | ND             | 680 J          | ND             | 530 J          | ND             | ND             | ND             |
| Anthracene                 | 50000          | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Carbazole                  | NA             | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Di-n-butylphthalate        | 8100           | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Fluoranthene               | 50000          | 1300 J         | 1000 J         | ND             | ND             | 600 J          | ND             | ND             | ND             |
| Pyrene                     | 50000          | 1200 J         | 880 J          | ND             | ND             | 900 J          | ND             | ND             | ND             |
| Butylbenzylphthalate       | 50000          | ND             | 6000           | ND             | ND             | ND             | ND             | ND             | ND             |
| Benzo(a)anthracene         | 224            | ND             | NA             | ND             | ND             | ND             | ND             | 460 J          | ND             |
| Chrysene                   | 400            | 1000 J         | NA             | ND             | ND             | ND             | ND             | ND             | ND             |
| Bis(2-Ethylhexyl)phthalate | 50000          | 8700           | 1200 J         | 1300 J         | 16000          | 1300 J         | 540 J          | ND             | ND             |
| Benzo(b)fluoranthene       | 1100           | 1600 J         | 940 J          | ND             | ND             | ND             | ND             | ND             | ND             |
| Benzo(k)fluoranthene       | 1100           | 1400 J         | 830 J          | ND             | ND             | ND             | ND             | ND             | ND             |
| Benzo(a)pyrene             | 61             | 1400 J         | 920 J          | ND             | ND             | ND             | ND             | ND             | ND             |
| Indeno(1,2,3,-cd)pyrene    | 3200           | 1700 J         | 770 J          | ND             | ND             | ND             | ND             | ND             | ND             |
| Dibenzo(a,h)anthracene     | 14             | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Benzo(g,h,i)perylene       | 50000          | 1900 J         | 920 J          | ND             | ND             | ND             | ND             | ND             | ND             |

MDL - Method Detection Limit

Conc. - Concentration

µg/kg - micrograms per kilogram

ND - Not Detected

J - Detected Below the MDL

TAGM - New York State Department of Environmental Conservation-Technical &amp; Administrative Guidance Memorandum

NA - Not Applicable

Shading indicates above TAGM

Table 2 (Continued)  
Summarized BNAs Sampling Results  
Lackawanna Foundry Site  
Lackawanna, New York  
May 1999

| Sample Number<br>Location  | TAGM           | A16666<br>16   | A16667<br>17   | A16668<br>18   | A16669<br>19   | A16670<br>20   | A16671<br>21   | A16672<br>22   | A16673<br>23   |
|----------------------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|
| Analyte                    | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg |
| Phenol                     | 30             | 840 J          | 910 J          | <b>11000</b>   | <b>4000</b>    | <b>18000</b>   | 5600 J         | ND             | 2400 J         |
| 2-Methylphenol             | 100            | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Naphthalene                | 13000          | ND             | ND             | 850 J          | 580            | <b>4400</b>    | ND             | ND             | 1100 J         |
| 2-Methylnaphthalene        | 36400          | ND             | ND             | ND             | 100 J          | 1100 J         | ND             | ND             | 850 J          |
| Acenaphthylene             | 41000          | ND             | ND             | ND             | ND             | ND             | ND             | 420 J          | 1300 J         |
| Acenaphthene               | 50000          | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Dibenzofuran               | 62000          | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Fluorene                   | 50000          | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Phenanthrene               | 50000          | ND             | ND             | ND             | ND             | ND             | ND             | ND             | 2200 J         |
| Anthracene                 | 50000          | ND             | ND             | ND             | ND             | ND             | ND             | ND             | 1400 J         |
| Carbazole                  | NA             | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Di-n-butylphthalate        | 8100           | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Fluoranthene               | 50000          | ND             | 860 J          | ND             | 74 J           | 1600 J         | 3400 J         | 4700           | 8500           |
| Pyrene                     | 50000          | ND             | 740 J          | ND             | ND             | 1400 J         | 3400 J         | 4100           | 8800           |
| Butylbenzylphthalate       | 224            | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Benzo(a)anthracene         | 400            | ND             | 410 J          | ND             | ND             | 840 J          | 2100 J         | <b>2600</b>    | <b>5200</b>    |
| Chrysene                   | 400            | ND             | 510 J          | ND             | ND             | 900 J          | 2600 J         | <b>2400</b>    | <b>5400</b>    |
| Bis(2-Ethylhexyl)phthalate | 50000          | ND             | ND             | ND             | 180 J          | 760 J          | 1700 J         | 430 J          | ND             |
| Benzo(b)fluoranthene       | 1100           | ND             | 580 J          | ND             | ND             | 910 J          | 3500 J         | <b>2700</b>    | <b>5700</b>    |
| Benzo(k)fluoranthene       | 1100           | ND             | 520 J          | ND             | ND             | 890 J          | 3400 J         | <b>2500</b>    | <b>5700</b>    |
| Benzo(a)pyrene             | 61             | ND             | 570 J          | ND             | ND             | 980 J          | 3700 J         | <b>3100</b>    | <b>5300</b>    |
| Indeno(1,2,3,-cd)pyrene    | 3200           | ND             | 420 J          | ND             | ND             | 660 J          | 3100 J         | 1600 J         | 3100 J         |
| Dibenzo(a,h)anthracene     | 14             | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Benzo(g,h,i)perylene       | 50000          | ND             | 470 J          | ND             | ND             | 770 J          | 3600 J         | 1700 J         | 3200 J         |

MDL - Method Detection Limit

Conc. - Concentration

µg/kg - micrograms per kilogram

ND - Not Detected

J - Detected Below the MDL

TAGM - New York State Department of Environmental Conservation-Technical & Administrative Guidance Memorandum

NA - Not Applicable, Shading indicates above TAGM

Table 3  
 Summarized TAL Metals Sampling Results  
 Lackawanna Foundry Site  
 Lackawanna, New York  
 May 1999

| Sample Number<br>Location | TAGM           | C16650<br>1    | C16651<br>2    | C16652<br>2-D  | C16653<br>3    | C16654<br>4    | C16655<br>5    | C16656<br>6    | C16657<br>7    |
|---------------------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|
| Analyte                   | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg |
| Aluminum                  | *              | 11000          | 1300           | 1200           | 7800           | 13000          | 12000          | 4600           | 5300           |
| Antimony                  | *              | ND             | ND             | ND             | ND             | ND             | ND             | 6.4            | ND             |
| Arsenic                   | 7.5*           | 14             | 1.6            | 2.5            | 11             | 13             | 15             | 20             | 6.3            |
| Barium                    | 300*           | 160            | 18             | 17             | 140            | 150            | 130            | 270            | 120            |
| Beryllium                 | 0.16*          | 0.93           | ND             | ND             | 0.81           | 2.2            | 0.85           | ND             | ND             |
| Cadmium                   | 1*             | 3.2            | 0.48           | 0.66           | 3.7            | 2.4            | 2.4            | 20             | 3.6            |
| Calcium                   | *              | 21000          | 1900           | 3300           | 94000          | 120000         | 14000          | 21000          | 18000          |
| Chromium                  | 10*            | 45             | 25             | 15             | 33             | 28             | 29             | 190            | 55             |
| Cobalt                    | 30*            | 12             | 2.1            | 1.2            | 7.7            | 6.2            | 14             | 20             | 6.7            |
| Copper                    | 25*            | 74             | 57             | 35             | 75             | 67             | 200            | 440            | 110            |
| Iron                      | 2000*          | 43000          | 24000          | 13000          | 46000          | 30000          | 35000          | 200000         | 47000          |
| Lead                      | *              | 300            | 44             | 38             | 200            | 260            | 180            | 610            | 2100           |
| Magnesium                 | *              | 4800           | 600            | 560            | 7200           | 11000          | 4800           | 3100           | 3300           |
| Manganese                 | *              | 1200           | 250            | 140            | 800            | 1300           | 700            | 1800           | 1100           |
| Mercury                   | 0.1            | 0.21           | 0.06           | 0.1            | 0.11           | 0.16           | 0.11           | 1.1            | 0.22           |
| Nickel                    | 13*            | 42             | 15             | 9.2            | 28             | 32             | 41             | 160            | 110            |
| Potassium                 | *              | 1500           | ND             | ND             | 920            | 1500           | 1200           | 630            | 570            |
| Selenium                  | 2*             | 1.0            | ND             | ND             | 0.79           | 0.98           | 0.58           | 0.77           | ND             |
| Silver                    | *              | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Sodium                    | *              | 110            | ND             | 64             | 140            | 520            | 79             | 410            | 230            |
| Thallium                  | *              | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Vanadium                  | 150*           | 27             | 5.2            | 3.7            | 17             | 18             | 23             | 20             | 20             |
| Zinc                      | 20*            | 760            | 140            | 79             | 640            | 540            | 340            | 1100           | 1000           |

- See TAGM Memorandum Appendix A

DL - Method Detection Limit

nc. - Concentration

µg - micrograms per kilogram

ND - Not Detected

DL - Detected Below the MDL

-New York State Department of Environmental Conservation-Technical & Administrative Guidance Memorandum

DL indicates above TAGM

Table 3 (Continued)  
 Summarized TAL Metals Sampling Results  
 Lackawanna Foundry Site  
 Lackawanna, New York  
 May 1999

| Sample Number<br>Location | TAGM           | C16666<br>16   | C16667<br>17   | C16668<br>18   | C16669<br>19   | C16670<br>20   | C16671<br>21   | C16672<br>22   | C16673<br>23   |
|---------------------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|
| Analyte                   | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg |
| Aluminum                  | *              | 1300           | 5900           | 790            | 1400           | 2400           | 12000          | 15000          | 9300           |
| Antimony                  | *              | ND             | 6.5            | ND             | ND             | ND             | ND             | ND             | ND             |
| Arsenic                   | 7.5*           | 1.5            | 7.1            | 4.4            | 6.7            | 6.5            | <b>20</b>      | <b>77</b>      | <b>8.4</b>     |
| Barium                    | 300*           | 15             | 180            | 9.0            | 29             | 130            | 200            | 170            | 94             |
| Beryllium                 | 0.16*          | ND             | ND             | ND             | ND             | ND             | <b>0.96</b>    | <b>1.2</b>     | <b>0.66</b>    |
| Cadmium                   | 1*             | ND             | ND             | ND             | 0.59           | <b>2.3</b>     | <b>13</b>      | <b>2.3</b>     | <b>1.5</b>     |
| Calcium                   | *              | 1200           | 130000         | 1200           | 850            | 3800           | 23000          | 34000          | 12000          |
| Chromium                  | 10*            | 6.1            | <b>33</b>      | 8.0            | <b>97</b>      | <b>130</b>     | <b>95</b>      | <b>32</b>      | <b>24</b>      |
| Cobalt                    | 30*            | 1.4            | 47             | ND             | 4.4            | 12             | 15             | 11             | 8.0            |
| Copper                    | 25*            | <b>44</b>      | <b>240</b>     | <b>28</b>      | <b>300</b>     | <b>720</b>     | <b>270</b>     | <b>47</b>      | <b>52</b>      |
| Iron                      | 2000*          | <b>16000</b>   | <b>25000</b>   | <b>5800</b>    | <b>58000</b>   | <b>93000</b>   | <b>110000</b>  | <b>32000</b>   | <b>30000</b>   |
| Lead                      | *              | 58             | 300            | 21             | 140            | 280            | 660            | 180            | 160            |
| Magnesium                 | *              | 410            | 19000          | 340            | 410            | 1000           | 4700           | 5900           | 3600           |
| Manganese                 | *              | 210            | 840            | 75             | 760            | 690            | 1600           | 1200           | 560            |
| Mercury                   | 0.1            | ND             | 0.06           | 0.04           | ND             | 1.3            | <b>0.77</b>    | <b>0.73</b>    | <b>0.13</b>    |
| Nickel                    | 13*            | 9.5            | 38             | 2.7            | 57             | 150            | 100            | 36             | 28             |
| Potassium                 | *              | ND             | 470            | ND             | ND             | 280            | 900            | 1500           | 1100           |
| Selenium                  | 2*             | ND             | ND             | ND             | ND             | ND             | 1.0            | ND             | 0.57           |
| Silver                    | *              | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Sodium                    | *              | ND             | 330            | ND             | 68             | 68             | 360            | 120            | 130            |
| Thallium                  | *              | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Vanadium                  | 150*           | 3.1            | 11             | 2.6            | 14             | 21             | 36             | 27             | 19             |
| Zinc                      | 20*            | <b>98</b>      | <b>1300</b>    | <b>55</b>      | <b>310</b>     | <b>290</b>     | <b>1000</b>    | <b>1100</b>    | <b>660</b>     |

\* - See TAGM Memorandum Appendix A

MDL - Method Detection Limit

Conc. - Concentration

µg/kg - micrograms per kilogram

ND - Not Detected

\* - Detected Below the MDL

TAGM - New York State Department of Environmental Conservation Technical & Administrative Guidance Memorandum


Shading indicates above TAGM



REFERENCE NO. 4

Personal Logbook  
 Issued to: Donna M. Janda  
 Roy F. Weston, Inc.  
 West Chester, PA  
 Emp. # 11644

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regarding  
wgs +  
13007).  
estrogen  
2P5  
15  
September

40 Oct 2000 SAT member Donna Janda  
1030 arrives at Lackawanna Foundry  
Site, 2 Elm Street, Lackawanna,  
Crie County, NY to meet with  
EPA ~~Work~~ work Assignment  
Manager (WAM) Dennis Mumhall,  
and On-Scene Coordinator (OSC)  
Kevin Mathers.

1045 WAM Mumhall onsite  
1100 OSC Mathers onsite. OSC  
Mathers begins to discuss history  
of site and Removal Action which  
began in 1999. The Lackawanna  
site is the location of a former  
foundry (approx. 2 acres) that  
produced specialty cast iron  
molds for over 60 years. The  
current property owner inherited  
the site from his father. Foundry  
operations steadily decreased  
over the past 30 years. Operations  
ceased completely in 1977. In 1997,  
NYSDEC inspected the site + had  
the property owner conduct an  
inventory of all on-site containers  
and to sample PCB transformers.  
The property owner provided the  
inventory but was subsequently  
unable to conduct a cleanup  
due to financial reasons. NYSDEC  
requested EPA assistance. In  
March 1999, EPA Area Emergency  
Assessment procedures (initiated

<sup>(25)</sup>  
 fence along the perimeter of the site, inventoried + moved all containers of hazardous substances into the foundry building, and repacked the leaking drums + containers. In April 1999, EPA Environmental Response Team (ERT) + Weston Response Engineering + Analytical Contract (REAC) collected surface + subsurface soil samples for an extent of contamination study. Arsenic + PCBs were the main contaminants of concern at this site. EPA began a removal action in October 1999 to address the issues found during previous investigations. The buildings were demolished + up to 2 feet of surface soil was excavated from across the property. Excavation activities were completed in April 2000. Restoration activities, including placement of 2 feet of clean topsoil and restoration of the wetland area in the middle of the site, were completed in June 2000. The purpose of the Brownfields investigation is to determine if ~~any~~ <sup>any</sup> subsurface soil + GW are contaminated to what extent (if so). The City ofackawanna ~~is~~ <sup>is</sup> intends to make this area a park if it meets the criteria for a brownfields site.

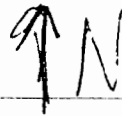
1136

1135

1130 EPA + SAT Conduct a site walk -  
 through to determine soil sampling  
 locations & monitoring well installa-  
 tion location. ~~the site~~ There are  
 no buildings on site. The site  
 is well vegetated with a small  
 pond & associated "wetlands" area  
 The central portion of the site.  
 The site is mainly flat with  
 a hilly area in the north section  
 of the property that slopes down to  
 Smokes Creek. According to  
 OSC Mathews, there is a concrete  
 buck pad buried approx 2-3 feet  
 underground along the eastern  
 portion of the site. ~~the~~ The  
 area beneath this pad was not  
 sampled previously & he would like  
 samples collected from under the  
 pad.

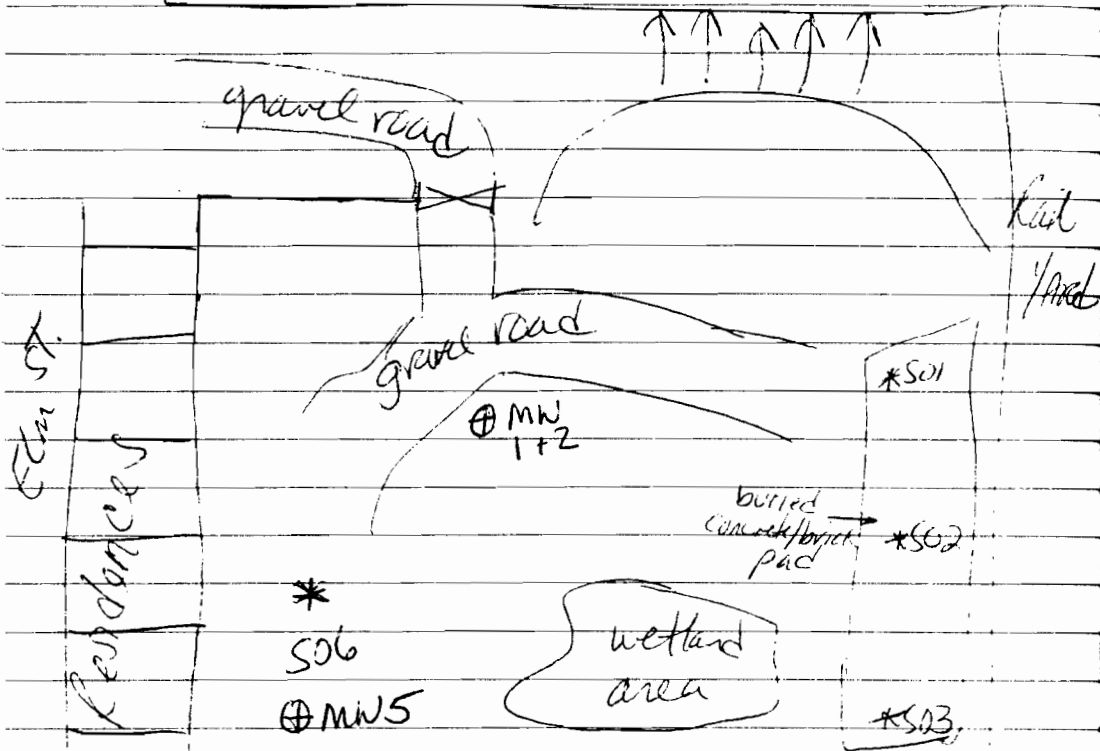
1135 Smokes Creek lies directly north  
 of the site. ~~an~~ active rail  
 yard lies to the east, a City  
 of Hackawanna recreational area  
 lies to the south & residential  
 properties lie to the west of the  
 Hackawanna site. Site Sketch is  
 on next page. No areas of  
 stressed vegetation were observed.  
 The site is enclosed by a fence  
 with a locked gate. However,  
 there is limited access for pedestrians  
 through a gap between the gate & fence.

# Site Sketch

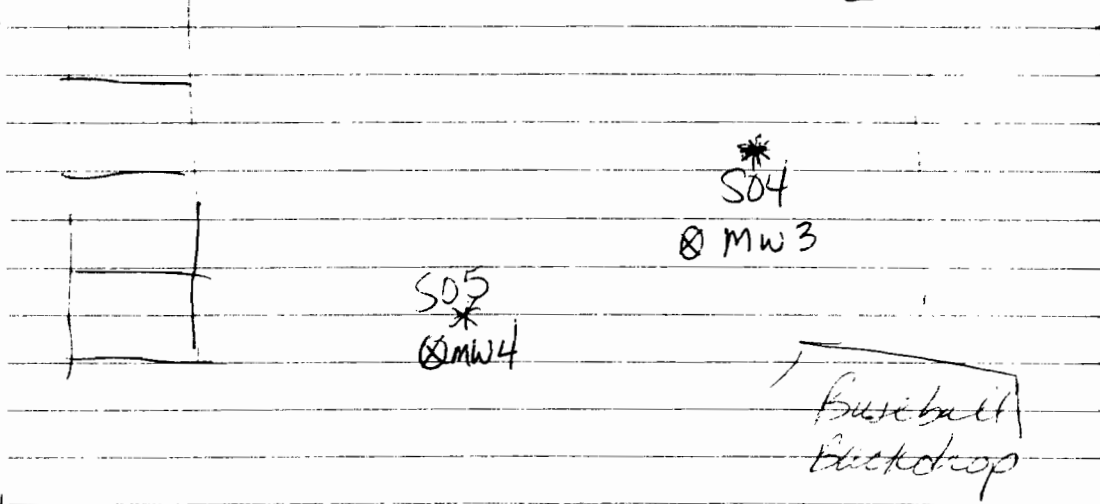


SMOKE'S CREEK

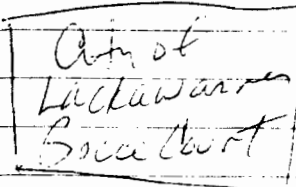
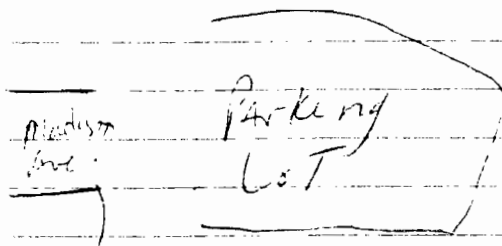
1200



1220



1240



NOT TO SCALE

↑ N

1200 Subsurface soil sample locations chosen (see map on previous page). Sample locations denoted S01-S06. One sample will be collected from locations S01-S03 for a total of 3 samples. For sample locations S04 to S06 samples will be collected at 2 foot intervals, beginning at 2-4 foot depth until either a 10 foot depth or groundwater is reached. Depth to GW at the Lackawanna Site is approx. 8-10 feet according to WC Mathies. Also, according to OSC Mathies, no one in the Lackawanna area uses groundwater for drinking. No one is on private wells.

1220 WAM Marshall would like a total of 5 monitoring wells installed at various locations throughout the site. Tentative locations are depicted on site sketch on previous page.

1240 OSC Mathies gives SAT members under copies of reports regarding the Lackawanna Foundry sites. The reports include analytical res. Hs for removal Action + REAC sampling events.

*D. J. Marshall*

1/2 in

1/4 in

\*S01

\*S02

\*S03

Baseball  
field drop

of  
warren  
court

SCALE

REFERENCE NO. 5



## New York State Department of Environmental Conservation

## Division of Environmental Remediation, Region 9

270 Michigan Avenue, Buffalo, New York, 14203-2999

Phone: (716) 851-7220 • FAX: (716) 851-7226

Website: www.dec.state.ny.us

FOIL

Releasable Non-Releasable John P. Cahill  
Commissioner

July 9, 1999

Mr. Kevin Matheis  
On-Scene Coordinator  
USEPA, Region II  
Federal Building, Rm. 1114  
111 West Huron Street  
Buffalo, New York 14202

Dear Mr. Matheis:

|                   |               |         |           |            |   |
|-------------------|---------------|---------|-----------|------------|---|
| Post-It® Fax Note | 7671          | Date    | 9-27-99   | # of pages | 3 |
| To                | KEVIN MATHEIS | From    | M. DOSTER |            |   |
| Co./Dept.         | EPA           | Co.     | DEC       |            |   |
| Phone #           |               | Phone # |           |            |   |
| Fax #             | 551-5944      | Fax #   |           |            |   |

LACKAWANNA FOUNDRY SITE  
LACKAWANNA (C), ERIE COUNTY  
REQUEST FOR ADDITIONAL SAMPLING

As we had discussed during our telephone conversation of July 7, 1999, the Department would like to ask your assistance in furthering the work at Lackawanna Foundry to close data gaps from previous investigations.

Currently no data exists to determine if local groundwater has been affected by the contaminants found at the site. It would greatly enhance the knowledge of the site if five (5) shallow monitoring wells could be installed to the top of the bedrock. Monitoring wells drilled at nearby sites, such as the Lehigh Industrial Park site, have shown auger refusals at 15 to 16 feet. Split spoon samples could be obtained from this drilling effort and yield valuable information to determine if contamination is at depth.

We had discussed the value of digging test pits in selected areas. Since capacitors containing PCB oils were hidden in the floor of the facility, it is necessary to learn if anything has been covered by the piles surrounding the property.

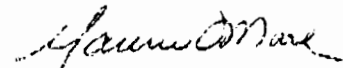
Additional sampling of the surface soils in the playground areas are needed to determine if contamination exists beyond that found in the previous investigations.

Sampling of the stream bank of Smokes Creek would help determine if the foundry sands have been placed into the stream channel and if so, how far have they been placed.

Finally, a sample of the accumulated water and sediment in the cellar area is necessary to determine potential for contamination existing at the facility to enter the groundwater from surface runoff and infiltration.

Enclosed is a figure to point out the location of the above items. It is the desire of this Department to see the site returned to beneficial use. Currently, the site presents the possibility of a "brownfields" restoration. Any assistance your agency can give at this site is welcomed and appreciated. If there is an opportunity to pursue the above items please let me know and I will lend any assistance I can. If you have any questions, please call me at 716/851-7220.

Sincerely,



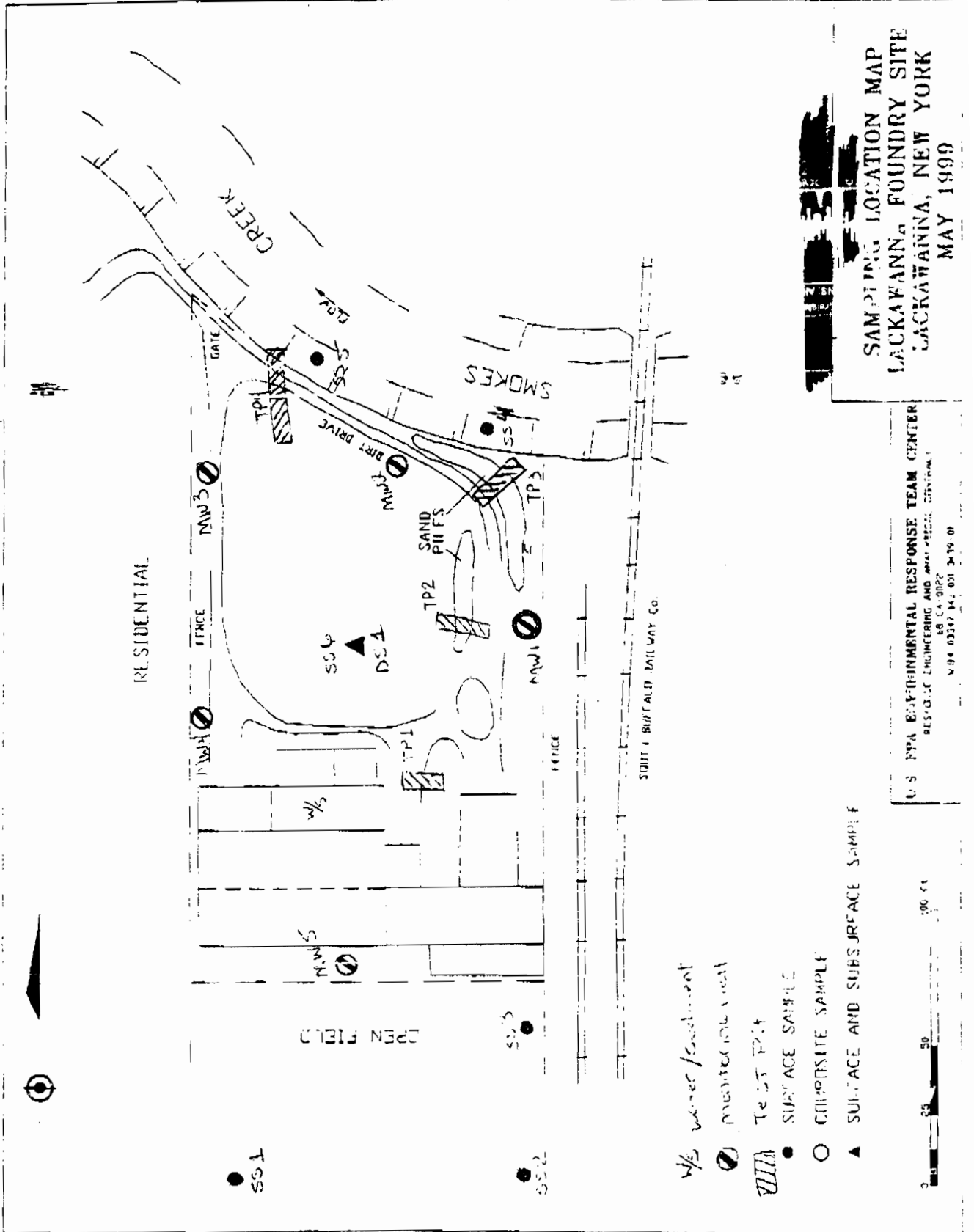
Maurice F. Moore  
Project Manager  
Hazardous Waste Remediation

MFM:lj

Enclosure

cc: Mr. Martin Doster, Regional Hazardous Waste Remediation Engineer

a:matheis.mfm



SAMPLING LOCATION MAP  
 LACKAWANNA FOUNDRY SITE  
 LACKAWANNA, NEW YORK  
 MAY 1999

U.S. EPA ENVIRONMENTAL RESPONSE TEAM CENTER  
 RESIDUAL ENGINEERING AND ANALYTICAL SERVICES  
 86 CA 2002  
 W 94 03347 347 031 3475 09

- w/s water/facility
- monitoring well
- TEST PIT
- SURFACE SAMPLE
- COMPOSITE SAMPLE
- SURFACE AND SUBSURFACE SAMPLE



REFERENCE NO. 6



FINAL REPORT  
SURFACE AND SUBSURFACE SOIL SAMPLING  
LACKAWANNA FOUNDRY SITE  
LACKAWANNA, NEW YORK  
MAY 1999

U.S. EPA Work Assignment No.: 3-439

Assessment Samples  
Fill material along  
Creek + Banks of  
Creek  
#16, 18, 19 + 20

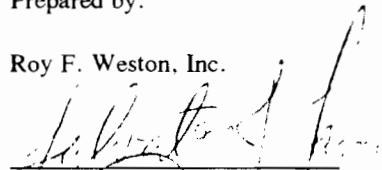


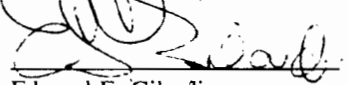
FINAL REPORT  
SURFACE AND SUBSURFACE SOIL SAMPLING  
LACKAWANNA FOUNDRY SITE  
LACKAWANNA, NEW YORK  
MAY 1999

U.S. EPA Work Assignment No.: 3-439  
Weston Work Order No.: 03347-143-001-3439-01  
U.S. EPA Contract No.: 68-C4-0022

Prepared by:

Roy F. Weston, Inc.

  
\_\_\_\_\_  
Salvatore J. Fama  
Task Leader

  
\_\_\_\_\_  
Edward F. Gilardi  
Program Manager

  
\_\_\_\_\_  
Date

  
\_\_\_\_\_  
Date

Prepared for:

U.S. EPA/ERTC

Philip Campagna  
Work Assignment Manager

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### APPENDICES

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|----------|-----------------------------------|

## 1.0 INTRODUCTION

### 1.1 Objective of this Study

The objective of this project was for the Response Engineering and Analytical Contract (REAC) to provide technical support to the United States Environmental Protection Agency/Environmental Response Team Center (U.S. EPA/ERTC) in conducting surface and subsurface soil sampling for volatile organic compounds (VOCs), pesticides, polychlorinated biphenyl (PCBs), base neutral acid extractables (BNA), target analyte list (TAL) metals, and cyanide at the Lackawanna Foundry Site in Lackawanna, New York.

### 1.2 Site Background

The Lackawanna Foundry is located on Madison Avenue off of State Route 5 in Lackawanna, New York. The 3-acre site is an old abandoned foundry. Currently, the site is being remediated by the U.S. EPA. The property is bordered by smoke creek to the north, an open field to the south, a rail road to the east, and residences to the west. A building constructed of metal, brick and wood occupies the south end of the site. The rest of the site is made up of sand piles and uneven ground which is loosely occupied by grass and shrubs. Soil sampling was conducted at 27 locations specified by the U.S. EPA work assignment manager (WAM) to identify and quantitate potential contamination by the target compounds.

## 2.0 METHODOLOGY

### 2.1 Surface Soil Sampling

The top layer of soil or debris were removed with a dedicated stainless steel spoon. Using another dedicated stainless steel spoon, a thin layer of soil was collected and the sample was placed into the appropriate sample jar. Samples and locations were documented on field sample data sheets and a site map provided in Figure 1.

### 2.2 Subsurface Soil Sampling

A bucket auger was used to bore a 3-foot hole. The auger was removed from the boring and soil from the tip of the auger was placed into the appropriate sample jar. Samples and locations were documented on field sample data sheets and a site map.

### 2.3 Composite Surface Sampling

A stainless steel trowel and mixing bucket were used to collect composite surface samples within the building. The samples were mixed, then placed into the appropriate sample jar. Samples and locations were documented on field data sheets and a site map.

### 2.4 Analytical Methods

VOCs were analyzed according to SW-846, U.S. EPA Method 8260, "*Volatile Compounds By Gas Chromatography/Mass Spectrometry (GS/MS): Capillary Column Technique*". Pesticides/PCBs were analyzed according to SW-846, U.S. EPA Method 8080, "*Organochlorine Pesticides And PCBs*". BNAs were analyzed according to SW-846, U.S. EPA Method 8270, "*GC/MS For Semivolatile Organics: Capillary Column Technique*". TAL Metals were analyzed according to SW-846, U.S. EPA Method 3050, "*Acid Digestion Of Sediments, Sludges, Soils*". Cyanides were analyzed according to SW-846, U.S. EPA Method 9012, "*Total And Amenable Cyanide (colorimetric, Automated UV)*".



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### 3.0 RESULTS

#### 3.1 VOCs

Methylene chloride was detected below the method detection limit (MDL) in the method blanks. The only measurable VOC compound detected above the MDL was m&p-xylene at 13 micrograms per kilogram ( $\mu\text{g}/\text{kg}$ ). This concentration is well below the New York State Department of Environmental Conservation (NYSDEC) Technical & Administrative Guidance Memorandum (TAGM) (Appendix A) soil cleanup objectives and levels of 1200  $\mu\text{g}/\text{kg}$ . It was found within the building at sampling location 10.

#### 3.2 Pesticides/PCBs

Concentrations of p,p'-DDE, dieldrin, p,p'-DDT, aroclor 1242, aroclor 1248, aroclor 1254, and aroclor 1260, exceeding the MDL, were found at various sampling locations throughout the site. The highest concentration found was at location 27 (aroclor 1248 at 39,000  $\mu\text{g}/\text{kg}$ ) in sample A16679. This sample was a composite sample collected inside the building. Summarized results compared to the NYSDEC-TAGM soil cleanup objectives and levels can be found in table 1.

#### 3.3 BNAs

Concentrations of semi-volatile compounds were detected at various sampling locations. Phenol in particular, was detected at relatively high concentrations in 20 of the 27 sampling locations. The highest concentration was found at location 2-D. At this location, phenol was detected at 23,000  $\mu\text{g}/\text{kg}$  in sample A16652. Most of the phenol concentrations detected were above the NYSDEC-TAGM cleanup levels. Various identified and unidentified non-target compounds were also detected. Summarized results compared to the NYSDEC-TAGM soil cleanup objectives and levels can be found in Table 2.

#### 3.4 TAL Metals

All metals analyzed were detected above their MDLs with the exception of silver and thallium. Arsenic was detected at 77 mg/kg at location 22 (sample number C16672). Cadmium was detected at 20 mg/kg at location 6 (sample number C16656). Chromium was detected at 190 mg/kg in location 6 (sample number C16656). Lead was detected at 2100 mg/kg at location 7 (sample number 16657). Summarized results compared to listed NYSDEC-TAGM soil cleanup objectives and levels can be found in Table 3.

#### 3.5 Cyanides

No cyanide concentrations were detected above the MDL in any of the samples.

### 4.0 DISCUSSION OF RESULTS

#### 4.1 VOCs

VOC concentrations identified in the samples on site were all below the NYSDEC-TAGM soil clean-up level.

#### 4.2 Pesticides/PCBs

Pesticides concentrations identified in the samples collected on site, were all below the NYSDEC-TAGM clean up levels. PCB weathering indicates that the aroclors in question are present, but due to breakdown, most predominant peaks changed peak ratios. PCBs were detected above the NYDEC-TAGM criteria at location 2 and 2-D. PCBs were present at the surface and subsurface, however concentrations were below NYDEC-TAGM. PCBs were detected above the NYDEC-TAGM at locations 6, 9, 11, 12, 13, 14, 19, and location 27. PCB concentrations were 39 times greater than NYDEC-TAGM levels for surface soil at location 27, however, this sample was collected within the building. The NYDEC-TAGM refers to soil clean up levels.

#### 4.3 BNAs

BNAs concentrations identified in the samples collected on site exceeded the NYSDEC-TAGM clean up levels. Phenol in particular, was detected well above the 30 µg/kg guideline at locations 2, 2-D, 8, 9, 10, 11, 12, 13, 14, 18, 19, 20, 24, and 27. Benz(a)anthracene, chrysene, benzo(b)fluoranthene, benzo(k)fluoranthene, indeno(1,2,3,-cd)pyrene, and dibenzo(a,h)anthracene were also detected above the NYSDEC-TAGM criteria.

#### 4.4 TAL Metals

Arsenic, beryllium, cadmium, chromium, copper, iron, mercury, nickel, and zinc were detected above the NYSDEC-TAGM criteria. The memorandum states that measured levels should be compared to site background levels. According to the results, antimony, silver and thallium do not seem to be a concern at the site.

#### 4.5 Cyanides

No cyanides were detected above the MDL, hence they are not a concern in the sampling locations.

Table I  
Summarized Pesticide/PCB Sampling Results  
Lackawanna Foundry Site  
Lackawanna, New York  
May 1999

| Sample Number<br>Location | TAGM<br>Conc.<br>µg/kg | A16650<br>1    |              | A16651<br>2    |              | A16652<br>2D   |              | A16653<br>3    |              | A16654<br>4    |              | A16655<br>5    |              | A16656<br>6    |              | A16657<br>7    |              |
|---------------------------|------------------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|
|                           |                        | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg |
| Heptachlor                | 100                    | ND             | 5.8          | ND             | 3.5          | ND             | 3.5          | ND             | 4.6          | ND             | 4.1          | ND             | 4.1          | ND             | 3.9          | ND             | 4.4          |
| p,p'-DDE                  | 2100                   | 12             | 5.8          | ND             | 3.5          | ND             | 3.5          | 30             | 4.6          | 19             | 4.1          | 0.9 J          | 4.1          | 12             | 3.9          | 4.5            | 4.4          |
| Dieldrin                  | 44                     | ND             | 5.8          | ND             | 3.5          | ND             | 3.5          | ND             | 4.6          | ND             | 4.1          | ND             | 4.1          | ND             | 3.9          | 2.9 J          | 4.4          |
| p,p'-DDD                  | 2900                   | ND             | 5.8          | ND             | 3.5          | ND             | 3.7          | ND             | 4.6          | 6.9            | 4.1          | ND             | 4.1          | ND             | 3.9          | ND             | 4.4          |
| Endosulfan (II)           | 900                    | ND             | 5.8          | ND             | 3.5          | ND             | 3.5          | ND             | 4.6          | ND             | 4.1          | ND             | 4.1          | ND             | 3.9          | ND             | 4.4          |
| p,p'-DDT                  | 2100                   | 25             | 5.8          | ND             | 3.5          | ND             | 3.5          | 14             | 4.6          | 6.4            | 4.1          | ND             | 4.1          | ND             | 3.9          | 9.6            | 4.4          |
| Aroclor 1242              | Surf.<br>1000          | ND             | 73           | ND             | 3.5          | ND             | 46           | ND             | 57           | ND             | 51           | ND             | 52           | ND             | 3.9          | ND             | 55           |
| Aroclor 1248              | Surf.<br>10000         | ND             | 73           | 560            | 44           | 2700           | 46           | ND             | 57           | 120            | 51           | 15 J           | 52           | 770            | 49           | 590            | 55           |
| Aroclor 1254              | Sub.<br>10000          | ND             | 73           | ND             | 44           | ND             | 46           | 25 J           | 57           | 81             | 51           | 16JW           | 52           | 1200           | 49           | 200            | 55           |
| Aroclor 1260              | Sub.<br>10000          | ND             | 73           | ND             | 44           | ND             | 46           | 19 J           | 57           | 31 J           | 51           | 13 J           | 52           | 680            | 49           | 67             | 55           |

MDL - Method Detection Limit  
Conc. - Concentration  
µg/kg - micrograms per kilogram  
ND - Not Detected  
J - Detected Below the MDL

TAGM - New York State Department of Environmental Conservation-Technical & Administrative Guidance Memorandum  
Surf. - Surface  
Sub. - Subsurface  
NA - Not Applicable  
W - Weathering  
Shading indicates above NYSDEC-TAGM

Table I(Continued)  
Summarized Pesticide/PCB Sampling Results  
Lackawanna Foundry Site  
Lackawanna, New York  
May 1999

| Sample Number<br>Location | TAGM<br>Conc.<br>µg/kg | A16658<br>8    |              | A16659<br>9    |              | A16660<br>10   |              | A16661<br>11   |              | A16662<br>12   |              | A16663<br>13   |              | A16664<br>14   |              | A16665<br>15   |              |
|---------------------------|------------------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|
|                           |                        | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg |
| Heptachlor                | 100                    | ND             | 3.8          | ND             | 3.7          | ND             | 3.4          | ND             | 3.9          | ND             | 4.2          | ND             | 3.7          | ND             | 4.5          | ND             | 5.3          |
| p,p'-DDE                  | 2100                   | 2.5 J          | 3.8          | 17             | 3.7          | 4.7            | 3.4          | 3.9 J          | 3.9          | 28             | 4.2          | 13             | 3.7          | 93             | 4.5          | ND             | 5.3          |
| Dieldrin                  | 44                     | 2.0 J          | 3.8          | ND             | 3.7          | 2.7 J          | 3.4          | 1.1 J          | 3.9          | 6.3            | 4.2          | 2.6 J          | 3.7          | 2.4 J          | 4.5          | ND             | 5.3          |
| p,p'-DDD                  | 2900                   | ND             | 3.8          | ND             | 3.7          | ND             | 3.4          | ND             | 3.9          | ND             | 4.2          | ND             | 3.7          | ND             | 4.5          | ND             | 5.3          |
| Endosulfan (II)           | 900                    | ND             | 3.8          | ND             | 3.7          | ND             | 3.4          | ND             | 3.9          | ND             | 4.2          | ND             | 3.7          | ND             | 4.5          | ND             | 5.3          |
| p,p'-DDT                  | 2100                   | 5.3            | 3.8          | ND             | 3.7          | ND             | 3.4          | 5.2            | 3.9          | ND             | 4.2          | ND             | 3.7          | ND             | 4.5          | ND             | 5.3          |
| Aroclor 1242              | Surf.<br>1000          | ND             | 48           | ND             | 3.7          | 6500           | 42           | ND             | 49           | ND             | 52           | ND             | 46           | ND             | 4.5          | ND             | 5.3          |
| Aroclor 1248              |                        | 68             | 48           | 3500           | 46           | ND             | 42           | 1000           | 49           | 7400           | 52           | 3100           | 46           | 2800           | 57           | 42 J           | 67           |
| Aroclor 1254              | Sub.<br>10000          | 68             | 48           | 1300           | 46           | ND             | 42           | ND             | 49           | 81             | 52           | 1300           | 46           | ND             | 57           | ND             | 67           |
| Aroclor 1260              |                        | 32 J           | 48           | 320            | 46           | ND             | 42           | ND             | 49           | 31 J           | 52           | 240            | 46           | ND             | 57           | ND             | 67           |

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Surf - Surface

Sub - Subsurface

NA - Not Applicable

Shading indicates above TAGM

Table 1(Continued)  
Summarized Pesticide/PCB Sampling Results  
Lackawanna Foundry Site  
Lackawanna, New York  
May 1999

| Sample Number<br>Location | TAGM<br>Conc.<br>µg/kg | A16666<br>16   |              | A16667<br>17   |              | A16668<br>18   |              | A16669<br>19   |              | A16670<br>20   |              | A16671<br>21   |              | A16672<br>22   |              | A16673<br>23   |              |
|---------------------------|------------------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|
|                           |                        | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg |
| g-BHC                     | 300                    | ND             | 3.7          | ND             | 3.8          | ND             | 3.6          | ND             | 3.6          | 2.5 J          | 3.7          | ND             | 6.9          | ND             | 4.1          | ND             | 4.1          |
| Heptachlor                | 100                    | ND             | 3.7          | ND             | 3.8          | ND             | 3.6          | ND             | 3.6          | ND             | 3.7          | ND             | 6.9          | ND             | 4.1          | ND             | 4.1          |
| p,p'-DDE                  | 2100                   | 0.7 J          | 3.7          | 2.2 J          | 3.8          | ND             | 3.6          | 5.1            | 3.6          | ND             | 3.7          | 8.4            | 6.9          | 2.3 J          | 4.1          | 2.5 J          | 4.1          |
| Dieldrin                  | 44                     | ND             | 3.7          | 0.6 J          | 3.8          | ND             | 3.6          | 2.2 J          | 3.6          | ND             | 3.7          | 21             | 6.9          | ND             | 4.1          | ND             | 4.1          |
| p,p'-DDD                  | 2900                   | ND             | 3.7          | ND             | 3.8          | ND             | 3.6          | ND             | 3.6          | ND             | 3.7          | ND             | 6.9          | 1.6 J          | 4.1          | ND             | 4.1          |
| Endosulfan (II)           | 900                    | ND             | 3.7          | ND             | 3.8          | ND             | 3.6          | ND             | 3.6          | ND             | 3.7          | ND             | 6.9          | ND             | 4.1          | ND             | 4.1          |
| p,p'-DDT                  | 2100                   | ND             | 3.7          | ND             | 3.8          | ND             | 3.6          | 4.9            | 3.6          | ND             | 3.7          | ND             | 6.9          | 1.9 J          | 4.1          | ND             | 4.1          |
| Aroclor 1242              | Surf.<br>1000          | 340            | 47           | ND             | 3.8          | ND             | 45           | 2500           | 45           | ND             | 46           | ND             | 86           | ND             | 4.1          | ND             | 51           |
| Aroclor 1248              |                        | ND             | 47           | 620            | 47           | 77             | 45           | ND             | 45           | ND             | 46           | 320            | 86           | 11JW           | 51           | 110            | 51           |
| Aroclor 1254              | Sub.<br>10000          | ND             | 47           | ND             | 47           | 33 J           | 45           | ND             | 45           | ND             | 46           | 300            | 86           | 16JW           | 51           | ND             | 51           |
| Aroclor 1260              |                        | ND             | 47           | ND             | 47           | ND             | 45           | ND             | 45           | 3900           | 46           | 200            | 86           | ND             | 51           | 14 J           | 51           |

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Sub - Subsurface

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W - Weathering

Shading indicates above TAGM

Table 1(Continued)  
 Summarized Pesticide/PCB Sampling Results  
 Lackawanna Foundry Site  
 Lackawanna, New York  
 May 1999

| Sample Number<br>Location | TAGM          | A16674<br>24   |              | A16675<br>25   |              | A16676<br>25D  |              | A16677<br>26   |              | A16678<br>26-D |              | A16679<br>27   |              |
|---------------------------|---------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|
|                           |               | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg |
| g-BHC                     | 300           | ND             | 4.4          | ND             | 4.0          | ND             | 4.3          | ND             | 4.8          | ND             | 4.4          | ND             | 4.0          |
| Heptachlor                | 100           | ND             | 4.4          | ND             | 4.0          | ND             | 4.3          | ND             | 4.8          | ND             | 4.4          | ND             | 4.0          |
| p,p'-DDE                  | 2100          | 1.9 J          | 4.4          | 3.3 J          | 4.0          | ND             | 4.3          | 11             | 4.8          | ND             | 4.4          | ND             | 4.0          |
| Dieldrin                  | 44            | ND             | 4.4          | ND             | 4.0          | ND             | 4.3          | ND             | 4.8          | ND             | 4.4          | ND             | 4.0          |
| p,p'-DDD                  | 2900          | ND             | 4.4          | ND             | 4.0          | ND             | 4.3          | ND             | 4.8          | ND             | 4.4          | ND             | 4.0          |
| Endosulfan (II)           | 900           | ND             | 4.4          | ND             | 4.0          | ND             | 4.3          | ND             | 4.8          | ND             | 4.4          | ND             | 4.0          |
| p,p'-DDT                  | 2100          | ND             | 4.4          | ND             | 4.0          | ND             | 4.3          | 18             | 4.8          | ND             | 4.4          | ND             | 4.0          |
| Aroclor 1242              | Surf.<br>1000 | 340            | 55           | ND             | 4.0          | ND             | 54           | ND             | 60           | ND             | 54           | ND             | 50           |
| Aroclor 1248              |               | 120            | 55           | 270            | 50           | ND             | 54           | 170            | 60           | ND             | 54           | 39000          | 50           |
| Aroclor 1254              | Sub.<br>10000 | ND             | 55           | 83             | 50           | ND             | 54           | ND             | 60           | ND             | 54           | ND             | 50           |
| Aroclor 1260              |               | 10 J           | 55           | ND             | 50           | ND             | 54           | 42 J           | 60           | 3900           | 54           | ND             | 50           |

MDL - Method Detection Limit

Conc. - Concentration

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Shading indicates above TAGM

Table 2  
Summarized BNAs Sampling Results  
Lackawanna Foundry Site  
Lackawanna, New York  
May 1999

| Sample Number<br>Location  | TAGM           | A166501        | A16651<br>2    | A16652<br>2-D  | A16653<br>3    | A16654<br>4    | A16655<br>5    | A16656<br>6    | A16657<br>7    |
|----------------------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|
| Analyte                    | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg |
| Phenol                     | 30             | ND             | 16000          | 23000          | ND             | ND             | ND             | ND             | 1900 J         |
| 2-Methylphenol             | 100            | ND             | 370            | 410 J          | ND             | ND             | ND             | ND             | ND             |
| Naphthalene                | 13000          | ND             | 970            | 960 J          | ND             | ND             | 310 J          | ND             | ND             |
| 2-Methylnaphthalene        | 36400          | ND             | 360            | 640 J          | ND             | ND             | 270 J          | ND             | ND             |
| Acenaphthylene             | 41000          | ND             | ND             | ND             | ND             | ND             | 200 J          | ND             | ND             |
| Acenaphthene               | 50000          | ND             | 95 J           | ND             | ND             | ND             | 740            | ND             | ND             |
| Dibenzofuran               | 6200           | ND             | 100 J          | ND             | ND             | ND             | 490            | ND             | ND             |
| Fluorene                   | 50000          | ND             | 87 J           | ND             | ND             | ND             | 880            | ND             | ND             |
| Phenanthrene               | 50000          | ND             | 1400           | ND             | ND             | ND             | 7900           | ND             | ND             |
| Anthracene                 | 50000          | ND             | 220 J          | ND             | ND             | ND             | 1800           | ND             | ND             |
| Carbazole                  | NA             | ND             | 190 J          | ND             | ND             | ND             | 1200           | ND             | ND             |
| Di-n-butylphthalate        | 8100           | ND             | 150 J          | ND             | ND             | ND             | ND             | ND             | ND             |
| Fluoranthene               | 50000          | ND             | 2000           | 380 J          | 1000 J         | 2300 J         | 9900           | 2000 J         | 2400 J         |
| Pyrene                     | 50000          | ND             | 1500           | ND             | ND             | 1900 J         | 7800           | 1700 J         | 2200 J         |
| Benzo(a)anthracene         | 224            | ND             | 800            | ND             | ND             | 1100 J         | 5000           | 1000 J         | 1500 J         |
| Chrysene                   | 400            | ND             | 920            | ND             | ND             | 1400 J         | 5300           | 1100 J         | 1900 J         |
| Bis(2-Ethylhexyl)phthalate | 50000          | ND             | 640            | 1300 J         | 1000 J         | 1000 J         | 260 J          | 810 J          | ND             |
| Benzo(b)fluoranthene       | 1100           | ND             | 1100           | ND             | ND             | 1700 J         | 5800           | 1400 J         | 3100 J         |
| Benzo(k)fluoranthene       | 1100           | ND             | 880            | ND             | ND             | 1400 J         | 4500           | 1200 J         | 2400 J         |
| Benzo(a)pyrene             | 61             | ND             | 1,100          | ND             | ND             | 1700 J         | 5800           | 1300 J         | 3300 J         |
| Indeno(1,2,3,-cd)pyrene    | 3200           | ND             | 750            | ND             | ND             | 1,300 J        | 3400           | 990 J          | 3200 J         |
| Dibenzo(a,h)anthracene     | 14             | ND             | 300 J          | ND             | ND             | ND             | 1300           | ND             | 1200 J         |
| Benzo(g,h,i)perylene       | 50000          | ND             | 890            | ND             | ND             | 1600 J         | 3700           | 1100 J         | 3600 J         |

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Table 2 (Continued)  
Summarized BNAs Sampling Results  
Lackawanna Foundry Site  
Lackawanna, New York  
May 1999

| Sample Number<br>Location  | TAGM           | A16658<br>8    | A16659<br>9    | A16660<br>10   | A16661<br>11   | A16662<br>12   | A16663<br>13   | A16664<br>14   | A16665<br>15   |
|----------------------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|
| Analyte                    | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg |
| Phenol                     | 30             | 1500           | 4900           | 1800           | 9700           | 4700           | 2200           | 7000           | 1600           |
| 2-Methylphenol             | 100            | ND             | ND             | 570 J          | ND             | ND             | ND             | ND             | ND             |
| Naphthalene                | 13000          | ND             | ND             | 3100           | 2100 J         | 960 J          | 670 J          | 1400 J         | ND             |
| 2-Methylnaphthalene        | 36400          | ND             | ND             | 1500 J         | ND             | 500 J          | 400 J          | ND             | ND             |
| Acenaphthylene             | 41000          | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Acenaphthene               | 50000          | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Dibenzofuran               | 6200           | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Fluorene                   | 50000          | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Phenanthrene               | 50000          | ND             | ND             | 680 J          | ND             | 530 J          | ND             | ND             | ND             |
| Anthracene                 | 50000          | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Carbazole                  | NA             | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Di-n-butylphthalate        | 8100           | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Fluoranthene               | 50000          | 1300 J         | 1000 J         | ND             | ND             | 600 J          | ND             | ND             | ND             |
| Pyrene                     | 50000          | 1200 J         | 880 J          | ND             | ND             | 900 J          | ND             | ND             | ND             |
| Butylbenzylphthalate       | 50000          | ND             | 6000           | ND             | ND             | ND             | ND             | ND             | ND             |
| Benzo(a)anthracene         | 224            | ND             | NA             | ND             | ND             | ND             | ND             | 460 J          | ND             |
| Chrysene                   | 400            | 1000 J         | NA             | ND             | ND             | ND             | ND             | ND             | ND             |
| Bis(2-Ethylhexyl)phthalate | 50000          | 8700           | 1200 J         | 1300 J         | 16000          | 1300 J         | 540 J          | ND             | ND             |
| Benzo(b)fluoranthene       | 1100           | 1600 J         | 940 J          | ND             | ND             | ND             | ND             | ND             | ND             |
| Benzo(k)fluoranthene       | 1100           | 1400 J         | 830 J          | ND             | ND             | ND             | ND             | ND             | ND             |
| Benzo(a)pyrene             | 61             | 1400 J         | 920 J          | ND             | ND             | ND             | ND             | ND             | ND             |
| Indeno(1,2,3,-cd)pyrene    | 3200           | 1700 J         | 770 J          | ND             | ND             | ND             | ND             | ND             | ND             |
| Dibenzo(a,h)anthracene     | 14             | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Benzo(g,h,i)perylene       | 50000          | 1900 J         | 920 J          | ND             | ND             | ND             | ND             | ND             | ND             |

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Table 2 (Continued)  
Summarized BNAs Sampling Results  
Lackawanna Foundry Site  
Lackawanna, New York  
May 1999

| Sample Number<br>Location  | TAGM           | A16666<br>16   | A16667<br>17   | A16668<br>18   | A16669<br>19   | A16670<br>20   | A16671<br>21   | A16672<br>22   | A16673<br>23   |
|----------------------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|
| Analyte                    | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg |
| Phenol                     | 30             | 840 J          | 910 J          | <b>11000</b>   | <b>4000</b>    | <b>18000</b>   | 5600 J         | ND             | 2400 J         |
| 2-Methylphenol             | 100            | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Naphthalene                | 13000          | ND             | ND             | 850 J          | 580            | 4400           | ND             | ND             | 1100 J         |
| 2-Methylnaphthalene        | 36400          | ND             | ND             | ND             | 100 J          | 1100 J         | ND             | ND             | 850 J          |
| Acenaphthylene             | 41000          | ND             | ND             | ND             | ND             | ND             | ND             | 420 J          | 1300 J         |
| Acenaphthene               | 50000          | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Dibenzofuran               | 62000          | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Fluorene                   | 50000          | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Phenanthrene               | 50000          | ND             | ND             | ND             | ND             | ND             | ND             | ND             | 2200 J         |
| Anthracene                 | 50000          | ND             | ND             | ND             | ND             | ND             | ND             | ND             | 1400 J         |
| Carbazole                  | NA             | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Di-n-butylphthalate        | 8100           | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Fluoranthene               | 50000          | ND             | 860 J          | ND             | 74 J           | 1600 J         | 3400 J         | 4700           | 8500           |
| Pyrene                     | 50000          | ND             | 740 J          | ND             | ND             | 1400 J         | 3400 J         | 4100           | 8800           |
| Butylbenzylphthalate       | 224            | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Benzo(a)anthracene         | 400            | ND             | 410 J          | ND             | ND             | 840 J          | 2100 J         | <b>2600</b>    | <b>5200</b>    |
| Chrysene                   | 400            | ND             | 510 J          | ND             | ND             | 900 J          | 2600 J         | <b>2400</b>    | <b>5400</b>    |
| Bis(2-Ethylhexyl)phthalate | 50000          | ND             | ND             | ND             | 180 J          | 760 J          | 1700 J         | 430 J          | ND             |
| Benzo(b)fluoranthene       | 1100           | ND             | 580 J          | ND             | ND             | 910 J          | 3500 J         | <b>2700</b>    | <b>5700</b>    |
| Benzo(k)fluoranthene       | 1100           | ND             | 520 J          | ND             | ND             | 890 J          | 3400 J         | <b>2500</b>    | <b>5700</b>    |
| Benzo(a)pyrene             | 61             | ND             | 570 J          | ND             | ND             | 980 J          | 3700 J         | <b>3100</b>    | <b>5300</b>    |
| Indeno(1,2,3,-cd)pyrene    | 3200           | ND             | 420 J          | ND             | ND             | 660 J          | 3100 J         | 1600 J         | 3100 J         |
| Dibenzo(a,h)anthracene     | 14             | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Benzo(g,h,i)perylene       | 50000          | ND             | 470 J          | ND             | ND             | 770 J          | 3600 J         | 1700 J         | 3200 J         |

MDL - Method Detection Limit

Conc. - Concentration

µg/kg - micrograms per kilogram

ND - Not Detected

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NA - Not Applicable, Shading indicates above TAGM

Table 2 (Continued)  
Summarized BNAs Sampling Results  
Lackawanna Foundry Site  
Lackawanna, New York  
May 1999

| Sample Number<br>Location  | TAGM           | A16674<br>24   | A16675<br>25   | A16676<br>25-D | A16677<br>26   | A16678<br>26-D | A16679<br>27   |
|----------------------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|
| Analyte                    | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg | Conc.<br>µg/kg |
| Phenol                     | 30             | <b>4000</b>    | 990 J          | ND             | ND             | ND             | <b>16000</b>   |
| 2-Methylphenol             | 100            | ND             | ND             | ND             | ND             | ND             | 820 J          |
| 1,2,4-Trichlorobenzene     | NA             | ND             | ND             | ND             | ND             | ND             | 1700 J         |
| Naphthalene                | 13000          | 1400 J         | ND             | ND             | ND             | ND             | 2900           |
| 2-Methylnaphthalene        | 36400          | 790 J          | ND             | ND             | ND             | ND             | 810 J          |
| Acenaphthylene             | 41000          | ND             | ND             | ND             | ND             | 96 J           | ND             |
| Acenaphthene               | 50000          | ND             | ND             | ND             | ND             | ND             | ND             |
| Dibenzofuran               | 6200           | ND             | ND             | ND             | ND             | ND             | ND             |
| Fluorene                   | 50000          | ND             | ND             | ND             | ND             | ND             | ND             |
| Phenanthrene               | 50000          | ND             | ND             | ND             | ND             | 440            | ND             |
| Anthracene                 | 50000          | ND             | ND             | ND             | ND             | 92 J           | ND             |
| Carbazole                  | NA             | ND             | ND             | ND             | ND             | ND             | ND             |
| Di-n-butylphthalate        | 8100           | ND             | ND             | ND             | ND             | ND             | ND             |
| Fluoranthene               | 50000          | 1200 J         | 5800           | 100 J          | 3300 J         | 930            | 420 J          |
| Pyrene                     | 50000          | 940 J          | 4800           | ND             | 2900 J         | 810            | 500 J          |
| Butylbenzylphthalate       | 50000          | ND             | ND             | ND             | ND             | ND             | ND             |
| Benzo(a)anthracene         | 224            | 650 J          | 2700 J         | ND             | 1600 J         | <b>470</b>     | ND             |
| Chrysene                   | 400            | 910 J          | 3100 J         | ND             | 2000 J         | <b>500</b>     | ND             |
| Bis(2-Ethylhexyl)phthalate | 50000          | ND             | ND             | ND             | 1100 J         | ND             | 1000 J         |
| Benzo(b)fluoranthene       | 1100           | 1200 J         | 3600 J         | ND             | 2100 J         | 540            | ND             |
| Benzo(k)fluoranthene       | 1100           | 1000 J         | 3200 J         | ND             | 1900 J         | 520            | ND             |
| Benzo(a)pyrene             | 61             | 1100 J         | 3700 J         | ND             | 2200 J         | <b>600</b>     | ND             |
| Indeno(1,2,3,-cd)pyrene    | 3200           | 860 J          | 2400 J         | ND             | 1400 J         | 360 J          | ND             |
| Dibenzo(a,h)anthracene     | 14             | ND             | ND             | ND             | ND             | ND             | ND             |
| Benzo(g,h,i)perylene       | 50000          | 970 J          | 2700 J         | ND             | ND             | 400 J          | ND             |

TAGM - New York State Department of Environmental Conservation-Technical & Administrative Guidance Memorandum

Conc. - Concentration

µg/kg - micrograms per kilogram

ND - Not Detected

J - Detected Below the MDL

MDL - Method Detection Limit, Shading indicates above TAGM

Table 3  
Summarized TAL Metals Sampling Results  
Lackawanna Foundry Site  
Lackawanna, New York  
May 1999

| Sample Number<br>Location | TAGM           | C16650<br>1    | C16651<br>2    | C16652<br>2-D  | C16653<br>3    | C16654<br>4    | C16655<br>5    | C16656<br>6    | C16657<br>7    |
|---------------------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|
| Analyte                   | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg |
| Aluminum                  | *              | 11000          | 1300           | 1200           | 7800           | 13000          | 12000          | 4600           | 5300           |
| Antimony                  | *              | ND             | ND             | ND             | ND             | ND             | ND             | 6.4            | ND             |
| Arsenic                   | 7.5*           | <b>14</b>      | 1.6            | 2.5            | <b>11</b>      | <b>13</b>      | <b>15</b>      | <b>20</b>      | 6.3            |
| Barium                    | 300*           | 160            | 18             | 17             | 140            | 150            | 130            | 270            | 120            |
| Beryllium                 | 0.16*          | <b>0.93</b>    | ND             | ND             | <b>0.81</b>    | <b>2.2</b>     | <b>0.85</b>    | ND             | ND             |
| Cadmium                   | 1*             | <b>3.2</b>     | 0.48           | 0.66           | <b>3.7</b>     | <b>2.4</b>     | <b>2.4</b>     | <b>20</b>      | <b>3.6</b>     |
| Calcium                   | *              | 21000          | 1900           | 3300           | 94000          | 120000         | 14000          | 21000          | 18000          |
| Chromium                  | 10*            | <b>45</b>      | <b>25</b>      | <b>15</b>      | <b>33</b>      | <b>28</b>      | <b>29</b>      | <b>190</b>     | <b>55</b>      |
| Cobalt                    | 30*            | 12             | 2.1            | 1.2            | 7.7            | 6.2            | 14             | 20             | 6.7            |
| Copper                    | 25*            | 74             | 57             | 35             | 75             | 67             | 200            | 440            | 110            |
| Iron                      | 2000*          | <b>43000</b>   | <b>24000</b>   | <b>13000</b>   | <b>46000</b>   | <b>30000</b>   | <b>35000</b>   | <b>200000</b>  | <b>47000</b>   |
| Lead                      | *              | 300            | 44             | 38             | 200            | 260            | 180            | 610            | 2100           |
| Magnesium                 | *              | 4800           | 600            | 560            | 7200           | 11000          | 4800           | 3100           | 3300           |
| Manganese                 | *              | 1200           | 250            | 140            | 800            | 1300           | 700            | 1800           | 1100           |
| Mercury                   | 0.1            | <b>0.21</b>    | 0.06           | <b>0.1</b>     | <b>0.11</b>    | <b>0.16</b>    | <b>0.11</b>    | <b>1.1</b>     | <b>0.22</b>    |
| Nickel                    | 13*            | <b>42</b>      | <b>15</b>      | 9.2            | <b>28</b>      | <b>32</b>      | <b>41</b>      | <b>160</b>     | <b>110</b>     |
| Potassium                 | *              | 1500           | ND             | ND             | 920            | 1500           | 1200           | 630            | 570            |
| Selenium                  | 2*             | 1.0            | ND             | ND             | 0.79           | 0.98           | 0.58           | 0.77           | ND             |
| Silver                    | *              | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Sodium                    | *              | 110            | ND             | 64             | 140            | 520            | 79             | 410            | 230            |
| Thallium                  | *              | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Vanadium                  | 150*           | 27             | 5.2            | 3.7            | 17             | 18             | 23             | 20             | 20             |
| Zinc                      | 20*            | <b>760</b>     | <b>140</b>     | <b>79</b>      | <b>640</b>     | <b>540</b>     | <b>340</b>     | <b>1100</b>    | <b>1000</b>    |

\* - See TAGM Memorandum Appendix A

MDL - Method Detection Limit

Conc. - Concentration

µg/kg - micrograms per kilogram

ND - Not Detected

J - Detected Below the MDL

TAGM - New York State Department of Environmental Conservation-Technical & Administrative Guidance Memorandum

Shading indicates above TAGM

Table 3 (Continued)  
Summarized TAL Metals Sampling Results  
Lackawanna Foundry Site  
Lackawanna, New York  
May 1999

| Sample Number<br>Location | TAGM           | C16658<br>8    | C16659<br>9    | C16660<br>10   | C16661<br>11   | C16662<br>12   | C16663<br>13   | C16664<br>14   | C16665<br>15   |
|---------------------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|
| Analyte                   | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg |
| Aluminum                  | *              | 3900           | 5800           | 2100           | 740            | 4200           | 4600           | 2300           | 2200           |
| Antimony                  | *              | ND             | ND             | ND             | ND             | ND             | 13             | ND             | ND             |
| Arsenic                   | 7.5*           | 6.5            | <b>16</b>      | 1.3            | 1.8            | 7.0            | <b>16</b>      | 4.4            | 2.7            |
| Barium                    | 300*           | 83             | 170            | 22             | 14             | 84             | 100            | 39             | 50             |
| Beryllium                 | 0.16*          | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Cadmium                   | 1*             | <b>2.4</b>     | <b>4.4</b>     | 0.59           | <b>4.2</b>     | <b>2.0</b>     | <b>3.4</b>     | 0.92           | 0.48           |
| Calcium                   | *              | 8300           | 14000          | 2200           | 1100           | 21000          | 1200           | 24000          | 1600           |
| Chromium                  | 10*            | <b>36</b>      | <b>63</b>      | <b>14</b>      | <b>13</b>      | <b>100</b>     | <b>79</b>      | <b>86</b>      | <b>14</b>      |
| Cobalt                    | 30*            | 8.5            | 8.4            | 2.0            | 1.7            | 7.0            | 11             | 3.6            | 0.98           |
| Copper                    | 25*            | <b>81</b>      | <b>200</b>     | <b>41</b>      | <b>33</b>      | <b>270</b>     | <b>370</b>     | <b>93</b>      | <b>100</b>     |
| Iron                      | 2000*          | <b>73000</b>   | <b>76000</b>   | <b>19000</b>   | <b>22000</b>   | <b>56000</b>   | <b>140000</b>  | <b>39000</b>   | <b>11000</b>   |
| Lead                      | *              | 230            | 340            | 34             | 48             | 150            | 540            | 110            | 58             |
| Magnesium                 | *              | 1800           | 4300           | 610            | 240            | 2400           | 480            | 3400           | 460            |
| Manganese                 | *              | 790            | 1100           | 180            | 160            | 810            | 770            | 2500           | 120            |
| Mercury                   | 0.1            | <b>0.13</b>    | <b>0.95</b>    | <b>0.27</b>    | 0.04           | <b>0.17</b>    | 0.07           | <b>0.12</b>    | ND             |
| Nickel                    | 13*            | <b>41</b>      | <b>160</b>     | <b>15</b>      | <b>13</b>      | <b>75</b>      | <b>72</b>      | <b>25</b>      | <b>31</b>      |
| Potassium                 | *              | 540            | 390            | 500            | ND             | 430            | 500            | 260            | ND             |
| Selenium                  | 2*             | ND             | ND             | ND             | ND             | ND             | 0.39           | ND             | ND             |
| Silver                    | *              | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Sodium                    | *              | 230            | 270            | 1500           | ND             | 210            | 200            | 92             | 72             |
| Thallium                  | *              | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Vanadium                  | 150*           | 14             | 16             | 4.2            | 2.7            | 14             | 16             | 43             | 5.4            |
| Zinc                      | 20*            | <b>450</b>     | <b>550</b>     | <b>42</b>      | <b>81</b>      | <b>280</b>     | <b>660</b>     | <b>190</b>     | <b>110</b>     |

\* - See TAGM Memorandum Appendix A

MDL - Method Detection Limit

Conc. - Concentration

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J - Detected Below the MDL

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Shading indicates above TAGM

Table 3 (Continued)  
Summarized TAL Metals Sampling Results  
Lackawanna Foundry Site  
Lackawanna, New York  
May 1999

| Sample Number<br>Location | TAGM           | C16666<br>16   | C16667<br>17   | C16668<br>18   | C16669<br>19   | C16670<br>20   | C16671<br>21   | C16672<br>22   | C16673<br>23   |
|---------------------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|
| Analyte                   | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg |
| Aluminum                  | *              | 1300           | 5900           | 790            | 1400           | 2400           | 12000          | 15000          | 9300           |
| Antimony                  | *              | ND             | 6.5            | ND             | ND             | ND             | ND             | ND             | ND             |
| Arsenic                   | 7.5*           | 1.5            | 7.1            | 4.4            | 6.7            | 6.5            | <b>20</b>      | <b>77</b>      | <b>8.4</b>     |
| Barium                    | 300*           | 15             | 180            | 9.0            | 29             | 130            | 200            | 170            | 94             |
| Beryllium                 | 0.16*          | ND             | ND             | ND             | ND             | ND             | <b>0.96</b>    | <b>1.2</b>     | <b>0.66</b>    |
| Cadmium                   | 1*             | ND             | ND             | ND             | 0.59           | <b>2.3</b>     | <b>13</b>      | <b>2.3</b>     | <b>1.5</b>     |
| Calcium                   | *              | 1200           | 130000         | 1200           | 850            | 3800           | 23000          | 34000          | 12000          |
| Chromium                  | 10*            | 6.1            | <b>33</b>      | 8.9            | <b>97</b>      | <b>130</b>     | <b>95</b>      | <b>32</b>      | <b>24</b>      |
| Cobalt                    | 30*            | 1.4            | 47             | ND             | 4.4            | 12             | 15             | 11             | 8.0            |
| Copper                    | 25*            | <b>44</b>      | <b>240</b>     | <b>28</b>      | <b>300</b>     | <b>720</b>     | <b>270</b>     | <b>47</b>      | <b>52</b>      |
| Iron                      | 2000*          | <b>16000</b>   | <b>25000</b>   | <b>5800</b>    | <b>58000</b>   | <b>93000</b>   | <b>110000</b>  | <b>32000</b>   | <b>30000</b>   |
| Lead                      | *              | 58             | 300            | 21             | 140            | 280            | 660            | 180            | 160            |
| Magnesium                 | *              | 410            | 19000          | 340            | 410            | 1000           | 4700           | 5900           | 3600           |
| Manganese                 | *              | 210            | 840            | 75             | 760            | 690            | 1600           | 1200           | 560            |
| Mercury                   | 0.1            | ND             | 0.06           | 0.04           | ND             | <b>1.3</b>     | <b>0.77</b>    | <b>0.73</b>    | <b>0.13</b>    |
| Nickel                    | 13*            | 9.5            | 38             | 2.7            | 57             | 150            | 100            | 36             | 28             |
| Potassium                 | *              | ND             | 470            | ND             | ND             | 280            | 900            | 1500           | 1100           |
| Selenium                  | 2*             | ND             | ND             | ND             | ND             | ND             | 1.0            | ND             | 0.57           |
| Silver                    | *              | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Sodium                    | *              | ND             | 330            | ND             | 68             | 68             | 360            | 120            | 130            |
| Thallium                  | *              | ND             | ND             | ND             | ND             | ND             | ND             | ND             | ND             |
| Vanadium                  | 150*           | 3.1            | 11             | 2.6            | 14             | 21             | 36             | 27             | 19             |
| Zinc                      | 20*            | <b>98</b>      | <b>1300</b>    | <b>55</b>      | <b>310</b>     | <b>290</b>     | <b>1000</b>    | <b>1100</b>    | <b>660</b>     |

\* - See TAGM Memorandum Appendix A

MDL - Method Detection Limit

Conc. - Concentration

µg/kg - micrograms per kilogram

ND - Not Detected

J - Detected Below the MDL

TAGM - New York State Department of Environmental Conservation Technical & Administrative Guidance Memorandum

Shading indicates above TAGM

Table 3 (Continued)  
Summarized TAL Metals Sampling Results  
Lackawanna Foundry Site  
Lackawanna, New York  
May 1999

| Sample Number<br>Location | TAGM           | C16674<br>24   | C16675<br>25   | C16676<br>25-D | C16677<br>26   | C16678<br>26-D | C16679<br>27   |
|---------------------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|
| Analyte                   | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg | Conc.<br>mg/kg |
| Aluminum                  | *              | 7300           | 12000          | 17000          | 14000          | 14000          | 2900           |
| Antimony                  | *              | ND             | ND             | ND             | ND             | ND             | ND             |
| Arsenic                   | 7.5*           | 5              | <b>21</b>      | <b>25</b>      | <b>66</b>      | <b>48</b>      | <b>19</b>      |
| Barium                    | 300*           | 64             | 130            | 140            | 250            | 280            | 59             |
| Beryllium                 | 0.16*          | ND             | <b>0.73</b>    | <b>0.97</b>    | <b>1.3</b>     | <b>0.67</b>    | ND             |
| Cadmium                   | 1*             | 1.2            | 2.3            | ND             | 5.2            | 1.3            | 0.84           |
| Calcium                   | *              | 10000          | 39000          | 11000          | 53000          | 120000         | 15000          |
| Chromium                  | 10*            | <b>18</b>      | <b>31</b>      | <b>25</b>      | <b>51</b>      | <b>23</b>      | <b>95</b>      |
| Cobalt                    | 30*            | 4.8            | 9.7            | 15             | 9.7            | 7.3            | 8.7            |
| Copper                    | 25*            | <b>49</b>      | <b>65</b>      | <b>27</b>      | <b>110</b>     | <b>34</b>      | <b>280</b>     |
| Iron                      | 2000*          | <b>25000</b>   | <b>29000</b>   | <b>32000</b>   | <b>45000</b>   | <b>20000</b>   | <b>78000</b>   |
| Lead                      | *              | 130            | 230            | 40             | 740            | 100            | 140            |
| Magnesium                 | *              | 3500           | 5400           | 5400           | 6300           | 7300           | 2500           |
| Manganese                 | *              | 390            | 710            | 610            | 1500           | 570            | 770            |
| Mercury                   | 0.1            | <b>0.13</b>    | <b>0.17</b>    | 0.08           | <b>1.6</b>     | <b>0.17</b>    | 0.09           |
| Nickel                    | 13*            | <b>18</b>      | <b>38</b>      | <b>41</b>      | <b>39</b>      | <b>22</b>      | <b>89</b>      |
| Potassium                 | *              | 1100           | 1400           | 1500           | 1600           | 1400           | 320            |
| Selenium                  | 2*             | ND             | ND             | ND             | 1.6            | ND             | ND             |
| Silver                    | *              | ND             | ND             | ND             | ND             | ND             | ND             |
| Sodium                    | *              | 250            | 170            | ND             | 190            | 190            | 120            |
| Thallium                  | *              | ND             | ND             | ND             | ND             | ND             | ND             |
| Vanadium                  | 150*           | 14             | 23             | 27             | 29             | 20             | 15             |
| Zinc                      | 20*            | <b>350</b>     | <b>760</b>     | <b>150</b>     | <b>1600</b>    | <b>330</b>     | <b>250</b>     |

\* - See TAGM Memorandum Appendix A

MDL - Method Detection Limit

Conc. - Concentration

µg/kg - micrograms per kilogram

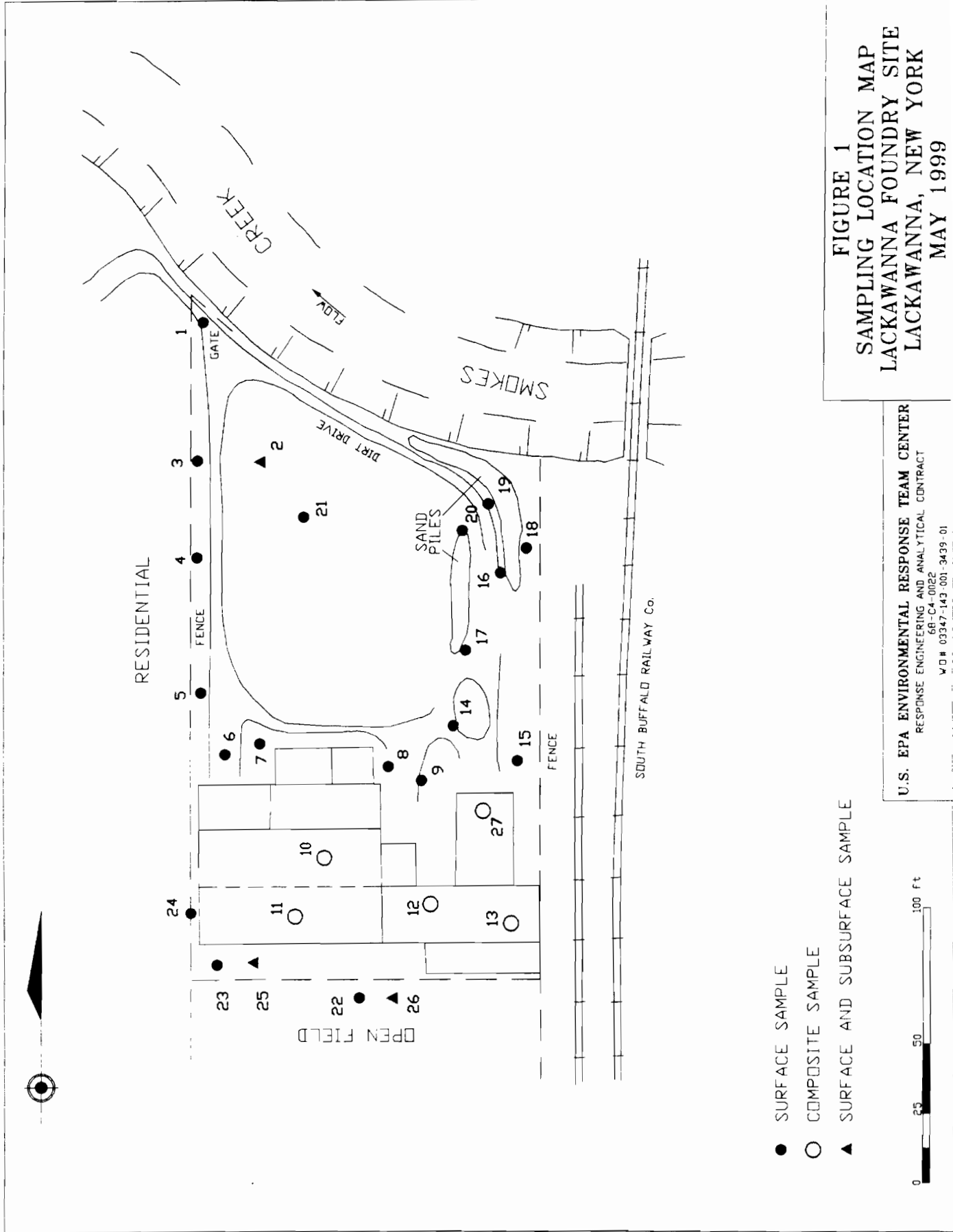
ND - Not Detected

J - Detected Below the MDL

TAGM - New York State Department of Environmental Conservation Technical & Administrative Guidance Memorandum  
Shading indicates above TAGM

# Figures





U.S. EPA ENVIRONMENTAL RESPONSE TEAM CENTER  
 RESPONSE ENGINEERING AND ANALYTICAL CONTRACT  
 68-C4-0022  
 WDB 03347-143-001-3439-01

# Appendix A



## New York State Department of Environmental Conservation

## MEMORANDUM

TO: Regional Hazardous Waste Remediation Engineers, Bureau Directors and Section Chiefs  
 FROM: Michael J. O'Toole, Jr., Director, Division of Hazardous Waste Remediation  
 SUBJECT: Revised TAGM - Determination of Soil Cleanup Objectives and Cleanup Levels

DATE: JAN 24 1994

Attached is the revised Division Technical and Administrative Guidance Memorandum (TAGM) on Determination of Soil Cleanup Objectives and Cleanup Levels in its final form. The changes are to the Tables of Appendix A. They are minor in nature and do not change the content of the TAGM. The changes include:

1. Alphabetizing contaminants in Table 1 through Table 4.
2. The addition of a few contaminants to Table 1 and Table 2.
3. Table 4 has been revised to indicate that background levels for lead vary widely and provide a range for undeveloped and developed areas. It also has been revised to indicate that site-specific form(s) of cyanide should be considered when establishing soil cleanup objectives for cyanide.

If you have any questions, please contact Ajay Shroff of my staff at (518)485-8792.

## Attachment

|                |                                      |
|----------------|--------------------------------------|
| cc: T. Jorling | J. Davis                             |
| J. Lacey       | J. Kelleher                          |
| M. Gerstman    | J. Colquhoun                         |
| A. DeBarbieri  | D. Persson                           |
| E. Sullivan    | J. Printup                           |
| T. Donovan     | M. Bermingham                        |
| C. Sullivan    | D. Johnson                           |
| J. Eckl        | M. Kadlecek                          |
| R. Davies      | Regional Directors                   |
| R. Dana        | Regional Engineers                   |
| C. Goddard     | Regional Solid and Haz. Waste Engrs. |
| C. Costopoulos | Regional Citizen Participation Spec. |
| P. Counterman  |                                      |





MEMORANDUM

TO:  
FROM:  
SUBJECT:  
DATE:

Regional Haz. Waste Remediation Engineers, Bureau Dirs. & Section Chiefs  
Michael J. O'Toole, Jr., Director, Div. of Hazardous Waste Remediation  
DIVISION TECHNICAL AND ADMINISTRATIVE GUIDANCE MEMORANDUM:  
DETERMINATION OF SOIL CLEANUP OBJECTIVES AND CLEANUP LEVELS

JAN 24 1994

The cleanup goal of the Department is to restore inactive hazardous waste sites to predisposal conditions, to the extent feasible and authorized by law. However, it is recognized that restoration to predisposal conditions will not always be feasible.

1. INTRODUCTION:

This TAGM provides a basis and procedure to determine soil cleanup levels at individual Federal Superfund, State Superfund, 1986 EQBA Title 3 and Responsible Party (RP) sites, when the Director of the DHWR determines that cleanup of a site to predisposal conditions is not possible or feasible.

The process starts with development of soil cleanup objectives by the Technology Section for the contaminants identified by the Project Managers. The Technology Section uses the procedure described in this TAGM to develop soil cleanup objectives. Attainment of these generic soil cleanup objectives will, at a minimum, eliminate all significant threats to human health and/or the environment posed by the inactive hazardous waste site. Project Managers should use these cleanup objectives in selecting alternatives in the Feasibility Study (FS). Based on the proposed selected remedial technology (outcome of FS), final site specific soil cleanup levels are established in the Record of Decision (ROD) for these sites.

It should be noted that even after soil cleanup levels are established in the ROD, these levels may prove to be unattainable when remedial construction begins. In that event, alternative remedial actions or institutional controls may be necessary to protect the environment.

2. BASIS FOR SOIL CLEANUP OBJECTIVES:

The following alternative bases are used to determine soil cleanup objectives:

- (a) Human health based levels that correspond to excess lifetime

cancer risks of one in a million for Class A<sup>1</sup> and B<sup>2</sup> carcinogens, or one in 100,000 for Class C<sup>3</sup> carcinogens. These levels are contained in USEPA's Health Effects Assessment Summary Tables (HEASTs) which are compiled and updated quarterly by the NYSDEC's Division of Hazardous Substances Regulation;

- (b) Human health based levels for systemic toxicants, calculated from Reference Doses (RfDs). RfDs are an estimate of the daily exposure an individual (including sensitive individuals) can experience without appreciable risk of health effects during a lifetime. An average scenario of exposure in which children ages one to six (who exhibit the greatest tendency to ingest soil) is assumed. An intake rate of 0.2 gram/day for a five-year exposure period for a 16-kg child is assumed. These levels are contained in USEPA's Health Effects Assessment Summary Tables (HEASTs) which are compiled and updated quarterly by the NYSDEC's Division of Hazardous Substances Regulation;
- (c) Environmental concentrations which are protective of groundwater/drinking water quality; based on promulgated or proposed New York State Standards;
- (d) Background values for contaminants; and
- (e) Detection limits.

A recommendation on the appropriate cleanup objective is based on the criterion that produces the most stringent cleanup level using criteria a, b, and c for organic chemicals, and criteria a, b, and d for heavy metals. If criteria a and/or b are below criterion d for a contaminant, its background value should be used as the cleanup objective. However, cleanup objectives developed using this approach must be, at a minimum, above the method detection limit (MDL) and it is preferable to have the soil cleanup objectives above the Contract Required Quantitation Limit (CRQL) as defined by NYSDEC. If the cleanup objective of a compound is "non-detectable", it should mean that it is not detected at the MDL. Efforts should be made to obtain the best MDL detection possible when selecting a laboratory and analytical protocol.

The water/soil partitioning theory is used to determine soil cleanup objectives which would be protective of groundwater/drinking water quality for its best use. This theory is conservative in nature and assumes that contaminated soil and groundwater are in direct contact. This theory is based upon the ability of organic matter in soil to adsorb organic chemicals. The approach predicts the maximum amount of contamination that may remain in soil so that leachate from the contaminated soil will not violate groundwater and/or drinking water

standards.

- (1) Class A are proved human carcinogens
- (2) Class B are probable human carcinogens
- (3) Class C are possible human carcinogens

This approach is not used for heavy metals, which do not partition appreciably into soil organic matter. For heavy metals, eastern USA or New York State soil background values may be used as soil cleanup objectives. A list of values that have been tabulated is attached.

Soil background data near the site, if available, is preferable and should be used as the cleanup objective for such metals. Background samples should be free from the influences of this site and any other source of contaminants. Ideal background samples may be obtained from uncontaminated upgradient and upwind locations.

### 3. DETERMINATION OF SOIL CLEANUP GOALS FOR ORGANICS IN SOIL FOR PROTECTION OF WATER QUALITY

Protection of water quality from contaminated soil is a two-part problem. The first is predicting the amount of contamination that will leave the contaminated media as leachate. The second part of the problem is to determine how much of that contamination will actually contribute to a violation of groundwater standards upon reaching and dispersing into groundwater. Some of the contamination which initially leaches out of soil will be absorbed by other soil before it reaches groundwater. Some portion will be reduced through natural attenuation or other mechanism.

#### PART A: PARTITION THEORY MODEL

There are many test and theoretical models which are used to predict leachate quality given a known value of soil contamination. The Water-Soil Equilibrium Partition Theory is used as a basis to determine soil standard or contamination limit for protection of water quality by most of the models currently in use. It is based on the ability of organic carbon in soil to adsorb contamination. Using a water quality value which may not be exceeded in leachate and the partition coefficient method, the equilibrium concentration ( $C_s$ ) will be expressed in the same units as the water standards. The following expression is used:

$$\text{Allowable Soil Concentration } C_s = f \times K_{oc} \times C_w \dots (1)$$

Where:  $f$  = fraction of organic carbon of the natural soil medium.

Koc = partition coefficient between water and soil media. Koc can be estimated by the following equation:

$$\log Koc = 3.64 - 0.55 \log S$$

S = water solubility in ppm

Cw = appropriate water quality value from TOGS 1.1.1

Most Koc and S values are listed in the Exhibit A-1 of the USEPA Superfund Public Health Evaluation Manual (EPA/540/1-86/060). The Koc values listed in this manual should be used for the purpose. If the Koc value for a contaminant is not listed, it should be estimated using the above mentioned equation.

#### PART-B: PROCEDURE FOR DETERMINATION OF SOIL CLEANUP OBJECTIVES

When the contaminated soil is in the unsaturated zone above the water table, many mechanisms are at work that prevent all of the contamination that would leave the contaminated soil from impacting groundwater. These mechanisms occur during transport and may work simultaneously. They include the following: (1) volatility, (2) sorption and desorption, (3) leaching and diffusion, (4) transformation and degradation, and (5) change in concentration of contaminants after reaching and/or mixing with the groundwater surface. To account for these mechanisms, a correction factor of 100 is used to establish soil cleanup objectives. This value of 100 for the correction is consistent with the logic used by EPA in its Dilution Attenuation Factor (DAF) approach for EP Toxicity and TCLP. (Federal Register/Vol. 55, No. 61, March 29, 1990/Pages 11826-27). Soil cleanup objectives are calculated by multiplying the allowable soil concentration by the correction factor. If the contaminated soil is very close (<3' - 5') to the groundwater table or in the groundwater, extreme caution should be exercised when using the correction factor of 100 (one hundred) as this may not give conservative cleanup objectives. For such situations the Technology Section should be consulted for site-specific cleanup objectives.

Soil cleanup objectives are limited to the following maximum values. These values are consistent with the approach promulgated by the States of Washington and Michigan.

- 1) Total VOCs  $\leq$  10 ppm.
- 2) Total Semi VOCs  $\leq$  500 ppm.
- 3) Individual Semi VOCs  $\leq$  50 ppm.
- 4) Total Pesticides  $\leq$  10 ppm.

One concern regarding the semi-volatile compounds is that some of these compounds are so insoluble that their Cs values are fairly large. Experience (Draft TOGS on Petroleum

Contaminated Soil Guidance) has shown that soil containing some of these insoluble substances at high concentrations can exhibit a distinct odor even though the substance will not leach from the soil. Hence any time a soil exhibits a discernible odor nuisance, it shall not be considered clean even if it has met the numerical criteria.

4. DETERMINATION OF FINAL CLEANUP LEVELS:

Recommended soil cleanup objectives should be utilized in the development of final cleanup levels through the Feasibility Study (FS) process. During the FS, various alternative remedial actions developed during the Remedial Investigation (RI) are initially screened and narrowed down to the list of potential alternative remedial actions that will be evaluated in detail. These alternative remedial actions are evaluated using the criteria discussed in TAGM 4030, Selection of Remedial Actions at Inactive Hazardous Waste Sites, revised May 15, 1990, and the preferred remedial action will be selected. After the detailed evaluation of the preferred remedial action, the final cleanup levels which can be actually achieved using the preferred remedial action must be established. Remedy selection, which will include final cleanup levels, is the subject of TAGM 4030.

Recommended soil cleanup objectives that have been calculated by the Technology Section are presented in Appendix A. These objectives are based on a soil organic carbon content of 1% (0.01) and should be adjusted for the actual organic carbon content if it is known. For determining soil organic carbon content, use attached USEPA method (Appendix B). Please contact the Technology Section, Bureau of Program Management for soil cleanup objectives not included in Appendix A.

Attachments

|                |                                      |
|----------------|--------------------------------------|
| cc: T. Jorling | J. Davis                             |
| J. Lacey       | J. Kelleher                          |
| M. Gerstman    | J. Colquhoun                         |
| A. DeBarbieri  | D. Persson                           |
| E. Sullivan    | A. Carlson                           |
| T. Donovan     | M. Birmingham                        |
| C. Sullivan    | D. Johnson                           |
| J. Eckl        | B. Hogan                             |
| R. Davies      | Regional Directors                   |
| R. Dana        | Regional Engineers                   |
| C. Goddard     | Regional Solid and Haz. Waste Engrs. |
| E. McCandless  | Regional Citizen Participation Spec. |
| P. Counterman  |                                      |



APPENDIX A  
TABLE 1  
Recommended soil cleanup objectives (mg/kg or ppm)  
Volatile Organic Contaminants

| Contaminant  | Partition coefficient<br>Koc | Groundwater Standards/<br>Criteria Cw<br>ug/l or ppb. | a                                     | b  | USEPA Health Based<br>(ppm) |                       | CRQL<br>(ppb) | ***<br>Rec. soil<br>Cleanup Obj<br>(ppm) |
|--|------------------------------|---|---------------------------------------|--|-----------------------------|-----------------------|---------------|--|
|  |                              |   | Allowable<br>Soil conc.<br>ppm.<br>Cs | Soil Cleanup<br>objectives to<br>Protect GW<br>Quality (ppm) | Carcinogens                 | Systemic<br>Toxicants |               |  |
| Acetone  | 2.2                          | 50  | 0.0011                                | 0.11   | N/A                         | 8,000                 | 10            | 0.                                       |
| Benzene  | 83                           | 0.7   | 0.0006                                | 0.06   | 24                          | N/A                   | 5             | 0.                                       |
| Benzoic Acid   | 54*                          | 50  | 0.027                                 | 2.7  | N/A                         | 300,000               | 5             | 2.                                       |
| 2-Butanone   | 4.5*                         | 50  | 0.003                                 | 0.3  | N/A                         | 4,000                 | 10            | 0.                                       |
| Carbon Disulfide   | 54*                          | 50  | 0.027                                 | 2.7  | N/A                         | 8,000                 | 5             | 2.                                       |
| Carbon Tetrachloride                                     | 110*                         | 5   | 0.006                                 | 0.6  | 5.4                         | 60                    | 5             | 0.                                       |
| Chlorobenzene  | 330                          | 5   | 0.017                                 | 1.7  | N/A                         | 2,000                 | 5             | 1.                                       |
| Chloroethane   | 37*                          | 50  | 0.019                                 | 1.9  | N/A                         | N/A                   | 10            | 1.                                       |
| Chloroform   | 31                           | 7   | 0.003                                 | 0.30   | 114                         | 800                   | 5             | 0.                                       |
| Bromochloroethane  | N/A                          | 50  | N/A                                   | N/A  | N/A                         | N/A                   | 5             | N/A                                      |
| 1,2-Dichlorobenzene                                      | 1,700                        | 4.7   | 0.079                                 | 7.9  | N/A                         | N/A                   | 330           | 7.                                       |
| 1,3-Dichlorobenzene                                      | 310 *                        | 5   | 0.0155                                | 1.55   | N/A                         | N/A                   | 330           | 1.                                       |
| 1,4-Dichlorobenzene                                      | 1,700                        | 5   | 0.085                                 | 8.5  | N/A                         | N/A                   | 330           | 8.                                       |
| 1,1-Dichloroethane                                       | 30                           | 5   | 0.002                                 | 0.2  | N/A                         | N/A                   | 5             | 0.                                       |
| 1,2-Dichloroethane                                       | 14                           | 5   | 0.001                                 | 0.1  | 7.7                         | N/A                   | 5             | 0.                                       |
| trans-Dichloroethane                                     | 65                           | 5   | 0.004                                 | 0.4  | 12                          | 700                   | 5             | 0.                                       |
| cis-Dichloroethane(trans)                                | 59                           | 5   | 0.003                                 | 0.3  | N/A                         | 2,000                 | 5             | 0.                                       |
| β-dichloropropene  | 51                           | 5   | 0.003                                 | 0.3  | N/A                         | N/A                   | 5             | 0.                                       |
| Ethylbenzene   | 1,100                        | 5   | 0.055                                 | 5.5  | N/A                         | 8,000                 | 5             | 5.                                       |
| 1,1,1,2,2,2-Freon(1,1,2 Trichloro-1,2,2 Trifluoroethane) | 1,230*                       | 5   | 0.060                                 | 6.0  | N/A                         | 200,000               | 5             | 6.                                       |
| Methylene chloride                                       | 21                           | 5   | 0.001                                 | 0.1  | 93                          | 5,000                 | 5             | 0.                                       |
| Methyl-2-Pentanone                                       | 19*                          | 50  | 0.01                                  | 1.0  | N/A                         | N/A                   | 10            | 1.                                       |
| 1,1,1-Trichloroethane                                    | 277                          | 5   | 0.014                                 | 1.4  | 14                          | 800                   | 5             | 1.                                       |
| 1,1,1-Trichloroethane                                    | 152                          | 5   | 0.0076                                | 0.76   | N/A                         | 7,000                 | 5             | 0.                                       |
| 1,1,2,2-Tetrachloroethane                                | 118                          | 5   | 0.006                                 | 0.6  | 35                          | N/A                   | 5             | 0.                                       |
| 2,3-trichloropropane                                     | 68                           | 5   | 0.0034                                | 0.34   | N/A                         | 80                    | 5             | 0.                                       |
| 2,4-Trichlorobenzene                                     | 670 *                        | 5   | 0.034                                 | 3.4  | N/A                         | N/A                   | 330           | 3.                                       |
| Toluene  | 300                          | 5   | 0.015                                 | 1.5  | N/A                         | 20,000                | 5             | 1.                                       |
| Trichloroethene  | 126                          | 5   | 0.007                                 | 0.70   | 64                          | N/A                   | 5             | 0.                                       |
| Vinyl chloride   | 57                           | 2   | 0.0012                                | 0.12   | N/A                         | N/A                   | 10            | 0.                                       |
| Xylenes  | 240                          | 5   | 0.012                                 | 1.2  | N/A                         | 200,000               | —             | 1.                                       |

a. Allowable Soil Concentration  $C_s = f \times C_w \times K_{oc}$

b. Soil cleanup objective =  $C_s \times \text{Correction Factor (CF)}$

N/A is not available

\* Partition coefficient is calculated by using the following equation:

$\log K_{oc} = -0.55 \log S + 3.64$ , where S is solubility in water in ppm.

All other Koc values are experimental values.

Correction Factor (CF) of 100 is used as per TAGM #4046

As per TAGM #4046, Total VOCs < 10 ppm.

Note: Soil cleanup objectives are developed for soil organic carbon content (f) of 1% , and should be adjusted for the actual soil organic carbon content if it is known.

APPENDIX A (cont.)

TABLE 2

Recommended Soil Cleanup Objectives (mg/kg or ppm)  
Semi-Volatile Organic Contaminants

| Contaminant                | Partition coefficient Koc | Groundwater Standards/ Criteria Cw ug/l or ppb. | a                            | b   | ** | USEPA Health Based (ppm) |                    | CRQL (ppc) | Rec. soil Cleanup Obj. (ppm) |
|----------------------------|---------------------------|---|------------------------------|---|----|--------------------------|--------------------|------------|------------------------------|
|                            |                           |   | Allowable Soil conc. ppm. Cs | Soil Cleanup objectives to Protect GW Quality (ppm) |    | Carcinogens              | Systemic Toxicants |            |                              |
| Acenaphthene               | 4,600                     | 20  | 0.9                          | 90.0  |    | N/A                      | 5,000              | 330        | 50.0                         |
| Acenaphthylene             | 2,056*                    | 20  | 0.41                         | 41.0  |    | N/A                      | N/A                | 330        | 5.0                          |
| Aniline                    | 13.8                      | 5   | 0.001                        | 0.1   |    | 123                      | N/A                | 330        | 0.1                          |
| Anthracene                 | 14,000                    | 50  | 7.00                         | 700.0   |    | N/A                      | 20,000             | 330        | 50.0*                        |
| Benzo(a)anthracene         | 1,380,000                 | 0.002   | 0.03                         | 3.0   |    | 0.224                    | N/A                | 330        | 0.224 or                     |
| Benzo(a)pyrene             | 5,500,000                 | 0.002(MD)                                       | 0.110                        | 11.0  |    | 0.0609                   | N/A                | 330        | 0.061 or                     |
| Benzo(b)fluoranthene       | 550,000                   | 0.002   | 0.011                        | 1.1   |    | N/A                      | N/A                | 330        | 1.1                          |
| Benzo(g,h,i)perylene       | 1,600,000                 | 5   | 8.0                          | 800   |    | N/A                      | N/A                | 330        | 50.0*                        |
| Benzo(k)fluoranthene       | 550,000                   | 0.002   | 0.011                        | 1.1   |    | N/A                      | N/A                | 330        | 1.1                          |
| bis(2-ethylhexyl)phthalate | 8,706*                    | 50  | 4.35                         | 435.0   |    | 50                       | 2,000              | 330        | 50.0                         |
| Butylbenzylphthalate       | 2,430                     | 50  | 1.215                        | 122.0   |    | N/A                      | 20,000             | 330        | 50.0*                        |
| Chrysene                   | 200,000                   | 0.002   | 0.004                        | 0.4   |    | N/A                      | N/A                | 330        | 0.4                          |
| 4-Chloroaniline            | 43 ****                   | 5   | 0.0022                       | 0.22  |    | 200                      | 300                | 330        | 0.220 or                     |
| 4-Chloro-3-methylphenol    | 47                        | 5   | 0.0024                       | 0.24  |    | N/A                      | N/A                | 330        | 0.240 or                     |
| 2-Chlorophenol             | 15*                       | 50  | 0.008                        | 0.8   |    | N/A                      | 400                | 330        | 0.8                          |
| Dibenzofuran               | 1,230*                    | 5   | 0.062                        | 6.2   |    | N/A                      | N/A                | 330        |                              |
| Benzo(a,h)anthracene       | 33,000,000                | 50  | 1,650                        | 165,000   |    | 0.0143                   | N/A                | 330        | 0.014 or                     |
| 3'-Dichlorobenzidine       | N/A                       | N/A   | N/A                          | N/A   |    | N/A                      | N/A                | N/A        | N/A                          |
| 2,4-Dichlorophenol         | 380                       | 1   | 0.004                        | 0.4   |    | N/A                      | 200                | 330        | 0.4                          |
| 2,4-Dinitrophenol          | 38                        | 5   | 0.002                        | 0.2   |    | N/A                      | 200                | 1,600      | 0.200 or                     |
| 2,6-Dinitrotoluene         | 198*                      | 5   | 0.01                         | 1.0   |    | 1.03                     | N/A                | 330        | 1.0                          |
| Diethylphthalate           | 142                       | 50  | 0.071                        | 7.1   |    | N/A                      | 60,000             | 330        | 7.1                          |
| Dimethylphthalate          | 40                        | 50  | 0.020                        | 2.0   |    | N/A                      | 80,000             | 330        | 2.0                          |
| Di-n-butyl phthalate       | 162*                      | 50  | 0.081                        | 8.1   |    | N/A                      | 8,000              | 330        | 8.1                          |
| Di-n-octyl phthalate       | 2,346*                    | 50  | 1.2                          | 120.0   |    | N/A                      | 2,000              | 330        | 50.0*                        |
| Fluoranthene               | 36,000                    | 50  | 19                           | 1900.0  |    | N/A                      | 3,000              | 330        | 50.0*                        |
| Fluorene                   | 7,300                     | 50  | 3.5                          | 350.0   |    | N/A                      | 3,000              | 330        | 50.0                         |
| Hexachlorobenzene          | 3,900                     | 0.35  | 0.014                        | 1.4   |    | 0.41                     | 60                 | 330        | 0.4                          |
| Indeno(1,2,3-cd)pyrene     | 1,600,000                 | 0.002   | 0.032                        | 3.2   |    | N/A                      | N/A                | 330        | 3.2                          |
| Isophorone                 | 88.31*                    | 50  | 0.044                        | 4.40  |    | 1,707                    | 20,000             | 330        | 4.4                          |
| 2-methylnaphthalene        | 727*                      | 50  | 0.354                        | 36.4  |    | N/A                      | N/A                | 330        | 36.4                         |
| 2-Methylphenol             | 15                        | 5   | 0.001                        | 0.1   |    | N/A                      | N/A                | 330        | 0.100 or                     |
| 4-Methylphenol             | 17                        | 50  | 0.009                        | 0.9   |    | N/A                      | 4,000              | 330        | 0.9                          |
| Naphthalene                | 1,300                     | 10  | 0.130                        | 13.0  |    | N/A                      | 300                | 330        | 13.0                         |
| Nitrobenzene               | 36                        | 5   | 0.002                        | 0.2   |    | N/A                      | 40                 | 330        | 0.200 or                     |
| 2-Nitroaniline             | 86                        | 5   | 0.0043                       | 0.43  |    | N/A                      | N/A                | 1,600      | 0.430 or                     |
| 2-Nitrophenol              | 65                        | 5   | 0.0033                       | 0.33  |    | N/A                      | N/A                | 330        | 0.330 or                     |
| 4-Nitrophenol              | 21                        | 5   | 0.001                        | 0.1   |    | N/A                      | N/A                | 1,600      | 0.100 or                     |
| 3-Nitroaniline             | 93                        | 5   | 0.005                        | 0.5   |    | N/A                      | N/A                | 1,600      | 0.500 or                     |
| Pentachlorophenol          | 1,022                     | 1   | 0.01                         | 1.0   |    | N/A                      | 2,000              | 1,600      | 1.0 or MD                    |
| Phenanthrene               | 4,365*                    | 50  | 2.20                         | 220.0   |    | N/A                      | N/A                | 330        | 50.0                         |
| Phenol                     | 27                        | 1   | 0.0003                       | 0.03  |    | N/A                      | 50,000             | 330        | 0.03 or                      |
| Pyrene                     | 13,295*                   | 50  | 6.65                         | 665.0   |    | N/A                      | 2,000              | 330        | 50.0*                        |
| 1,4,5-Trichlorophenol      | 89*                       | 1   | 0.001                        | 0.1   |    | N/A                      | 8,000              | 330        |                              |

APPENDIX A  
NEW YORK STATE DEPARTMENT OF ENVIRONMENTAL CONSERVATION  
TECHNICAL & ADMINISTRATIVE GUIDANCE MEMORADUM  
LACKAWANNA FOUNDRY SITE  
LACKAWANNA, NEW YORK  
MAY 1999

- a. Allowable Soil Concentration  $C_s = f \times C_w \times K_{oc}$
- b. Soil cleanup objective =  $C_s \times$  Correction Factor (CF)

N/A is not available

MDL is Method Detection Limit

Partition coefficient is calculated by using the following equation:

$\log K_{oc} = -0.55 \log S + 3.64$ , where S is solubility in water in ppm. Other  $K_{oc}$  values are experimental values.

\*\* Correction Factor (CF) of 100 is used as per TAGM #4046

\*\*\* As per TAGM #4046, Total VOCs < 10 ppm., Total Semi-VOCs < 500 ppm. and Individual Semi-VOCs < 50 ppm.

\*\*  $K_{oc}$  is derived from the correlation  $K_{oc} = 0.63 K_{ow}$  ( Determining Soil Response Action Levels..... EPA/540/2-89/057 ).  $K_{ow}$  is obtained from the USEPA computer database 'MAIN'.

te: Soil cleanup objectives are developed for soil organic carbon content (f) of 1%, and should be adjusted for the actual soil organic carbon content if it is known.

APPENDIX A (cont.)

TABLE 3

Recommended soil cleanup objectives (mg/kg or ppm)  
Organic Pesticides / Herbicides and PCBs

| Contaminant                         | Partition coefficient Koc | Groundwater Standards/ Criteria Cw ug/l or pob. | a                            | b   | **          | USEPA Health Based (ppm) |                         | CRCL (pob) | ***                       |
|-------------------------------------|---------------------------|---|------------------------------|---|-------------|--------------------------|-------------------------|------------|---------------------------|
|                                     |                           |   | Allowable Soil conc. ppm. Cs | Soil Cleanup objectives to Protect GW Quality (ppm) | Carcinogens | Systemic Toxicants       | Req. soil Cleanup (ppm) |            |                           |
| Aldrin                              | 96,000                    | ND(<0.01)                                       | 0.005                        | 0.5   |             | 0.041                    | 2                       | 8          | 0.04                      |
| alpha - BHC                         | 3,800                     | ND(<0.05)                                       | 0.002                        | 0.2   |             | 0.111                    | N/A                     | 8          | 0.11                      |
| Beta - BHC                          | 3,800                     | ND(<0.05)                                       | 0.002                        | 0.2   |             | 3.89                     | N/A                     | 8          | 0.2                       |
| delta - BHC                         | 6,600                     | ND(<0.05)                                       | 0.003                        | 0.3   |             | N/A                      | N/A                     | 8          | 0.3                       |
| Chlordane                           | 21,305*                   | 0.1   | 0.02                         | 2.0   |             | 0.54                     | 50                      | 80         | 0.54                      |
| 2,4-D                               | 104*                      | 4.4   | 0.005                        | 0.5   |             | N/A                      | 800                     | 800        | 0.5                       |
| 4,4'-DDD                            | 770,000*                  | ND(<0.01)                                       | 0.077                        | 7.7   |             | 2.9                      | N/A                     | 16         | 2.9                       |
| 4,4'-DDE                            | 440,000*                  | ND(<0.01)                                       | 0.0440                       | 4.4   |             | 2.1                      | N/A                     | 16         | 2.1                       |
| 4,4'-DDT                            | 243,000*                  | ND(<0.01)                                       | 0.025                        | 2.5   |             | 2.1                      | 40                      | 16         | 2.1                       |
| Dibenzo-P-dioxins(PCDD)             |                           |   |                              |   |             |                          |                         |            |                           |
| 2,3,7,8 TCDD                        | 1709800                   | 0.000035  | 0.0006                       | 0.06  |             | N/A                      | N/A                     | N/A        | N/A                       |
| Dieldrin                            | 10,700*                   | ND(<0.01)                                       | 0.0010                       | 0.1   |             | 0.044                    | 4                       | 16         | 0.044                     |
| Endosulfan I                        | 8,168*                    | 0.1   | 0.009                        | 0.9   |             | N/A                      | N/A                     | 16         | 0.9                       |
| Endosulfan II                       | 8,031*                    | 0.1   | 0.009                        | 0.9   |             | N/A                      | N/A                     | 16         | 0.9                       |
| Endosulfan Sulfate                  | 10,032*                   | 0.1   | 0.01                         | 1.0   |             | N/A                      | N/A                     | 16         | 1.0                       |
| Endrin                              | 9,157*                    | ND(<0.01)                                       | 0.001                        | 0.1   |             | N/A                      | 20                      | 8          | 0.1                       |
| Endrin ketone                       | N/A                       | N/A   | N/A                          | N/A   |             | N/A                      | N/A                     | N/A        | N/A                       |
| gamma - BHC (Lindane)               | 1,080                     | ND(<0.05)                                       | 0.0006                       | 0.06  |             | 5.4                      | 20                      | 8          | 0.06                      |
| gamma - chlordane                   | 140,000                   | 0.1   | 0.14                         | 14.0  |             | 0.54                     | 5                       | 80         | 0.54                      |
| heptachlor                          | 12,000                    | ND(<0.01)                                       | 0.0010                       | 0.1   |             | 0.16                     | 40                      | 8          | 0.10                      |
| heptachlor epoxide                  | 220                       | ND(<0.01)                                       | 0.0002                       | 0.02  |             | 0.077                    | 0.8                     | 8          | 0.02                      |
| methoxychlor                        | 25,637                    | 35.0  | 9.0                          | 900   |             | N/A                      | 400                     | 80         | 9.0                       |
| metolene                            | N/A                       | N/A   | N/A                          | N/A   |             | N/A                      | N/A                     | N/A        | N/A                       |
| parathion                           | 760                       | 1.5   | 0.012                        | 1.2   |             | N/A                      | 500                     | 8          | 1.2                       |
| PCBs                                | 17,510*                   | 0.1   | 0.1                          | 10.0  |             | 1.0                      | N/A                     | 160        | 1.0(Surface<br>10(sub-sur |
| Polychlorinated dibenzofurans(PCDF) |                           |   |                              |   |             |                          |                         |            |                           |
|                                     | N/A                       | N/A   | N/A                          | N/A   |             | N/A                      | N/A                     | N/A        | N/A                       |
| dicofol                             | 2,600                     | 0.26  | 0.007                        | 0.7   |             | N/A                      | 600                     | 330        | 0.7                       |
| 2,4,5-T                             | 53                        | 35  | 0.019                        | 1.9   |             | N/A                      | 200                     | 330        | 1.9                       |

a. Allowable Soil Concentration Cs = f x Cw x Koc

b. Soil cleanup objective = Cs x Correction Factor (CF)

N/A is not available

Partition coefficient is calculated by using the following equation:

$\log Koc = -0.55 \log S + 3.64$ , where S is solubility in water in ppm.

All other Koc values are experimental values.

Correction Factor (CF) of 100 is used as per TAGM #4046

\* As per TAGM #4046, Total Pesticides < 10 ppm.

Note: Soil cleanup objectives are developed for soil organic carbon content (f) of 1% (5% for PCBs as per PCB guidance document), and should be adjusted for the actual soil organic carbon content if it is known.

## APPENDIX A

Rev. 12/93

TABLE 4

Recommended Soil Cleanup Objectives (mg/kg or ppm) for Heavy Metals

| Contaminants | Protect Water Quality ppm | Eastern USA Background ppm | CRDL mg/kg or ppm | Rec.soil Cleanup Objct. (ppm) |
|--------------|---------------------------|----------------------------|-------------------|-------------------------------|
| Aluminum     | N/A                       | 33,000                     | 2.0               | SB                            |
| Antimony     | N/A                       | N/A                        | 0.6               | SB                            |
| Arsenic      | N/A                       | 3-12 **                    | 0.1               | 7.5 or SB                     |
| Bismuth      | N/A                       | 15-600                     | 2.0               | 300 or SB                     |
| Beryllium    | N/A                       | 0-1.75                     | 0.05              | 0.15(HEAST) or SB             |
| Cadmium      | N/A                       | 0.1-1                      | 0.05              | 1 or SB                       |
| Chromium     | N/A                       | 130 - 35,000 **            | 50.0              | SB                            |
| Cobalt       | N/A                       | 1.5-40 **                  | 0.1               | 10 or SB                      |
| Cobalt       | N/A                       | 2.5-60 **                  | 0.5               | 30 or SB                      |
| Copper       | N/A                       | 1-50                       | 0.25              | 25 or SB                      |
| Cyanide      | N/A                       | N/A                        | 0.1               | ***                           |
| Iron         | N/A                       | 2,000 - 550,000            | 1.0               | 2,000 or SB                   |
| Lead         | N/A                       | ****                       | 0.03              | SB****                        |
| Magnesium    | N/A                       | 100 - 5,000                | 50.0              | SB                            |
| Manganese    | N/A                       | 50 - 5,000                 | 0.15              | SB                            |
| Mercury      | N/A                       | 0.001-0.2                  | 0.002             | 0.1                           |
| Nickel       | N/A                       | 0.5-25                     | 0.4               | 13 or SB                      |
| Potassium    | N/A                       | 8,500 - 43,000 **          | 50.0              | SB                            |
| Selenium     | N/A                       | 0.1-3.9                    | 0.05              | 2 or SB                       |
| Silver       | N/A                       | N/A                        | 0.1               | SB                            |
| Sodium       | N/A                       | 6,000 - 8,000              | 50.0              | SB                            |
| Thallium     | N/A                       | N/A                        | 0.1               | SB                            |
| Vanadium     | N/A                       | 1-300                      | 0.5               | 150 or SB                     |
| Zinc         | N/A                       | 9-50                       | 0.2               | 20 or SB                      |

Note: Some forms of metal salts such as Aluminum Phosphide, Calcium Cyanide, Potassium Cyanide, Copper cyanide, Silver cyanide, Sodium cyanide, Zinc phosphide, Thallium salts, Vanadium pentoxide, and Chromium (VI) compounds are more toxic in nature. Please refer to the USEPA HEASTs database to find cleanup objectives if such metal salts are present in soil.

SB is site background

N/A is not available

CRDL is contract required detection limit which is approx. 10 times the CRDL for water.

New York State background

\*\* Some forms of Cyanide are complex and very stable while other forms are pH dependent and hence are very unstable. Site-specific form(s) of Cyanide should be taken into consideration when establishing soil cleanup objective.

\*\*\* Background levels for lead vary widely. Average levels in undeveloped, rural areas may range from 4-61 ppm. Average background levels in metropolitan or suburban areas or near highways are much higher and typically range from 200-500 ppm.

\*Recommended soil cleanup objectives are average background concentrations as reported in a 1984 survey of reference material by E. Carol McGovern, NYSDEC.

## TOTAL ORGANIC CARBON (TOC)

### USE AND LIMITATIONS

Total organic carbon is a measure of the total amount of nonvolatile, volatile, partially volatile, and particulate organic compounds in a sample. Total organic carbon is independent of the oxidation state of the organic compounds and is not a measure of the organically bound and inorganic elements that can contribute to the biochemical and chemical oxygen demand tests.

Because inorganic carbon (e.g., carbonates, bicarbonates, free CO<sub>2</sub>) will interfere with total organic carbon determinations, samples should be treated to remove inorganic carbon before being analyzed.

### FIELD PROCEDURES

#### Collection

Samples can be collected in glass or plastic containers. A minimum sample size of 25 g is recommended. If unrepresentative material is to be removed from the sample, it should be removed in the field under the supervision of the chief scientist and noted on the field log sheet.

#### Processing

Samples should be stored frozen and can be held for up to 6 mo under that condition. Excessive temperatures should not be used to thaw samples.

### LABORATORY PROCEDURES

#### Analytical Procedures

##### • Equipment

- Induction furnace  
e.g., Leco WR-12, Dohrmann DC-50, Coleman CH analyzer,  
Perkin Elmer 240 elemental analyzer, Carlo-Erba 1106
- Analytical balance  
0.1 mg accuracy
- Desiccator
- Combustion boats
- 10 percent hydrochloric acid (HCl)
- Cupric oxide fines (or equivalent material)
- Benzoic acid or other carbon source as a standard.

Conventional Sediment Variables  
Total Organic Carbon (TOC)  
March 1986

• Equipment preparation

- Clean combustion boats by placing them in the induction furnace at 950° C. After being cleaned, combustion boats should not be touched with bare hands.
- Cool boats to room temperature in a desiccator.
- Weigh each boat to the nearest 0.1 mg.

• Sample preparation

- Allow frozen samples to warm to room temperature.
- Homogenize each sample mechanically, incorporating any overlying water.
- Transfer a representative aliquot (5-10 g) to a clean container.

• Analytical procedures

- Dry samples to constant weight at 70 ± 2° C. The drying temperature is relatively low to minimize loss of volatile organic compounds.
- Cool dried samples to room temperature in a desiccator.
- Grind sample using a mortar and pestle to break up aggregates.
- Transfer a representative aliquot (0.2-0.5 g) to a clean, preweighed combustion boat.
- Determine sample weight to the nearest 0.1 mg.
- Add several drops of HCl to the dried sample to remove carbonates. Wait until the effervescing is completed and add more acid. Continue this process until the incremental addition of acid causes no further effervescence. Do not add too much acid at one time as this may cause loss of sample due to frothing. Exposure of small samples (i.e., 1-10 mg) having less than 50 percent carbonate to an HCl atmosphere for 24-48 h has been shown to be an effective means of removing carbonates (Hedges and Stern 1984). If this method is used for sample sizes greater than 10 mg, its effectiveness should be demonstrated by the user.
- Dry the HCl-treated sample to constant weight at 70 ± 2° C.
- Cool to room temperature in a desiccator.
- Add previously ashed cupric oxide fines or equivalent material (e.g., alumina oxide) to the sample in the combustion boat.
- Combust the sample in an induction furnace at a minimum temperature of 950 ± 10° C.

• Calculations

- If an ascarite-filled tube is used to capture CO<sub>2</sub>, the carbon content of the sample can be calculated as follows:

$$\text{Percent carbon} = \frac{A(0.2729)(100)}{g}$$



Conventional Sediment Variables  
Total Organic Carbon (TOC)  
March 1986

Where:

- A = the weight (g) of CO<sub>2</sub> determined by weighing the ascarite tube before and after combustion
- B = dry weight (g) of the unacidified sample in the combustion boat
- 0.2729 = the ratio of the molecular weight of carbon to the molecular weight of carbon dioxide

A silica gel trap should be placed before the ascarite tube to catch any moisture driven off during sample combustion. Additional silica gel should be placed at the exit end of the ascarite tube to trap any water that might be formed by reaction of the trapped CO<sub>2</sub> with the NaOH in the ascarite.

- If an elemental analyzer is used, the amount of CO<sub>2</sub> will be measured by a thermal conductivity detector. The instrument should be calibrated daily using an empty boat blank as the zero point and at least two standards. Standards should bracket the expected range of carbon concentrations in the samples.

#### QA/QC Procedures

It is critical that each sample be thoroughly homogenized in the laboratory before a subsample is taken for analysis. Laboratory homogenization should be conducted even if samples were homogenized in the field.

Dried samples should be cooled in a desiccator and held there until they are weighed. If a desiccator is not used, the sediment will accumulate ambient moisture and the sample weight will be overestimated. A color-indicating desiccant is recommended so that spent desiccant can be detected easily. Also, the seal on the desiccator should be checked periodically and, if necessary, the ground glass rims should be greased or the "O" rings should be replaced.

It is recommended that triplicate analyses be conducted on one of every 20 samples, or on one sample per batch if less than 20 samples are analyzed. A method blank should be analyzed at the same frequency as the triplicate analyses. The analytical balance should be inspected daily and calibrated at least once per week. The carbon analyzer should be calibrated daily with freshly prepared standards. A standard reference material should be analyzed at least once for each major survey.

#### DATA REPORTING REQUIREMENTS

Total organic carbon should be reported as a percentage of the dry weight of the unacidified sample to the nearest 0.1 unit. The laboratory should report the results of all samples (including QA replicates, method

# Appendix B

APPENDIX B  
SOIL SAMPLING FIELD DATA SHEETS  
LACKAWANNA FOUNDRY SITE  
LACKAWANNA, NEW YORK  
MAY 1999

FIELD DATA SHEET

16650

REAC, EDISON, NJ
(908) 321-4200
EPA CONTRACT 68-C4-0022

Date: 4/14/99
Samplers: Fanni Magdon Metz
Site Name: Lackenbach Foundry
Sample Location: West Fence line
Chain of Custody No.:
REAC Task Leader: Fanni
EPA WAM:
Work Assignment No.:

Table with 5 columns: SITE DESCRIPTION, SOIL TYPE, SURFACE WATER, STREAM, BOTTOM. Includes categories like landfill, industrial, commercial, residential, hedgerows, etc.

Table with 5 columns: SAMPLE TYPE, DEVICE, SAMPLE INFORMATION, WEATHER PARAMETERS. Includes categories like surface water, groundwater, sediment, soil, etc.

ANALYSES TO BE PERFORMED

SAMPLE PREPARATION

- ORGANICS
A. halogenated & aromatic volatiles
B. volatiles
C. trihalomethanes
D. pesticides/PCB
E. PCB
F. base neutral/acid extractables
G. pesticides, drinking water
H. herbicides, drinking water
I. other

- OTHER ANALYSES
A. total cyanide
B. total phenol
C. petroleum hydrocarbons
D. pH
E. alkalinity
F. hardness
G. total dissolved solids
H. total suspended solids
I. sulfate
J. TOC
K. grain size
L. percent moisture
M. other

- CONTAINER
glass jar
plastic jar
acetate core
plastic bag
plastic bucket
other

- PRESERVATIVES
HNO3
NaOH
Zn Acetate
HCl
Na2SO4
other

- INORGANICS
A. metals, priority pollutant
B. metals, TAL
C. metals scan (ICP)
D. metals, other

- STORAGE
wet ice
dry ice
ambient

- RCRA
A. TCLP
B. ignitability
C. corrosivity pH
D. reactivity
E. other

COMMENTS

**FIELD DATA SHEET**

16651

**REAC, EDISON, NJ  
(908) 321-4200  
EPA CONTRACT 68-C4-0022**

Date: 4/14/96      Samplers: Fama, Melo, Maza      Chain of Custody No.: \_\_\_\_\_  
 Site Name: Lackawanna      REAC Task Leader: Fama  
 Time: 1100      Sample Location: #2 surface middle well      EPA WAM: \_\_\_\_\_  
 Work Assignment No.: \_\_\_\_\_

| SITE DESCRIPTION  |            |                   | SOIL TYPE   |              | SURFACE WATER |       | STREAM   |            | BOTTOM |             |
|-------------------|------------|-------------------|-------------|--------------|---------------|-------|----------|------------|--------|-------------|
| landfill          | old field  | upland palustrine | rock        | clay         | color         | _____ | width    | _____      | rock   | silt        |
| <u>industrial</u> | wooded     | lowland riverine  | gravel      | muck         | odor          | _____ | depth    | _____      | rubble | clay        |
| commercial        | famland    | lacustrine        | <u>sand</u> | loam         | flow          | _____ | velocity | _____ cm/s | gravel | organic     |
| residential       | gully      |                   | <u>silt</u> | peat         | direction     | _____ | pools    | _____ %    | shell  | other _____ |
| hedgerows         | floodplain |                   | color       | <u>BROWN</u> |               |       | riffles  | _____ %    | sand   |             |

| SAMPLE TYPE   |          | DEVICE   |       | SAMPLE INFORMATION |              | WEATHER PARAMETERS  |             |
|---------------|----------|----------|-------|--------------------|--------------|---------------------|-------------|
| surface water | effluent | kemmerer | ponar | color              | <u>BROWN</u> | pH                  | _____       |
| groundwater   | sludge   | trowel   | other | odor               | _____        | ORP                 | _____       |
| potable water | leachate | bucket   |       | temp               | <u>50°F</u>  | salinity            | _____       |
| sediment      | waste    | auger    |       | DO                 | _____        | sample depth        | _____       |
| <u>SOIL</u>   | other    | ekman    |       | cond               | _____        | tide stage          | _____       |
|               |          |          |       |                    |              | ambient temp        | <u>40°F</u> |
|               |          |          |       |                    |              | barometric pressure | _____       |
|               |          |          |       |                    |              | relative humidity   | <u>30%</u>  |
|               |          |          |       |                    |              | weather conditions  | _____       |

**ANALYSES TO BE PERFORMED**

**ORGANICS**

- A. halogenated & aromatic volatiles
- B. volatiles
- C. trihalomethanes
- D. pesticides/PCB
- E. PCB
- F. base neutral/acid extractables
- G. pesticides, drinking water
- H. herbicides, drinking water
- I. other \_\_\_\_\_

**INORGANICS**

- A. metals, priority pollutant
- B. metals, TAL
- C. metals scan (ICP)
- D. metals, other \_\_\_\_\_

**RCRA**

- A. TCLP
- B. ignitability
- C. corrosivity \_\_\_\_\_ pH \_\_\_\_\_
- D. reactivity
- E. other \_\_\_\_\_

**OTHER ANALYSES**

- A. total cyanide
- B. total phenol
- C. petroleum hydrocarbons
- D. pH
- E. alkalinity
- F. hardness
- G. total dissolved solids
- H. total suspended solids
- I. sulfate
- J. TOC
- K. grain size
- L. percent moisture
- M. other \_\_\_\_\_

**SAMPLE PREPARATION**

**CONTAINER**

- glass jar
- plastic jar
- acetate core
- plastic bag
- plastic bucket
- other \_\_\_\_\_

**PRESERVATIVES**

- HNO<sub>3</sub>
- NaOH
- Zn Acetate
- HCl
- Na<sub>2</sub>SO<sub>4</sub>
- other \_\_\_\_\_

**STORAGE**

- wet ice
- dry ice
- ambient

**COMMENTS**



**FIELD DATA SHEET**

16652

**REAC, EDISON, NJ  
(908) 321-4200  
EPA CONTRACT 68-C4-0022**

Date: 4/14/86      Samplers: Fama, M. et al, Magan      Chain of Custody No.: \_\_\_\_\_  
 Time: 1100      Site Name: Leakage from Foundation      REAC Task Leader: Fama  
 Sample Location: #20 sub surface      EPA WAM: Campanelli      Work Assignment No.: 3-439

| SITE DESCRIPTION  |            |                   | SOIL TYPE          |      | SURFACE WATER   |                     | STREAM |             | BOTTOM |  |
|-------------------|------------|-------------------|--------------------|------|-----------------|---------------------|--------|-------------|--------|--|
| landfill          | old field  | upland palustrine | rock               | clay | color _____     | width _____         | rock   | silt        |        |  |
| <u>industrial</u> | wooded     | lowland riverine  | gravel             | muck | odor _____      | depth _____         | rubble | clay        |        |  |
| commercial        | farmland   | lacustrine        | <u>sand</u>        | loam | flow _____      | velocity _____ cm/s | gravel | organic     |        |  |
| residential       | gully      |                   | silt               | peat | direction _____ | pools _____ %       | shell  | other _____ |        |  |
| hedgerows         | floodplain |                   | color <u>Brown</u> |      |                 | riffles _____ %     | sand   |             |        |  |

| SAMPLE TYPE   |             | DEVICE       |   | SAMPLE INFORMATION |                    | WEATHER PARAMETERS        |  |
|---------------|-------------|--------------|---|--------------------|--------------------|---------------------------|--|
| surface water | effluent    | kemmerer     | ponar  | color <u>Brown</u> | pH _____           | ambient temp <u>40°F</u>  |  |
| groundwater   | sludge      | trowel       | other  | odor _____         | ORP _____          | barometric pressure _____ |  |
| potable water | leachate    | bucket       |   | temp _____         | salinity _____     | relative humidity _____   |  |
| sediment      | waste       | <u>auger</u> |   | DO _____           | sample depth _____ | weather conditions _____  |  |
| <u>soil</u>   | other _____ | ekman        |   | cond _____         | tide stage _____   |                           |  |

**ANALYSES TO BE PERFORMED**

- ORGANICS**
- A. halogenated & aromatic volatiles
  - B. volatiles
  - C. trihalomethanes
  - D. pesticides/PCB
  - E. PCB
  - F. base neutral/acid extractables
  - G. pesticides, drinking water
  - H. herbicides, drinking water
  - I. other \_\_\_\_\_

- INORGANICS**
- A. metals, priority pollutant
  - B. metals, TAL
  - C. metals scan (ICP)
  - D. metals, other \_\_\_\_\_

- RCRA**
- A. TCLP
  - B. ignitability
  - C. corrosivity \_\_\_\_\_ pH \_\_\_\_\_
  - D. reactivity
  - E. other \_\_\_\_\_

**COMMENTS**

**SAMPLE PREPARATION**

- CONTAINER**
- glass jar
  - plastic jar      NaOH
  - acetate core
  - plastic bag
  - plastic bucket
  - other \_\_\_\_\_
- PRESERVATIVES**
- HNO<sub>3</sub>
  - Zn Acetate
  - HCl
  - Na<sub>2</sub>SO<sub>4</sub>
  - other \_\_\_\_\_

- STORAGE**
- Wet ice
  - dry ice
  - ambient

# FIELD DATA SHEET

16653

**REAC, EDISON, NJ  
(908) 321-4200  
EPA CONTRACT 68-C4-0022**

Date: 4/14/96 Samplers: Farr, Metc, Mize Chain of Custody No.: \_\_\_\_\_  
 Time: 1100 Site Name: Lackawanna Foundry REAC Task Leader: Farr  
 Sample Location: #3 West Fence EPA WAM: Compton  
 Work Assignment No.: 3-439

| SITE DESCRIPTION  |            |                   | SOIL TYPE        |       | SURFACE WATER |       | STREAM   |            | BOTTOM |         |
|-------------------|------------|-------------------|------------------|-------|---------------|-------|----------|------------|--------|---------|
| landfill          | old field  | upland palustrine | rock             | clay  | color         | _____ | width    | _____      | rock   | silt    |
| <u>industrial</u> | wooded     | lowland riverine  | gravel           | muck  | odor          | _____ | depth    | _____      | rubble | clay    |
| commercial        | farmland   | lacustrine        | <u>sand loam</u> | _____ | flow          | _____ | velocity | _____ cm/s | gravel | organic |
| residential       | gully      | _____             | silt             | peat  | direction     | _____ | pools    | _____ %    | shell  | other   |
| hedgerows         | floodplain | _____             | color            | _____ | _____         | _____ | riffles  | _____ %    | sand   | _____   |

| SAMPLE TYPE     |          | DEVICE   |                   | SAMPLE INFORMATION |              | WEATHER PARAMETERS  |             |
|-----------------|----------|----------|-------------------|--------------------|--------------|---------------------|-------------|
| surface water   | effluent | kemmerer | ponar             | color              | <u>Black</u> | pH                  | _____       |
| groundwater     | sludge   | trowel   | other <u>span</u> | odor               | _____        | ORP                 | _____       |
| potable water   | leachate | bucket   | _____             | temp               | <u>50°F</u>  | salinity            | _____       |
| <u>sediment</u> | waste    | auger    | _____             | DO                 | _____        | sample depth        | _____       |
| <u>soil</u>     | other    | ekman    | _____             | cond               | _____        | tide stage          | _____       |
| _____           | _____    | _____    | _____             | _____              | _____        | ambient temp        | <u>40°F</u> |
| _____           | _____    | _____    | _____             | _____              | _____        | barometric pressure | _____       |
| _____           | _____    | _____    | _____             | _____              | _____        | relative humidity   | _____       |
| _____           | _____    | _____    | _____             | _____              | _____        | weather conditions  | _____       |

**ANALYSES TO BE PERFORMED**

**SAMPLE PREPARATION**

**ORGANICS**

**OTHER ANALYSES**

**CONTAINER**

**PRESERVATIVES**

- A. halogenated & aromatic volatiles
- B. volatiles
- C. trihalomethanes
- D. pesticides/PCB
- E. PCB
- F. base neutral/acid extractables
- G. pesticides. drinking water
- H. herbicides. drinking water
- I. other \_\_\_\_\_

- A. total cyanide
- B. total phenol
- C. petroleum hydrocarbons
- D. pH
- E. alkalinity
- F. hardness
- G. total dissolved solids
- H. total suspended solids
- I. sulfate
- J. TOC
- K. grain size
- L. percent moisture
- M. other \_\_\_\_\_

- glass jar
- plastic jar
- acetate core
- plastic bag
- plastic bucket
- other \_\_\_\_\_

NaOH

- HNO<sub>3</sub>
- Zn Acetate
- HCl
- Na<sub>2</sub>SO<sub>4</sub>
- other \_\_\_\_\_

**INORGANICS**

**STORAGE**

- wet ice
- dry ice
- ambient

- A. metals. priority pollutant
- B. metals. TAL
- C. metals scan (ICP)
- D. metals. other \_\_\_\_\_

**RCRA**

- A. TCLP
- B. ignitability
- C. corrosivity \_\_\_\_\_ pH \_\_\_\_\_
- D. reactivity
- E. other \_\_\_\_\_

**COMMENTS**

# FIELD DATA SHEET

16654

**REAC, EDISON, NJ**  
**(908) 321-4200**  
**EPA CONTRACT 68-C4-0022**

Date: 4/14/66 Samplers: \_\_\_\_\_ Chain of Custody No.: \_\_\_\_\_  
 Time: 11:00 Site Name: Leckie Run REAC Task Leader: Fams  
 Sample Location: #4 Fence line EPA WAM: Compton  
 Work Assignment No.: \_\_\_\_\_

| SITE DESCRIPTION  |            |                   | SOIL TYPE   |              | SURFACE WATER |       | STREAM   |            | BOTTOM |         |
|-------------------|------------|-------------------|-------------|--------------|---------------|-------|----------|------------|--------|---------|
| landfill          | old field  | upland palustrine | rock        | clay         | color         | _____ | width    | _____      | rock   | silt    |
| <u>Industrial</u> | wooded     | lowland riverine  | gravel      | <u>muck</u>  | odor          | _____ | depth    | _____      | rubble | clay    |
| Commercial        | farmland   | lacustrine        | <u>sand</u> | loam         | flow          | _____ | velocity | _____ cm/s | gravel | organic |
| residential       | gully      |                   | silt        | peat         | direction     | _____ | pools    | _____ %    | shell  | other   |
| hedgerows         | floodplain |                   | color       | <u>Black</u> |               |       | riffles  | _____ %    | sand   |         |

| SAMPLE TYPE     |          | DEVICE   |                   | SAMPLE INFORMATION |       | WEATHER PARAMETERS |       |
|-----------------|----------|----------|-------------------|--------------------|-------|--------------------|-------|
| surface water   | effluent | kemmerer | ponar             | color              | _____ | pH                 | _____ |
| groundwater     | sludge   | trowel   | other <u>Spaw</u> | odor               | _____ | ORP                | _____ |
| potable water   | leachate | bucket   |                   | temp               | _____ | salinity           | _____ |
| <u>sediment</u> | waste    | auger    |                   | DO                 | _____ | sample depth       | _____ |
| <u>soil</u>     | other    | ekman    |                   | cond               | _____ | tide stage         | _____ |

**ANALYSES TO BE PERFORMED**

**ORGANICS**

- A. halogenated & aromatic volatiles
- B. volatiles
- C. trihalomethanes
- D. pesticides/PCB
- E. PCB
- F. base neutral/acid extractables
- G. pesticides, drinking water
- H. herbicides, drinking water
- I. other \_\_\_\_\_

**INORGANICS**

- A. metals, priority pollutant
- B. metals, TAL
- C. metals scan (ICP)
- D. metals, other \_\_\_\_\_

**RCRA**

- A. TCLP
- B. ignitability
- C. corrosivity \_\_\_\_\_ pH \_\_\_\_\_
- D. reactivity
- E. other \_\_\_\_\_

**OTHER ANALYSES**

- A. total cyanide
- B. total phenol
- C. petroleum hydrocarbons
- D. pH
- E. alkalinity
- F. hardness
- G. total dissolved solids
- H. total suspended solids
- I. sulfate
- J. TOC
- K. grain size
- L. percent moisture
- M. other \_\_\_\_\_

**SAMPLE PREPARATION**

**CONTAINER**

- glass jar
- plastic jar
- acetate core
- plastic bag
- plastic bucket
- other \_\_\_\_\_

**PRESERVATIVES**

- HNO<sub>3</sub>
- NaOH
- Zn Acetate
- HCl
- Na<sub>2</sub>SO<sub>4</sub>
- other \_\_\_\_\_

**STORAGE**

- wet ice
- dry ice
- ambient

**COMMENTS**



**FIELD DATA SHEET**

16655

**REAC, EDISON, NJ  
(908) 321-4200  
EPA CONTRACT 68-C4-0022**

Date: 4/14/99 Samplers: Fame, Metz, Major Chain of Custody No.: \_\_\_\_\_  
 Site Name: LACKAWANNA Foundry REAC Task Leader: Fame  
 Time: 1100 Sample Location: #5 Fence Line EPA WAM: Campaign  
 Work Assignment No.: 7-439

| SITE DESCRIPTION  |            |                   | SOIL TYPE   |              | SURFACE WATER |       | STREAM   |            | BOTTOM |         |
|-------------------|------------|-------------------|-------------|--------------|---------------|-------|----------|------------|--------|---------|
| landfill          | old field  | upland palustrine | rock        | clay         | color         | _____ | width    | _____      | rock   | silt    |
| <u>Industrial</u> | wooded     | lowland riverine  | gravel      | muck         | odor          | _____ | depth    | _____      | rubble | clay    |
| commercial        | farmland   | lacustrine        | <u>sand</u> | loam         | flow          | _____ | velocity | _____ cm/s | gravel | organic |
| residential       | gully      |                   | silt        | peat         | direction     | _____ | pools    | _____ %    | shell  | other   |
| hedgerows         | floodplain |                   | color       | <u>Black</u> |               |       | riffles  | _____ %    | sand   |         |

| SAMPLE TYPE   |          | DEVICE   |                    | SAMPLE INFORMATION |       | WEATHER PARAMETERS |       |
|---------------|----------|----------|--------------------|--------------------|-------|--------------------|-------|
| surface water | effluent | kemmerer | ponar              | color              | _____ | pH                 | _____ |
| groundwater   | sludge   | trowel   | other <u>spoon</u> | odor               | _____ | ORP                | _____ |
| potable water | leachate | bucket   |                    | temp               | _____ | salinity           | _____ |
| sediment      | waste    | auger    |                    | DO                 | _____ | sample depth       | _____ |
| <u>Soil</u>   | other    | ekman    |                    | cond               | _____ | tide stage         | _____ |

**ANALYSES TO BE PERFORMED**

- ORGANICS**
- A. halogenated & aromatic volatiles
  - B. volatiles
  - C. trihalomethanes
  - D. pesticides/PCB
  - E. PCB
  - F. base neutral/acid extractables
  - G. pesticides, drinking water
  - H. herbicides, drinking water
  - I. other \_\_\_\_\_

- INORGANICS**
- A. metals, priority pollutant
  - B. metals, TAL
  - C. metals scan (ICP)
  - D. metals, other \_\_\_\_\_

- RCRA**
- A. TCLP
  - B. ignitability
  - C. corrosivity \_\_\_\_\_ pH \_\_\_\_\_
  - D. reactivity
  - E. other \_\_\_\_\_

- OTHER ANALYSES**
- A. total cyanide
  - B. total phenol
  - C. petroleum hydrocarbons
  - D. pH
  - E. alkalinity
  - F. hardness
  - G. total dissolved solids
  - H. total suspended solids
  - I. sulfate
  - J. TOC
  - K. grain size
  - L. percent moisture
  - M. other \_\_\_\_\_

**SAMPLE PREPARATION**

- |                  |                                 |
|------------------|---------------------------------|
| <b>CONTAINER</b> | <b>PRESERVATIVES</b>            |
| glass jar        | HNO <sub>3</sub>                |
| plastic jar      | NaOH                            |
| acetate core     | Zn Acetate                      |
| plastic bag      | HCl                             |
| plastic bucket   | Na <sub>2</sub> SO <sub>4</sub> |
| other _____      | other _____                     |

- STORAGE**
- wet ice
  - dry ice
  - ambient

**COMMENTS**

**FIELD DATA SHEET**

16656

**REAC, EDISON, NJ  
(908) 321-4200  
EPA CONTRACT 68-C4-0022**

Date: 4/14/99 Samplers: Fame Metz, Nader Chain of Custody No.: \_\_\_\_\_  
 Time: 1100 Site Name: LOCKAWANA FOUNDATION REAC Task Leader: Fame  
 Sample Location: #6 FARM ROAD NEAR WHITE STAGE NW EPA WAM: Compton  
 Work Assignment No.: 3-439

| SITE DESCRIPTION  |            |                   | SOIL TYPE     |              | SURFACE WATER |       | STREAM   |            | BOTTOM |         |
|-------------------|------------|-------------------|---------------|--------------|---------------|-------|----------|------------|--------|---------|
| landfill          | old field  | upland palustrine | rock          | clay         | color         | _____ | width    | _____      | rock   | silt    |
| <u>industrial</u> | wooded     | lowland riverine  | <u>gravel</u> | muck         | odor          | _____ | depth    | _____      | rubble | clay    |
| commercial        | farmland   | lacustrine        | sand          | loam         | flow          | _____ | velocity | _____ cm/s | gravel | organic |
| residential       | gully      |                   | silt          | peat         | direction     | _____ | pools    | _____ %    | shell  | other   |
| hedgerows         | floodplain |                   | color         | <u>Black</u> |               |       | riffles  | _____ %    | sand   |         |

| SAMPLE TYPE   |          | DEVICE   |                   | SAMPLE INFORMATION |       | WEATHER PARAMETERS |       |
|---------------|----------|----------|-------------------|--------------------|-------|--------------------|-------|
| surface water | effluent | kemmerer | ponar             | color              | _____ | pH                 | _____ |
| groundwater   | sludge   | trowel   | other <u>poon</u> | odor               | _____ | ORP                | _____ |
| potable water | leachate | bucket   |                   | temp               | _____ | salinity           | _____ |
| sediment      | waste    | auger    |                   | DO                 | _____ | sample depth       | _____ |
| <u>soil</u>   | other    | ekman    |                   | cond               | _____ | tide stage         | _____ |

**ANALYSES TO BE PERFORMED**

**SAMPLE PREPARATION**

**ORGANICS**

- A. halogenated & aromatic volatiles
- B. volatiles
- C. trihalomethanes
- D. pesticides/PCB
- E. PCB
- F. base neutral/acid extractables
- G. pesticides, drinking water
- H. herbicides, drinking water
- I. other \_\_\_\_\_

**OTHER ANALYSES**

- A. total cyanide
- B. total phenol
- C. petroleum hydrocarbons
- D. pH
- E. alkalinity
- F. hardness
- G. total dissolved solids
- H. total suspended solids
- I. sulfate
- J. TOC
- K. grain size
- L. percent moisture
- M. other \_\_\_\_\_

**CONTAINER**

- glass jar
- plastic jar
- acetate core
- plastic bag
- plastic bucket
- other \_\_\_\_\_

**NaOH**

**PRESERVATIVES**

- HNO<sub>3</sub>
- Zn Acetate
- HCl
- Na<sub>2</sub>SO<sub>4</sub>
- other \_\_\_\_\_

**STORAGE**

- wet ice
- dry ice
- ambient

**INORGANICS**

- A. metals, priority pollutant
- B. metals, TAL
- C. metals scan (ICP)
- D. metals, other \_\_\_\_\_

**RCRA**

- A. TCLP
- B. ignitability
- C. corrosivity \_\_\_\_\_ pH \_\_\_\_\_
- D. reactivity
- E. other \_\_\_\_\_

**COMMENTS**

# FIELD DATA SHEET

16657

**REAC, EDISON, NJ**  
**(908) 321-4200**  
**EPA CONTRACT 68-C4-0022**

Date: 4/14/99 Samplers: Fama, Metz, Maja Chain of Custody No.: \_\_\_\_\_  
 Site Name: Lackawanna Foundry REAC Task Leader: Fama  
 Time: 1100 Sample Location: #7 North Side Center 5' away from building EPA WAM: Compagn  
 Work Assignment No.: 3-439

| SITE DESCRIPTION  |            |                   | SOIL TYPE   |                  | SURFACE WATER |       | STREAM   |            | BOTTOM |         |
|-------------------|------------|-------------------|-------------|------------------|---------------|-------|----------|------------|--------|---------|
| landfill          | old field  | upland palustrine | rock        | clay             | color         | _____ | width    | _____      | rock   | silt    |
| <u>industrial</u> | wooded     | lowland riverine  | gravel      | muck             | odor          | _____ | depth    | _____      | rubble | clay    |
| commercial        | farmland   | lacustrine        | sand        | loam             | flow          | _____ | velocity | _____ cm/s | gravel | organic |
| residential       | gully      |                   | <u>silt</u> | peat             | direction     | _____ | pools    | _____ %    | shell  | other   |
| hedgerows         | floodplain |                   | color       | <u>black-wet</u> |               |       | riffles  | _____ %    | sand   |         |

| SAMPLE TYPE   |          | DEVICE   |                    | SAMPLE INFORMATION |       | WEATHER PARAMETERS |       |
|---------------|----------|----------|--------------------|--------------------|-------|--------------------|-------|
| surface water | effluent | kemmerer | ponar              | color              | _____ | pH                 | _____ |
| groundwater   | sludge   | trowel   | other <u>spoon</u> | odor               | _____ | ORP                | _____ |
| potable water | leachate | bucket   |                    | temp               | _____ | salinity           | _____ |
| sediment      | waste    | auger    |                    | DO                 | _____ | sample depth       | _____ |
| <u>soil</u>   | other    | ekman    |                    | cond               | _____ | tide stage         | _____ |

**ANALYSES TO BE PERFORMED**

- ORGANICS**
- A. halogenated & aromatic volatiles
  - B. volatiles
  - C. trihalomethanes
  - D. pesticides/PCB
  - E. PCB
  - F. base neutral/acid extractables
  - G. pesticides, drinking water
  - H. herbicides, drinking water
  - I. other \_\_\_\_\_

- INORGANICS**
- A. metals, priority pollutant
  - B. metals, TAL
  - C. metals scan (ICP)
  - D. metals, other \_\_\_\_\_

- RCRA**
- A. TCLP
  - B. ignitability
  - C. corrosivity \_\_\_\_\_ pH \_\_\_\_\_
  - D. reactivity
  - E. other \_\_\_\_\_

- OTHER ANALYSES**
- A. total cyanide
  - B. total phenol
  - C. petroleum hydrocarbons
  - D. pH
  - E. alkalinity
  - F. hardness
  - G. total dissolved solids
  - H. total suspended solids
  - I. sulfate
  - J. TOC
  - K. grain size
  - L. percent moisture
  - M. other \_\_\_\_\_

**SAMPLE PREPARATION**

- CONTAINER**
- glass jar
  - plastic jar
  - acetate core
  - plastic bag
  - plastic bucket
  - other \_\_\_\_\_
- PRESERVATIVES**
- HNO<sub>3</sub>
  - NaOH
  - Zn Acetate
  - HCl
  - Na<sub>2</sub>SO<sub>4</sub>
  - other \_\_\_\_\_

- STORAGE**
- wet ice
  - dry ice
  - ambient

**COMMENTS**

**FIELD DATA SHEET**

16658

**REAC, EDISON, NJ  
(908) 321-4200  
EPA CONTRACT 68-C4-0022**

Date: 4/14/96 Samplers: Fame Metz, Magan Chain of Custody No.: \_\_\_\_\_  
 Site Name: Lackawanna Foundry REAC Task Leader: Fame  
 Time: 1100 Sample Location: # 8 North Side 6' from building EPA WAM: Compagnie  
 Work Assignment No.: 3-439

| SITE DESCRIPTION  |            |                   | SOIL TYPE |              | SURFACE WATER |       | STREAM   |            | BOTTOM |         |
|-------------------|------------|-------------------|-----------|--------------|---------------|-------|----------|------------|--------|---------|
| landfill          | old field  | upland palustrine | rock      | clay         | color         | _____ | width    | _____      | rock   | silt    |
| <u>industrial</u> | wooded     | lowland riverine  | gravel    | muck         | odor          | _____ | depth    | _____      | rubble | clay    |
| commercial        | farmland   | lacustrine        | sand      | loam         | flow          | _____ | velocity | _____ cm/s | gravel | organic |
| residential       | gully      |                   | silt      | peat         | direction     | _____ | pools    | _____ %    | shell  | other   |
| hedgerows         | floodplain |                   | color     | <u>Black</u> |               |       | riffles  | _____ %    | sand   |         |

| SAMPLE TYPE   |          | DEVICE   |                    | SAMPLE INFORMATION |       | WEATHER PARAMETERS |       |
|---------------|----------|----------|--------------------|--------------------|-------|--------------------|-------|
| surface water | effluent | kemmerer | ponar              | color              | _____ | pH                 | _____ |
| groundwater   | sludge   | trowel   | other <u>ifoon</u> | odor               | _____ | ORP                | _____ |
| potable water | leachate | bucket   |                    | temp               | _____ | salinity           | _____ |
| sediment      | waste    | auger    |                    | DO                 | _____ | sample depth       | _____ |
| <u>soil</u>   | other    | ekman    |                    | cond               | _____ | tide stage         | _____ |

**ANALYSES TO BE PERFORMED**

**ORGANICS**

- A. halogenated & aromatic volatiles
- B. volatiles
- C. trihalomethanes
- D. pesticides/PCB
- E. PCB
- F. base neutral/acid extractables
- G. pesticides, drinking water
- H. herbicides, drinking water
- I. other \_\_\_\_\_

**INORGANICS**

- A. metals, priority pollutant
- B. metals, TAL
- C. metals scan (ICP)
- D. metals, other \_\_\_\_\_

**RCRA**

- A. TCLP
- B. ignitability
- C. corrosivity \_\_\_\_\_ pH \_\_\_\_\_
- D. reactivity
- E. other \_\_\_\_\_

**OTHER ANALYSES**

- A. total cyanide
- B. total phenol
- C. petroleum hydrocarbons
- D. pH
- E. alkalinity
- F. hardness
- G. total dissolved solids
- H. total suspended solids
- I. sulfate
- J. TOC
- K. grain size
- L. percent moisture
- M. other \_\_\_\_\_

**SAMPLE PREPARATION**

**CONTAINER**

- glass jar
- plastic jar
- acetate core
- plastic bag
- plastic bucket
- other \_\_\_\_\_

NaOH

**PRESERVATIVES**

- HNO<sub>3</sub>
- Zn Acetate
- HCl
- Na<sub>2</sub>SO<sub>4</sub>
- other \_\_\_\_\_

**STORAGE**

- wet ice
- dry ice
- ambient

**COMMENTS**

**FIELD DATA SHEET**

16559

**REAC, EDISON, NJ  
(908) 321-4200  
EPA CONTRACT 68-C4-0022**

Date: 4/14/99 Samplers: Fama, Metz, Majum Chain of Custody No.: \_\_\_\_\_  
 Site Name: Lackawanna Foundry REAC Task Leader: Fama  
 Time: 1100 Sample Location: #9 North side middle courtyard EPA WAM: Comp/99  
 Work Assignment No.: 3-439

| SITE DESCRIPTION  |            |                   | SOIL TYPE   |              | SURFACE WATER |       | STREAM   |            | BOTTOM |         |
|-------------------|------------|-------------------|-------------|--------------|---------------|-------|----------|------------|--------|---------|
| landfill          | old field  | upland palustrine | rock        | clay         | color         | _____ | width    | _____      | rock   | silt    |
| <u>Industrial</u> | wooded     | lowland riverine  | gravel      | muck         | odor          | _____ | depth    | _____      | rubble | clay    |
| commercial        | farmland   | lacustrine        | sand        | loam         | flow          | _____ | velocity | _____ cm/s | gravel | organic |
| residential       | gully      |                   | <u>silt</u> | peat         | direction     | _____ | pools    | _____ %    | shell  | other   |
| hedgerows         | floodplain |                   | color       | <u>Black</u> |               |       | riffles  | _____ %    | sand   |         |

| SAMPLE TYPE   |          | DEVICE   |                   | SAMPLE INFORMATION |       | WEATHER PARAMETERS |       |
|---------------|----------|----------|-------------------|--------------------|-------|--------------------|-------|
| surface water | effluent | kemmerer | ponar             | color              | _____ | pH                 | _____ |
| groundwater   | sludge   | trowel   | other <u>SPUN</u> | odor               | _____ | ORP                | _____ |
| potable water | leachate | bucket   |                   | temp               | _____ | salinity           | _____ |
| sediment      | waste    | auger    |                   | DO                 | _____ | sample depth       | _____ |
| <u>soil</u>   | other    | ekman    |                   | cond               | _____ | tide stage         | _____ |

**ANALYSES TO BE PERFORMED**

- ORGANICS**
- A halogenated & aromatic volatiles
  - B volatiles
  - C trihalomethanes
  - D pesticides/PCB
  - E PCB
  - F base neutral/acid extractables
  - G pesticides, drinking water
  - H herbicides, drinking water
  - I other \_\_\_\_\_

- INORGANICS**
- A metals, priority pollutant
  - B metals, TAL
  - C metals scan (ICP)
  - D metals, other \_\_\_\_\_

- RCRA**
- A TCLP
  - B ignitability
  - C corrosivity \_\_\_\_\_ pH \_\_\_\_\_
  - D reactivity
  - E other \_\_\_\_\_

**SAMPLE PREPARATION**

- CONTAINER**
- glass jar
  - plastic jar
  - acetate core
  - plastic bag
  - plastic bucket
  - other \_\_\_\_\_
- NaOH**
- PRESERVATIVES**
- HNO<sub>3</sub>
  - Zn Acetate
  - HCl
  - Na<sub>2</sub>SO<sub>4</sub>
  - other \_\_\_\_\_

**STORAGE**

- wet ice
- dry ice
- ambient

**COMMENTS**

# FIELD DATA SHEET

16660

**REAC, EDISON, NJ**  
**(908) 321-4200**  
**EPA CONTRACT 68-C4-0022**

Date: 4/14/99 Samplers: Fams, Metz, Mager Chain of Custody No.: \_\_\_\_\_  
 Time: 1100 Site Name: Lackawanna Foundry REAC Task Leader: Fams  
 Sample Location: #10 inside building North Area EPA WAM: Compass  
 Work Assignment No.: 3-439

| SITE DESCRIPTION  |            |                   | SOIL TYPE |              | SURFACE WATER |       | STREAM   |            | BOTTOM |         |
|-------------------|------------|-------------------|-----------|--------------|---------------|-------|----------|------------|--------|---------|
| landfill          | old field  | upland palustrine | rock      | clay         | color         | _____ | width    | _____      | rock   | silt    |
| <u>industrial</u> | wooded     | lowland riverine  | gravel    | muck         | odor          | _____ | depth    | _____      | rubble | clay    |
| commercial        | farmland   | lacustrine        | sand      | loam         | flow          | _____ | velocity | _____ cm/s | gravel | organic |
| residential       | gully      |                   | silt      | peat         | direction     | _____ | pools    | _____ %    | shell  | other   |
| hedgerows         | floodplain |                   | color     | <u>Black</u> |               |       | riffles  | _____ %    | sand   |         |

| SAMPLE TYPE     |                        | DEVICE        |       | SAMPLE INFORMATION |       |              | WEATHER PARAMETERS |                     |       |
|-----------------|------------------------|---------------|-------|--------------------|-------|--------------|--------------------|---------------------|-------|
| surface water   | effluent               | kemmerer      | ponar | color              | _____ | pH           | _____              | ambient temp        | _____ |
| groundwater     | sludge                 | <u>trowel</u> | other | odor               | _____ | ORP          | _____              | barometric pressure | _____ |
| potable water   | leachate               | bucket        |       | temp               | _____ | salinity     | _____              | relative humidity   | _____ |
| <u>sediment</u> | waste                  | auger         |       | DO                 | _____ | sample depth | _____              | weather conditions  | _____ |
| soil            | other <u>Composite</u> | ekman         |       | cond               | _____ | tide stage   | _____              |                     |       |

**ANALYSES TO BE PERFORMED**

**ORGANICS**

- A. halogenated & aromatic volatiles
- B. volatiles
- C. trihalomethanes
- D. pesticides/PCB
- E. PCB
- F. base neutral/acid extractables
- G. pesticides, drinking water
- H. herbicides, drinking water
- I. other \_\_\_\_\_

**INORGANICS**

- A. metals, priority pollutant
- B. metals, TAL
- C. metals scan (ICP)
- D. metals, other \_\_\_\_\_

**RCRA**

- A. TCLP
- B. ignitability
- C. corrosivity \_\_\_\_\_ pH \_\_\_\_\_
- D. reactivity
- E. other \_\_\_\_\_

**OTHER ANALYSES**

- A. total cyanide
- B. total phenol
- C. petroleum hydrocarbons
- D. pH
- E. alkalinity
- F. hardness
- G. total dissolved solids
- H. total suspended solids
- I. sulfate
- J. TOC
- K. grain size
- L. percent moisture
- M. other \_\_\_\_\_

**SAMPLE PREPARATION**

**CONTAINER**

- glass jar
- plastic jar
- acetate core
- plastic bag
- plastic bucket
- other \_\_\_\_\_

NaOH

**PRESERVATIVES**

- HNO<sub>3</sub>
- Zn Acetate
- HCl
- Na<sub>2</sub>SO<sub>4</sub>
- other \_\_\_\_\_

**STORAGE**

- wet ice
- dry ice
- ambient

**COMMENTS**

FIELD DATA SHEET

16661

REAC, EDISON, NJ
(908) 321-4200
EPA CONTRACT 68-C4-0022

Date: 4/14/66 Samplers: Metz
Site Name: Luckenbach Foundry
Sample Location: #11 inside SW Area under conveyor belt
Chain of Custody No.:
REAC Task Leader: FGM
EPA WAM: Campaign
Work Assignment No.: 3-439

Table with 5 columns: SITE DESCRIPTION, SOIL TYPE, SURFACE WATER, STREAM, BOTTOM. Includes categories like landfill, industrial, commercial, residential, hedgerows, etc.

Table with 4 columns: SAMPLE TYPE, DEVICE, SAMPLE INFORMATION, WEATHER PARAMETERS. Includes categories like surface water, groundwater, potable water, sediment, soil, etc.

ANALYSES TO BE PERFORMED

- ORGANICS
A halogenated & aromatic volatiles
B volatiles
C trihalomethanes
D pesticides/PCB
E PCB
F base neutral/acid extractables
G pesticides, drinking water
H herbicides, drinking water
I other

- OTHER ANALYSES
A. total cyanide
B. total phenol
C. petroleum hydrocarbons
D. pH
E. alkalinity
F. hardness
G. total dissolved solids
H. total suspended solids
I. sulfate
J. TOC
K. grain size
L. percent moisture
M. other

SAMPLE PREPARATION

- CONTAINER
glass jar
plastic jar
acetate core
plastic bag
plastic bucket
other
PRESERVATIVES
HNO3
NaOH
Zn Acetate
HCl
Na2SO4
other
STORAGE
wet ice
dry ice
ambient

- INORGANICS
A metals, priority pollutant
B metals, TAL
C metals scan (ICP)
D metals, other

- RCRA
A TCLP
B ignitability
C corrosivity pH
D reactivity
E other

COMMENTS
under conveyor belt

**FIELD DATA SHEET**

16662

**REAC, EDISON, NJ  
(908) 321-4200  
EPA CONTRACT 68-C4-0022**

Date: 4/14/69 Samplers: Metz Chain of Custody No.: \_\_\_\_\_  
 Time: 1100 Site Name: Locke Wanne Foundry REAC Task Leader: Fams  
 Sample Location: #12 SE Low area EPA WAM: Campagna  
 Work Assignment No.: 3-439

| SITE DESCRIPTION  |            |                   | SOIL TYPE |                  | SURFACE WATER |       | STREAM   |            | BOTTOM |         |
|-------------------|------------|-------------------|-----------|------------------|---------------|-------|----------|------------|--------|---------|
| landfill          | old field  | upland palustrine | rock      | clay             | color         | _____ | width    | _____      | rock   | silt    |
| <u>industrial</u> | wooded     | lowland riverine  | gravel    | muck             | odor          | _____ | depth    | _____      | rubble | clay    |
| commercial        | farmland   | lacustrine        | sand      | loam             | flow          | _____ | velocity | _____ cm/s | gravel | organic |
| residential       | gully      |                   | silt      | peat             | direction     | _____ | pools    | _____ %    | shell  | other   |
| hedgerows         | floodplain |                   | color     | <u>Black-wet</u> |               |       | riffles  | _____ %    | sand   |         |

| SAMPLE TYPE   |                         | DEVICE   |       | SAMPLE INFORMATION |       | WEATHER PARAMETERS |       |
|---------------|-------------------------|----------|-------|--------------------|-------|--------------------|-------|
| surface water | effluent                | kemmerer | ponar | color              | _____ | pH                 | _____ |
| groundwater   | sludge                  | trowel   | other | odor               | _____ | ORP                | _____ |
| potable water | leachate                | bucket   |       | temp               | _____ | salinity           | _____ |
| sediment      | waste                   | auger    |       | DO                 | _____ | sample depth       | _____ |
| <u>soil</u>   | other <u>Comp's 1+0</u> | ekman    |       | cond               | _____ | tide stage         | _____ |

**ANALYSES TO BE PERFORMED**

- ORGANICS**
- A halogenated & aromatic volatiles
  - B volatiles
  - C trihalomethanes
  - D pesticides/PCB
  - E PCB
  - F base neutral/acid extractables
  - G pesticides, drinking water
  - H herbicides, drinking water
  - I other \_\_\_\_\_

- INORGANICS**
- A metals, priority pollutant
  - B metals, TAL
  - C metals scan (ICP)
  - D metals, other \_\_\_\_\_

- RCRA**
- A TCLP
  - B ignitability
  - C corrosivity \_\_\_\_\_ pH \_\_\_\_\_
  - D reactivity
  - E other \_\_\_\_\_

**COMMENTS**

**SAMPLE PREPARATION**

- |                  |                                 |
|------------------|---------------------------------|
| <b>CONTAINER</b> | <b>PRESERVATIVES</b>            |
| glass jar        | HNO <sub>3</sub>                |
| plastic jar      | NaOH                            |
| acetate core     | Zn Acetate                      |
| plastic bag      | HCl                             |
| plastic bucket   | Na <sub>2</sub> SO <sub>4</sub> |
| other            | other _____                     |

**STORAGE**

- J. TOC
- K. grain size
- L. percent moisture
- M. other \_\_\_\_\_



**FIELD DATA SHEET**

16663

**REAC, EDISON, NJ  
(908) 321-4200  
EPA CONTRACT 68-C4-0022**

Date: 9/14/99 Samplers: Metz Chain of Custody No.: \_\_\_\_\_  
 REAC Task Leader: Farms  
 Site Name: Lackawanna Foundry EPA WAM: Campaign  
 Time: 1:00 Sample Location: #13 SE under stack Work Assignment No.: 3-439

| SITE DESCRIPTION  |            |                   | SOIL TYPE                 | SURFACE WATER   | STREAM              | BOTTOM |             |
|-------------------|------------|-------------------|---------------------------|-----------------|---------------------|--------|-------------|
| landfill          | old field  | upland palustrine | rock clay                 | color _____     | width _____         | rock   | silt        |
| <u>industrial</u> | wooded     | lowland riverine  | gravel muck               | odor _____      | depth _____         | rubble | clay        |
| commercial        | farmland   | lacustrine        | sand loam                 | flow _____      | velocity _____ cm/s | gravel | organic     |
| residential       | gully      |                   | silt peat                 | direction _____ | pools _____ %       | shell  | other _____ |
| hedgerows         | floodplain |                   | color <u>Black + blue</u> |                 | riffles _____ %     | sand   |             |

| SAMPLE TYPE     |                      | DEVICE             | SAMPLE INFORMATION |                    | WEATHER PARAMETERS        |  |
|-----------------|----------------------|--------------------|--------------------|--------------------|---------------------------|--|
| surface water   | effluent             | kemmerer ponar     | color _____        | pH _____           | ambient temp _____        |  |
| groundwater     | sludge               | trowel other _____ | odor _____         | ORP _____          | barometric pressure _____ |  |
| potable water   | leachate             | bucket             | temp _____         | salinity _____     | relative humidity _____   |  |
| <u>sediment</u> | waste                | auger              | DO _____           | sample depth _____ | weather conditions _____  |  |
| <u>soil</u>     | other <u>Compost</u> | ekman              | cond _____         | tide stage _____   |                           |  |

**ANALYSES TO BE PERFORMED**

- ORGANICS**
- A. halogenated & aromatic volatiles
  - B. volatiles
  - C. trihalomethanes
  - D. pesticides/PCB
  - E. PCB
  - F. base neutral/acid extractables
  - G. pesticides, drinking water
  - H. herbicides, drinking water
  - I. other \_\_\_\_\_

- OTHER ANALYSES**
- A. total cyanide
  - B. total phenol
  - C. petroleum hydrocarbons
  - D. pH
  - E. alkalinity
  - F. hardness
  - G. total dissolved solids
  - H. total suspended solids
  - I. sulfate
  - J. TOC
  - K. grain size
  - L. percent moisture
  - M. other \_\_\_\_\_

- INORGANICS**
- A. metals, priority pollutant
  - B. metals, TAL
  - C. metals scan (ICP)
  - D. metals, other \_\_\_\_\_

- RCRA**
- A. TCLP
  - B. ignitability
  - C. corrosivity \_\_\_\_\_ pH \_\_\_\_\_
  - D. reactivity
  - E. other \_\_\_\_\_

**SAMPLE PREPARATION**

- CONTAINER**
- glass jar
  - plastic jar NaOH
  - acetate core
  - plastic bag
  - plastic bucket
  - other \_\_\_\_\_
- PRESERVATIVES**
- HNO<sub>3</sub>
  - Zn Acetate
  - HCl
  - Na<sub>2</sub>SO<sub>4</sub>
  - other \_\_\_\_\_

- STORAGE**
- wet ice
  - dry ice
  - ambient

**COMMENTS**

**FIELD DATA SHEET**

16664

**REAC, EDISON, NJ  
(908) 321-4200  
EPA CONTRACT 68-C4-0022**

Date: 4/14/99 Samplers: Fama Metz, Magan Chain of Custody No.: \_\_\_\_\_  
 Time: 1100 Site Name: Lackeyman Foundry REAC Task Leader: Fama  
 Sample Location: #14 low muddy driveway area EPA WAM: Compton  
 Work Assignment No.: 3-439

| SITE DESCRIPTION  |            |                   | SOIL TYPE |                  | SURFACE WATER |       | STREAM   |            | BOTTOM |         |
|-------------------|------------|-------------------|-----------|------------------|---------------|-------|----------|------------|--------|---------|
| landfill          | old field  | upland palustrine | rock      | clay             | color         | _____ | width    | _____      | rock   | silt    |
| <u>industrial</u> | wooded     | lowland riverine  | gravel    | muck             | odor          | _____ | depth    | _____      | rubble | clay    |
| commercial        | farmland   | lacustrine        | sand      | loam             | flow          | _____ | velocity | _____ cm/s | gravel | organic |
| residential       | gully      |                   | silt      | peat             | direction     | _____ | pools    | _____ %    | shell  | other   |
| hedgerows         | floodplain |                   | color     | <u>black-wet</u> |               |       | riffles  | _____ %    | sand   |         |

| SAMPLE TYPE   |          | DEVICE   |                    | SAMPLE INFORMATION |       | WEATHER PARAMETERS |       |
|---------------|----------|----------|--------------------|--------------------|-------|--------------------|-------|
| surface water | effluent | kemmerer | ponar              | color              | _____ | pH                 | _____ |
| groundwater   | sludge   | trowel   | other <u>spoon</u> | odor               | _____ | ORP                | _____ |
| potable water | leachate | bucket   |                    | temp               | _____ | salinity           | _____ |
| sediment      | waste    | auger    |                    | DO                 | _____ | sample depth       | _____ |
| <u>soil</u>   | other    | ekman    |                    | cond               | _____ | tide stage         | _____ |

**ANALYSES TO BE PERFORMED**

- ORGANICS**
- A. halogenated & aromatic volatiles
  - B. volatiles
  - C. trihalomethanes
  - D. pesticides/PCB
  - E. PCB
  - F. base neutral/acid extractables
  - G. pesticides, drinking water
  - H. herbicides, drinking water
  - I. other \_\_\_\_\_

- INORGANICS**
- A. metals, priority pollutant
  - B. metals, TAL
  - C. metals scan (ICP)
  - D. metals, other \_\_\_\_\_

- RCRA**
- A. TCLP
  - B. ignitability
  - C. corrosivity \_\_\_\_\_ pH \_\_\_\_\_
  - D. reactivity
  - E. other \_\_\_\_\_

- OTHER ANALYSES**
- A. total cyanide
  - B. total phenol
  - C. petroleum hydrocarbons
  - D. pH
  - E. alkalinity
  - F. hardness
  - G. total dissolved solids
  - H. total suspended solids
  - I. sulfate
  - J. TOC
  - K. grain size
  - L. percent moisture
  - M. other \_\_\_\_\_

**SAMPLE PREPARATION**

- CONTAINER**
- glass jar
  - plastic jar
  - acetate core
  - plastic bag
  - plastic bucket
  - other \_\_\_\_\_
- PRESERVATIVES**
- HNO<sub>3</sub>
  - NaOH
  - Zn Acetate
  - HCl
  - Na<sub>2</sub>SO<sub>4</sub>
  - other \_\_\_\_\_

- STORAGE**
- wet ice
  - dry ice
  - ambient

COMMENTS South of fence

**FIELD DATA SHEET**

16655

**REAC, EDISON, NJ  
(908) 321-4200  
EPA CONTRACT 68-C4-0022**

Date: 4/14/89 Samplers: Fama, Metz, Mayle Chain of Custody No.: \_\_\_\_\_  
 Site Name: Lackawanna Foundry REAC Task Leader: Fama  
 Time: 1100 Sample Location: #15 East grey material EPA WAM: Campaign  
 Work Assignment No.: 3-439

| SITE DESCRIPTION  |            |                   | SOIL TYPE |                   | SURFACE WATER |       | STREAM   |            | BOTTOM |         |
|-------------------|------------|-------------------|-----------|-------------------|---------------|-------|----------|------------|--------|---------|
| landfill          | old field  | upland palustrine | rock      | clay              | color         | _____ | width    | _____      | rock   | silt    |
| <u>industrial</u> | wooded     | lowland riverine  | gravel    | muck              | odor          | _____ | depth    | _____      | rubble | clay    |
| commercial        | farmland   | lacustrine        | sand      | loam              | flow          | _____ | velocity | _____ cm/s | gravel | organic |
| residential       | gully      |                   | silt      | peat              | direction     | _____ | pools    | _____ %    | shell  | other   |
| hedgerows         | floodplain |                   | color     | <u>grey-brown</u> |               |       | riffles  | _____ %    | sand   |         |

| SAMPLE TYPE     |          | DEVICE   |                   | SAMPLE INFORMATION |       | WEATHER PARAMETERS |       |
|-----------------|----------|----------|-------------------|--------------------|-------|--------------------|-------|
| surface water   | effluent | kemmerer | ponar             | color              | _____ | pH                 | _____ |
| groundwater     | sludge   | trowel   | other <u>spur</u> | odor               | _____ | ORP                | _____ |
| potable water   | leachate | bucket   |                   | temp               | _____ | salinity           | _____ |
| <u>sediment</u> | waste    | auger    |                   | DO                 | _____ | sample depth       | _____ |
| soil            | other    | ekman    |                   | cond               | _____ | tide stage         | _____ |

**ANALYSES TO BE PERFORMED**

- ORGANICS**
- A. halogenated & aromatic volatiles
  - B. volatiles
  - C. trihalomethanes
  - D. pesticides/PCB
  - E. PCB
  - F. base neutral/acid extractables
  - G. pesticides. drinking water
  - H. herbicides. drinking water
  - I. other \_\_\_\_\_

- INORGANICS**
- A. metals. priority pollutant
  - B. metals. TAL
  - C. metals scan (ICP)
  - D. metals. other \_\_\_\_\_

- RCRA**
- A. TCLP
  - B. ignitability
  - C. corrosivity \_\_\_\_\_ pH \_\_\_\_\_
  - D. reactivity
  - E. other \_\_\_\_\_

- OTHER ANALYSES**
- A. total cyanide
  - B. total phenol
  - C. petroleum hydrocarbons
  - D. pH
  - E. alkalinity
  - F. hardness
  - G. total dissolved solids
  - H. total suspended solids
  - I. sulfate
  - J. TOC
  - K. grain size
  - L. percent moisture
  - M. other \_\_\_\_\_

**SAMPLE PREPARATION**

- CONTAINER**
- glass jar
  - plastic jar NaOH
  - acetate core
  - plastic bag
  - plastic bucket
  - other \_\_\_\_\_
- PRESERVATIVES**
- HNO<sub>3</sub>
  - Zn Acetate
  - HCl
  - Na<sub>2</sub>SO<sub>4</sub>
  - other \_\_\_\_\_

**STORAGE**

- wet ice
- dry ice
- ambient

COMMENTS NE of building

FIELD DATA SHEET

16666

REAC, EDISON, NJ
(908) 321-4200
EPA CONTRACT 68-C4-0022

Date: 5/14/99 Samplers: Fama, Nayan, Metz
Site Name: Lackawanna Foundry
Time: 1100 Sample Location: #16 SW of East pile
Chain of Custody No.:
REAC Task Leader: Fama
EPA WAM: Campbell
Work Assignment No.: 3-439

Table with 7 columns: SITE DESCRIPTION, SOIL TYPE, SURFACE WATER, STREAM, BOTTOM. Includes categories like landfill, industrial, commercial, residential, hedgerows, old field, wooded, farmland, gully, floodplain, rock, clay, gravel, sand, silt, peat, color, color, muck, loam, flow, direction, width, depth, velocity, pools, riffles, rock, rubble, gravel, shell, sand, silt, clay, organic, other.

Table with 5 columns: SAMPLE TYPE, DEVICE, SAMPLE INFORMATION, WEATHER PARAMETERS. Includes categories like surface water, groundwater, potable water, sediment, soil, effluent, sludge, leachate, waste, other, kemmerer, trowel, bucket, auger, ekman, ponar, other, color, odor, temp, DO, cond, pH, ORP, salinity, sample depth, tide stage, ambient temp, barometric pressure, relative humidity, weather conditions.

ANALYSES TO BE PERFORMED

ORGANICS

- A. halogenated & aromatic volatiles
B. volatiles
C. trihalomethanes
D. pesticides/PCB
E. PCB
F. base neutral/acid extractables
G. pesticides, drinking water
H. herbicides, drinking water
I. other

OTHER ANALYSES

- A. total cyanide
B. total phenol
C. petroleum hydrocarbons
D. pH
E. alkalinity
F. hardness
G. total dissolved solids
H. total suspended solids
I. sulfate
J. TOC
K. grain size
L. percent moisture
M. other

SAMPLE PREPARATION

CONTAINER

- glass jar
plastic jar
acetate core
plastic bag
plastic bucket
other

PRESERVATIVES

- HNO3
NaOH
Zn Acetate
HCl
Na2SO4
other

STORAGE

- wet ice
dry ice
ambient

INORGANICS

- A. metals, priority pollutant
B. metals, TAL
C. metals scan (ICP)
D. metals, other

RCRA

- A. TCLP
B. ignitability
C. corrosivity pH
D. reactivity
E. other

COMMENTS

South west corner of east pile

**FIELD DATA SHEET**

16667

**REAC, EDISON, NJ  
(908) 321-4200  
EPA CONTRACT 68-C4-0022**

Date: 6/14/06      Samplers: Fama, Metz, Megan      Chain of Custody No.: \_\_\_\_\_  
 Time: 120      Site Name: Lackawanna Foundry      REAC Task Leader: Fama  
 Sample Location: #17 SE corner gw pile      EPA WAM: Campana      Work Assignment No.: 3-439

| SITE DESCRIPTION  |            |                   | SOIL TYPE          |      | SURFACE WATER   |                     | STREAM |        | BOTTOM      |  |
|-------------------|------------|-------------------|--------------------|------|-----------------|---------------------|--------|--------|-------------|--|
| landfill          | old field  | upland palustrine | rock               | clay | color _____     | width _____         |        | rock   | silt        |  |
| <u>industrial</u> | wooded     | lowland riverine  | gravel             | muck | odor _____      | depth _____         |        | rubble | clay        |  |
| commercial        | farmland   | lacustrine        | <u>sand</u>        | loam | flow _____      | velocity _____ cm/s |        | gravel | organic     |  |
| residential       | gully      |                   | silt               | peat | direction _____ | pools _____ %       |        | shell  | other _____ |  |
| hedgerows         | floodplain |                   | color <u>Brown</u> |      |                 | riffles _____ %     |        | sand   |             |  |

| SAMPLE TYPE   |             | DEVICE   |                   | SAMPLE INFORMATION |                    | WEATHER PARAMETERS        |  |
|---------------|-------------|----------|-------------------|--------------------|--------------------|---------------------------|--|
| surface water | effluent    | kemmerer | ponar             | color _____        | pH _____           | ambient temp _____        |  |
| groundwater   | sludge      | trowel   | other <u>span</u> | odor _____         | ORP _____          | barometric pressure _____ |  |
| potable water | leachate    | bucket   |                   | temp _____         | salinity _____     | relative humidity _____   |  |
| sediment      | waste       | auger    |                   | DO _____           | sample depth _____ | weather conditions _____  |  |
| <u>soil</u>   | other _____ | ekman    |                   | cond _____         | tide stage _____   |                           |  |

**ANALYSES TO BE PERFORMED**

- ORGANICS**
- A. halogenated & aromatic volatiles
  - B. volatiles
  - C. trihalomethanes
  - D. pesticides/PCB
  - E. PCB
  - F. base neutral/acid extractables
  - G. pesticides, drinking water
  - H. herbicides, drinking water
  - I. other \_\_\_\_\_
- INORGANICS**
- A. metals, priority pollutant
  - B. metals, TAL
  - C. metals scan (ICP)
  - D. metals, other \_\_\_\_\_

- OTHER ANALYSES**
- A. total cyanide
  - B. total phenol
  - C. petroleum hydrocarbons
  - D. pH
  - E. alkalinity
  - F. hardness
  - G. total dissolved solids
  - H. total suspended solids
  - I. sulfate
  - J. TOC
  - K. grain size
  - L. percent moisture
  - M. other \_\_\_\_\_

**SAMPLE PREPARATION**

- CONTAINER**
- glass jar
  - plastic jar
  - acetate core
  - plastic bag
  - plastic bucket
  - other \_\_\_\_\_
- STORAGE**
- wet ice
  - dry ice
  - ambient
- PRESERVATIVES**
- HNO<sub>3</sub>
  - NaOH
  - Zn Acetate
  - HCl
  - Na<sub>2</sub>SO<sub>4</sub>
  - other \_\_\_\_\_

- RCRA**
- A. TCLP
  - B. ignitability
  - C. corrosivity \_\_\_\_\_ pH \_\_\_\_\_
  - D. reactivity
  - E. other \_\_\_\_\_

COMMENTS Southeast corner west pile

FIELD DATA SHEET

16668

REAC, EDISON, NJ
(908) 321-4200
EPA CONTRACT 68-C4-0022

Date: 4/14/96
Time: 1100
Samplers: Farns, Metz, Magan
Site Name: Lackawanna
Sample Location: #18 east side of east pile
Chain of Custody No.:
REAC Task Leader: Farns
EPA WAM: Campaign
Work Assignment No.: 3-439

Table with 5 columns: SITE DESCRIPTION, SOIL TYPE, SURFACE WATER, STREAM, BOTTOM. Includes categories like landfill, industrial, commercial, residential, hedgerows, and soil types like rock, gravel, sand, silt, peat.

Table with 4 columns: SAMPLE TYPE, DEVICE, SAMPLE INFORMATION, WEATHER PARAMETERS. Includes sample types like surface water, groundwater, sediment, soil and various measurement parameters.

ANALYSES TO BE PERFORMED

ORGANICS

- A. halogenated & aromatic volatiles
B. volatiles
C. trihalomethanes
D. pesticides/PCB
E. PCB
F. base neutral/acid extractables
G. pesticides, drinking water
H. herbicides, drinking water
I. other

INORGANICS

- A. metals, priority pollutant
B. metals, TAL
C. metals scan (ICP)
D. metals, other

RCRA

- A. TCLP
B. ignitability
C. corrosivity pH
D. reactivity
E. other

OTHER ANALYSES

- A. total cyanide
B. total phenol
C. petroleum hydrocarbons
D. pH
E. alkalinity
F. hardness
G. total dissolved solids
H. total suspended solids
I. sulfate
J. TOC
K. grain size
L. percent moisture
M. other

SAMPLE PREPARATION

CONTAINER

- glass jar
plastic jar
acetate core
plastic bag
plastic bucket
other

STORAGE

- wet ice
dry ice
ambient

PRESERVATIVES

- HNO3
NaOH
Zn Acetate
HCl
Na2SO4
other

COMMENTS

east side of east pile

**FIELD DATA SHEET**

16659

**REAC, EDISON, NJ  
(908) 321-4200  
EPA CONTRACT 68-C4-0022**

Date: 7/14/02 Samplers: Fanni, Metz, Magan Chain of Custody No.: \_\_\_\_\_  
 Time: 1105 Site Name: Lackawanna Foundry REAC Task Leader: Fanni  
 Sample Location: #1 center west side of East pile EPA WAM: Campaign  
 Work Assignment No.: 3-439

| SITE DESCRIPTION |            |                   | SOIL TYPE |                  | SURFACE WATER |       | STREAM   |            | BOTTOM |         |
|------------------|------------|-------------------|-----------|------------------|---------------|-------|----------|------------|--------|---------|
| landfill         | old field  | upland palustrine | rock      | clay             | color         | _____ | width    | _____      | rock   | silt    |
| industrial       | wooded     | lowland riverine  | gravel    | muck             | odor          | _____ | depth    | _____      | rubble | clay    |
| commercial       | farmland   | lacustrine        | sand      | loam             | flow          | _____ | velocity | _____ cm/s | gravel | organic |
| residential      | gully      |                   | silt      | peat             | direction     | _____ | pools    | _____ %    | shell  | other   |
| hedgerows        | floodplain |                   | color     | <u>red-brown</u> |               |       | riffles  | _____ %    | sand   |         |

| SAMPLE TYPE   |          | DEVICE   |       | SAMPLE INFORMATION |       | WEATHER PARAMETERS |       |
|---------------|----------|----------|-------|--------------------|-------|--------------------|-------|
| surface water | effluent | kemmerer | ponar | color              | _____ | pH                 | _____ |
| groundwater   | sludge   | trowel   | other | odor               | _____ | ORP                | _____ |
| potable water | leachate | bucket   |       | temp               | _____ | salinity           | _____ |
| sediment      | waste    | auger    |       | DO                 | _____ | sample depth       | _____ |
| soil          | other    | ekman    |       | cond               | _____ | tide stage         | _____ |

**ANALYSES TO BE PERFORMED**

**ORGANICS**

- A halogenated & aromatic volatiles
- B volatiles
- C trihalomethanes
- D pesticides/PCB
- E PCB
- F base neutral/acid extractables
- G pesticides. drinking water
- H herbicides. drinking water
- I other \_\_\_\_\_

**INORGANICS**

- A metals. priority pollutant
- B metals. TAL
- C metals scan (ICP)
- D metals. other \_\_\_\_\_

**RCRA**

- A TCLP
- B ignitability
- C corrosivity \_\_\_\_\_ pH \_\_\_\_\_
- D reactivity
- E other \_\_\_\_\_

**OTHER ANALYSES**

- A. total cyanide
- B. total phenol
- C. petroleum hydrocarbons
- D. pH
- E. alkalinity
- F. hardness
- G. total dissolved solids
- H. total suspended solids
- I. sulfate
- J. TOC
- K. grain size
- L. percent moisture
- M. other \_\_\_\_\_

**SAMPLE PREPARATION**

**CONTAINER**

- glass jar
- plastic jar
- acetate core
- plastic bag
- plastic bucket
- other \_\_\_\_\_

NaOH

**PRESERVATIVES**

- HNO<sub>3</sub>
- Zn Acetate
- HCl
- Na<sub>2</sub>SO<sub>4</sub>
- other \_\_\_\_\_

**STORAGE**

- wet ice
- dry ice
- ambient

**COMMENTS**

*Center west side of East pile*

# FIELD DATA SHEET

16670

**REAC, EDISON, NJ  
(908) 321-4200  
EPA CONTRACT 68-C4-0022**

Date: 9/14/96 Samplers: Fama, Metz, Magan Chain of Custody No.: \_\_\_\_\_  
 Time: 1100 Site Name: Lackawanna Foundry REAC Task Leader: Fama  
 Sample Location: #20 North corner west pile EPA WAM: Compagne  
 Work Assignment No.: 3-439

| SITE DESCRIPTION                               |                                     |  | SOIL TYPE                                |                               | SURFACE WATER   |                     | STREAM                          |                                      | BOTTOM |  |
|--|-------------------------------------|--|--|-------------------------------|-----------------|---------------------|---------------------------------|--------------------------------------|--------|--|
| <input type="checkbox"/> landfill              | <input type="checkbox"/> old field  | <input type="checkbox"/> upland palustrine | <input type="checkbox"/> rock            | <input type="checkbox"/> clay | color _____     | width _____         | <input type="checkbox"/> rock   | <input type="checkbox"/> silt        |        |  |
| <input checked="" type="checkbox"/> industrial | <input type="checkbox"/> wooded     | <input type="checkbox"/> lowland riverine  | <input type="checkbox"/> gravel          | <input type="checkbox"/> muck | odor _____      | depth _____         | <input type="checkbox"/> rubble | <input type="checkbox"/> clay        |        |  |
| <input type="checkbox"/> commercial            | <input type="checkbox"/> farmland   | <input type="checkbox"/> lacustrine        | <input checked="" type="checkbox"/> sand | <input type="checkbox"/> loam | flow _____      | velocity _____ cm/s | <input type="checkbox"/> gravel | <input type="checkbox"/> organic     |        |  |
| <input type="checkbox"/> residential           | <input type="checkbox"/> gully      |  | <input type="checkbox"/> silt            | <input type="checkbox"/> peat | direction _____ | pools _____ %       | <input type="checkbox"/> shell  | <input type="checkbox"/> other _____ |        |  |
| <input type="checkbox"/> hedgerows             | <input type="checkbox"/> floodplain |  | color <u>Brown</u>                       |                               |                 | riffles _____ %     | <input type="checkbox"/> sand   |                                      |        |  |

| SAMPLE TYPE                              |                                      | DEVICE                            |   | SAMPLE INFORMATION |                    | WEATHER PARAMETERS        |  |
|--|--------------------------------------|-----------------------------------|---|--------------------|--------------------|---------------------------|--|
| <input type="checkbox"/> surface water   | <input type="checkbox"/> effluent    | <input type="checkbox"/> kemmerer | <input type="checkbox"/> ponar              | color _____        | pH _____           | ambient temp _____        |  |
| <input type="checkbox"/> groundwater     | <input type="checkbox"/> sludge      | <input type="checkbox"/> trowel   | <input type="checkbox"/> other <u>spoon</u> | odor _____         | ORP _____          | barometric pressure _____ |  |
| <input type="checkbox"/> potable water   | <input type="checkbox"/> leachate    | <input type="checkbox"/> bucket   |   | temp _____         | salinity _____     | relative humidity _____   |  |
| <input type="checkbox"/> sediment        | <input type="checkbox"/> waste       | <input type="checkbox"/> auger    |   | DO _____           | sample depth _____ | weather conditions _____  |  |
| <input checked="" type="checkbox"/> soil | <input type="checkbox"/> other _____ | <input type="checkbox"/> ekman    |   | cond _____         | tide stage _____   |                           |  |

**ANALYSES TO BE PERFORMED**

**ORGANICS**

- A. halogenated & aromatic volatiles
- B. volatiles
- C. trihalomethanes
- D. pesticides/PCB
- E. PCB
- F. base neutral/acid extractables
- G. pesticides, drinking water
- H. herbicides, drinking water
- I. other \_\_\_\_\_

**INORGANICS**

- A. metals, priority pollutant
- B. metals TAL
- C. metals scan (ICP)
- D. metals, other \_\_\_\_\_

**RCRA**

- A. TCLP
- B. ignitability
- C. corrosivity \_\_\_\_\_ pH \_\_\_\_\_
- D. reactivity
- E. other \_\_\_\_\_

**OTHER ANALYSES**

- A. total cyanide
- B. total phenol
- C. petroleum hydrocarbons
- D. pH
- E. alkalinity
- F. hardness
- G. total dissolved solids
- H. total suspended solids
- I. sulfate
- J. TOC
- K. grain size
- L. percent moisture
- M. other \_\_\_\_\_

**SAMPLE PREPARATION**

**CONTAINER**

- glass jar
- plastic jar
- acetate core
- plastic bag
- plastic bucket
- other \_\_\_\_\_

NaOH

**PRESERVATIVES**

- HNO<sub>3</sub>
- Zn Acetate
- HCl
- Na<sub>2</sub>SO<sub>4</sub>
- other \_\_\_\_\_

**STORAGE**

- wet ice
- dry ice
- ambient

**COMMENTS**

*North corner of west pile*



**FIELD DATA SHEET**

16671

**REAC, EDISON, NJ  
(908) 321-4200  
EPA CONTRACT 68-C4-0022**

Date: 4/14/06 Samplers: Fame, Metz, Major Chain of Custody No.: \_\_\_\_\_  
 Site Name: Lackawanna REAC Task Leader: Fame  
 Time: 1:00 Sample Location: #21 center pond area EPA WAM: Campaign  
 Work Assignment No.: 3-439

| SITE DESCRIPTION  |            |                   | SOIL TYPE |              | SURFACE WATER |       | STREAM   |            | BOTTOM |         |
|-------------------|------------|-------------------|-----------|--------------|---------------|-------|----------|------------|--------|---------|
| landfill          | old field  | upland palustrine | rock      | clay         | color         | _____ | width    | _____      | rock   | silt    |
| <u>industrial</u> | wooded     | lowland riverine  | gravel    | <u>muck</u>  | odor          | _____ | depth    | _____      | rubble | clay    |
| commercial        | farmland   | lacustrine        | sand      | loam         | flow          | _____ | velocity | _____ cm/s | gravel | organic |
| residential       | gully      |                   | silt      | peat         | direction     | _____ | pools    | _____ %    | shell  | other   |
| hedgerows         | floodplain |                   | color     | <u>Black</u> |               |       | riffles  | _____ %    | sand   |         |

| SAMPLE TYPE   |          | DEVICE        |       | SAMPLE INFORMATION |       | WEATHER PARAMETERS |       |
|---------------|----------|---------------|-------|--------------------|-------|--------------------|-------|
| surface water | effluent | kemmerer      | ponar | color              | _____ | pH                 | _____ |
| groundwater   | sludge   | <u>trowel</u> | other | odor               | _____ | ORP                | _____ |
| potable water | leachate | <u>bucket</u> |       | temp               | _____ | salinity           | _____ |
| sediment      | waste    | auger         |       | DO                 | _____ | sample depth       | _____ |
| <u>soil</u>   | other    | ekman         |       | cond               | _____ | tide stage         | _____ |

**ANALYSES TO BE PERFORMED**

**ORGANICS**

- A. halogenated & aromatic volatiles
- B. volatiles
- C. trihalomethanes
- D. pesticides/PCB
- E. PCB
- F. base neutral/acid extractables
- G. pesticides, drinking water
- H. herbicides, drinking water
- I. other \_\_\_\_\_

**INORGANICS**

- A. metals, priority pollutant
- B. metals, TAL
- C. metals scan (ICP)
- D. metals, other \_\_\_\_\_

**RCRA**

- A. TCLP
- B. ignitability
- C. corrosivity \_\_\_\_\_ pH \_\_\_\_\_
- D. reactivity
- E. other \_\_\_\_\_

**OTHER ANALYSES**

- A. total cyanide
- B. total phenol
- C. petroleum hydrocarbons
- D. pH
- E. alkalinity
- F. hardness
- G. total dissolved solids
- H. total suspended solids
- I. sulfate
- J. TOC
- K. grain size
- L. percent moisture
- M. other \_\_\_\_\_

**SAMPLE PREPARATION**

**CONTAINER**

- glass jar
- plastic jar
- acetate core
- plastic bag
- plastic bucket
- other \_\_\_\_\_

NaOH

**PRESERVATIVES**

- HNO<sub>3</sub>
- Zn Acetate
- HCl
- Na<sub>2</sub>SO<sub>4</sub>
- other \_\_\_\_\_

**STORAGE**

- wet ice
- dry ice
- ambient

**COMMENTS**

**FIELD DATA SHEET**

16672

**REAC, EDISON, NJ  
(908) 321-4200  
EPA CONTRACT 68-C4-0022**

Date: 4/14/99      Samplers: Fams, Metz, Major      Chain of Custody No.: \_\_\_\_\_  
 Time: 1100      Site Name: Lackawanna      REAC Task Leader: Fams  
 Sample Location: #22 12<sup>th</sup> South of building      EPA WAM: Campaign  
 Work Assignment No.: 3-439

| SITE DESCRIPTION  |            |                   | SOIL TYPE |              | SURFACE WATER |       | STREAM   |            | BOTTOM |         |
|-------------------|------------|-------------------|-----------|--------------|---------------|-------|----------|------------|--------|---------|
| landfill          | old field  | upland palustrine | rock      | clay         | color         | _____ | width    | _____      | rock   | silt    |
| <u>industrial</u> | wooded     | lowland riverine  | gravel    | muck         | odor          | _____ | depth    | _____      | rubble | clay    |
| commercial        | farmland   | lacustrine        | sand      | loam         | flow          | _____ | velocity | _____ cm/s | gravel | organic |
| residential       | gully      |                   | silt      | peat         | direction     | _____ | pools    | _____ %    | shell  | other   |
| hedgerows         | floodplain |                   | color     | <u>BLACK</u> |               |       | riffles  | _____ %    | sand   |         |

| SAMPLE TYPE   |          | DEVICE   |                   | SAMPLE INFORMATION |       | WEATHER PARAMETERS |       |
|---------------|----------|----------|-------------------|--------------------|-------|--------------------|-------|
| surface water | effluent | kemmerer | ponar             | color              | _____ | pH                 | _____ |
| groundwater   | sludge   | trowel   | other <u>SPAN</u> | odor               | _____ | ORP                | _____ |
| potable water | leachate | bucket   |                   | temp               | _____ | salinity           | _____ |
| sediment      | waste    | auger    |                   | DO                 | _____ | sample depth       | _____ |
| <u>soil</u>   | other    | ekman    |                   | cond               | _____ | tide stage         | _____ |

| ANALYSES TO BE PERFORMED            |  | SAMPLE PREPARATION        |                                 |
|-------------------------------------|--|---------------------------|---------------------------------|
| <b>ORGANICS</b>                     |  | <b>OTHER ANALYSES</b>     | <b>PRESERVATIVES</b>            |
| A. halogenated & aromatic volatiles |  | A. total cyanide          | HNO <sub>3</sub>                |
| B. volatiles                        |  | B. total phenol           | NaOH                            |
| C. trihalomethanes                  |  | C. petroleum hydrocarbons | Zn Acetate                      |
| D. pesticides/PCB                   |  | D. pH                     | HCl                             |
| E. PCB                              |  | E. alkalinity             | Na <sub>2</sub> SO <sub>4</sub> |
| F. base neutral/acid extractables   |  | F. hardness               | other _____                     |
| G. pesticides, drinking water       |  | G. total dissolved solids |                                 |
| H. herbicides, drinking water       |  | H. total suspended solids |                                 |
| I. other _____                      |  | I. sulfate                |                                 |
|                                     |  | J. TOC                    | <b>STORAGE</b>                  |
| <b>INORGANICS</b>                   |  | K. grain size             | wet ice                         |
| A. metals, priority pollutant       |  | L. percent moisture       | dry ice                         |
| B. metals, TAL                      |  | M. other _____            | ambient                         |
| C. metals scan (ICP)                |  |                           |                                 |
| D. metals, other _____              |  |                           |                                 |
| <b>RCRA</b>                         |  |                           |                                 |
| A. TCLP                             |  |                           |                                 |
| B. ignitability                     |  |                           |                                 |
| C. corrosivity _____ pH _____       |  |                           |                                 |
| D. reactivity                       |  |                           |                                 |
| E. other _____                      |  |                           |                                 |

COMMENTS

# FIELD DATA SHEET

16673

**REAC, EDISON, NJ**  
**(908) 321-4200**  
**EPA CONTRACT 68-C4-0022**

Date: 9/14/91      Samplers: Fanna, Magjan      Chain of Custody No.: \_\_\_\_\_  
 Time: 1100      Site Name: Leckwamm's      REAC Task Leader: Fanna  
 Sample Location: #23 12' southwest of building      EPA WAM: Campaign  
 Work Assignment No.: 3-439

| SITE DESCRIPTION  |            |                   | SOIL TYPE |              | SURFACE WATER |          | STREAM |        | BOTTOM  |  |
|-------------------|------------|-------------------|-----------|--------------|---------------|----------|--------|--------|---------|--|
| landfill          | old field  | upland palustrine | rock      | clay         | color         | width    |        | rock   | silt    |  |
| <u>industrial</u> | wooded     | lowland riverine  | gravel    | muck         | odor          | depth    |        | rubble | clay    |  |
| commercial        | farmland   | lacustrine        | sand      | loam         | flow          | velocity | cm/s   | gravel | organic |  |
| residential       | gully      |                   | silt      | peat         | direction     | pools    | %      | shell  | other   |  |
| hedgerows         | floodplain |                   | color     | <u>BROWN</u> |               | riffles  | %      | sand   |         |  |

| SAMPLE TYPE   |          | DEVICE   |                    | SAMPLE INFORMATION |              | WEATHER PARAMETERS  |  |
|---------------|----------|----------|--------------------|--------------------|--------------|---------------------|--|
| surface water | effluent | kemmerer | ponar              | color              | pH           | ambient temp        |  |
| groundwater   | sludge   | trowel   | other <u>2poon</u> | odor               | ORP          | barometric pressure |  |
| potable water | leachate | bucket   |                    | temp               | salinity     | relative humidity   |  |
| sediment      | waste    | auger    |                    | DO                 | sample depth | weather conditions  |  |
| <u>soil</u>   | other    | ekman    |                    | cond               | tide stage   |                     |  |

**ANALYSES TO BE PERFORMED**

- ORGANICS**
- A. halogenated & aromatic volatiles
  - B. volatiles
  - C. trihalomethanes
  - D. pesticides/PCB
  - E. PCB
  - F. base neutral/acid extractables
  - G. pesticides, drinking water
  - H. herbicides, drinking water
  - I. other \_\_\_\_\_

- INORGANICS**
- A. metals, priority pollutant
  - B. metals, TAL
  - C. metals scan (ICP)
  - D. metals, other \_\_\_\_\_

- RCRA**
- A. TCLP
  - B. ignitability
  - C. corrosivity \_\_\_\_\_ pH \_\_\_\_\_
  - D. reactivity
  - E. other \_\_\_\_\_

- OTHER ANALYSES**
- A. total cyanide
  - B. total phenol
  - C. petroleum hydrocarbons
  - D. pH
  - E. alkalinity
  - F. hardness
  - G. total dissolved solids
  - H. total suspended solids
  - I. sulfate
  - J. TOC
  - K. grain size
  - L. percent moisture
  - M. other \_\_\_\_\_

**SAMPLE PREPARATION**

- |                  |                                 |
|------------------|---------------------------------|
| <b>CONTAINER</b> | <b>PRESERVATIVES</b>            |
| glass jar        | HNO <sub>3</sub>                |
| plastic jar      | NaOH                            |
| acetate core     | Zn Acetate                      |
| plastic bag      | HCl                             |
| plastic bucket   | Na <sub>2</sub> SO <sub>4</sub> |
| other _____      | other _____                     |
| <b>STORAGE</b>   |                                 |
| wet ice          |                                 |
| dry ice          |                                 |
| ambient          |                                 |

**COMMENTS**

12' south of west side of building

**FIELD DATA SHEET**

16674

**REAC, EDISON, NJ  
(908) 321-4200  
EPA CONTRACT 68-C4-0022**

Date: 4/14/00 Samplers: Fame, Majan Chain of Custody No.: \_\_\_\_\_  
 Time: 1100 Site Name: Lackawanna Foundry REAC Task Leader: Fame  
 Sample Location: #24 5' from southwest corner EPA WAM: Campaign  
 Work Assignment No.: 3-439

| SITE DESCRIPTION  |            |                   | SOIL TYPE |              | SURFACE WATER |       | STREAM   |            | BOTTOM |         |
|-------------------|------------|-------------------|-----------|--------------|---------------|-------|----------|------------|--------|---------|
| landfill          | old field  | upland palustrine | rock      | clay         | color         | _____ | width    | _____      | rock   | silt    |
| <u>industrial</u> | wooded     | lowland riverine  | gravel    | muck         | odor          | _____ | depth    | _____      | rubble | clay    |
| commercial        | farmland   | lacustrine        | sand      | loam         | flow          | _____ | velocity | _____ cm/s | gravel | organic |
| residential       | gully      |                   | silt      | peat         | direction     | _____ | pools    | _____ %    | shell  | other   |
| hedgerows         | floodplain |                   | color     | <u>black</u> |               |       | riffles  | _____ %    | sand   |         |

| SAMPLE TYPE   |          | DEVICE   |                   | SAMPLE INFORMATION |       | WEATHER PARAMETERS |       |
|---------------|----------|----------|-------------------|--------------------|-------|--------------------|-------|
| surface water | effluent | kemmerer | ponar             | color              | _____ | pH                 | _____ |
| groundwater   | sludge   | trowel   | other <u>span</u> | odor               | _____ | ORP                | _____ |
| potable water | leachate | bucket   |                   | temp               | _____ | salinity           | _____ |
| sediment      | waste    | auger    |                   | DO                 | _____ | sample depth       | _____ |
| <u>soil</u>   | other    | ekman    |                   | cond               | _____ | tide stage         | _____ |

**ANALYSES TO BE PERFORMED**

- ORGANICS**
- A. halogenated & aromatic volatiles
  - B. volatiles
  - C. trihalomethanes
  - D. pesticides/PCB
  - E. PCB
  - F. base neutral/acid extractables
  - G. pesticides, drinking water
  - H. herbicides, drinking water
  - I. other \_\_\_\_\_

- INORGANICS**
- A. metals, priority pollutant
  - B. metals, TAL
  - C. metals scan (ICP)
  - D. metals, other \_\_\_\_\_

- RCRA**
- A. TCLP
  - B. ignitability
  - C. corrosivity \_\_\_\_\_ pH \_\_\_\_\_
  - D. reactivity
  - E. other \_\_\_\_\_

- OTHER ANALYSES**
- A. total cyanide
  - B. total phenol
  - C. petroleum hydrocarbons
  - D. pH
  - E. alkalinity
  - F. hardness
  - G. total dissolved solids
  - H. total suspended solids
  - I. sulfate
  - J. TOC
  - K. grain size
  - L. percent moisture
  - M. other \_\_\_\_\_

**SAMPLE PREPARATION**

- CONTAINER**
- glass jar
  - plastic jar
  - acetate core
  - plastic bag
  - plastic bucket
  - other \_\_\_\_\_
- PRESERVATIVES**
- HNO<sub>3</sub>
  - NaOH
  - Zn Acetate
  - HCl
  - Na<sub>2</sub>SO<sub>4</sub>
  - other \_\_\_\_\_

- STORAGE**
- wet ice
  - dry ice
  - ambient

**COMMENTS**

5' from southwest corner of building

**FIELD DATA SHEET**

16675

**REAC, EDISON, NJ  
(908) 321-4200  
EPA CONTRACT 68-C4-0022**

Date: 4/14/99 Samplers: Fama, Majan Chain of Custody No.: \_\_\_\_\_  
 Site Name: Lactanone Foundry REAC Task Leader: Fama  
 Time: 1105 Sample Location: #25 surface south of building EPA WAM: Campyng  
 Work Assignment No.: 7-439

| SITE DESCRIPTION |            |                   | SOIL TYPE |              | SURFACE WATER |          | STREAM |        | BOTTOM  |  |
|------------------|------------|-------------------|-----------|--------------|---------------|----------|--------|--------|---------|--|
| landfill         | old field  | upland palustrine | rock      | clay         | color         | width    | rock   | silt   |         |  |
| industrial       | wooded     | lowland riverine  | gravel    | muck         | odor          | depth    | rubble | clay   |         |  |
| commercial       | farmland   | lacustrine        | sand      | loam         | flow          | velocity | cm/s   | gravel | organic |  |
| residential      | gully      |                   | silt      | peat         | direction     | pools    | %      | shell  | other   |  |
| hedgerows        | floodplain |                   | color     | <u>Black</u> |               | riffles  | %      | sand   |         |  |

| SAMPLE TYPE   |          | DEVICE       |       | SAMPLE INFORMATION |              | WEATHER PARAMETERS  |  |
|---------------|----------|--------------|-------|--------------------|--------------|---------------------|--|
| surface water | effluent | kemmerer     | ponar | color              | pH           | ambient temp        |  |
| groundwater   | sludge   | trowel       | other | odor               | ORP          | barometric pressure |  |
| potable water | leachate | bucket       |       | temp               | salinity     | relative humidity   |  |
| sediment      | waste    | <u>auger</u> |       | DO                 | sample depth | weather conditions  |  |
| <u>soil</u>   | other    | ekman        |       | cond               | tide stage   |                     |  |

**ANALYSES TO BE PERFORMED**

- ORGANICS**
- A. halogenated & aromatic volatiles
  - B. volatiles
  - C. trihalomethanes
  - D. pesticides/PCB
  - E. PCB
  - F. base neutral/acid extractables
  - G. pesticides, drinking water
  - H. herbicides, drinking water
  - I. other \_\_\_\_\_

- INORGANICS**
- A. metals, priority pollutant
  - B. metals, TAL
  - C. metals scan (ICP)
  - D. metals, other \_\_\_\_\_

- RCRA**
- A. TCLP
  - B. ignitability
  - C. corrosivity \_\_\_\_\_ pH \_\_\_\_\_
  - D. reactivity
  - E. other \_\_\_\_\_

- OTHER ANALYSES**
- A. total cyanide
  - B. total phenol
  - C. petroleum hydrocarbons
  - D. pH
  - E. alkalinity
  - F. hardness
  - G. total dissolved solids
  - H. total suspended solids
  - I. sulfate
  - J. TOC
  - K. grain size
  - L. percent moisture
  - M. other \_\_\_\_\_

**SAMPLE PREPARATION**

- CONTAINER**
- glass jar
  - plastic jar
  - acetate core
  - plastic bag
  - plastic bucket
  - other \_\_\_\_\_
- PRESERVATIVES**
- HNO<sub>3</sub>
  - NaOH
  - Zn Acetate
  - HCl
  - Na<sub>2</sub>SO<sub>4</sub>
  - other \_\_\_\_\_
- STORAGE**
- wet ice
  - dry ice
  - ambient

COMMENTS: 12' south of building on south west side

**FIELD DATA SHEET**

16676

**REAC, EDISON, NJ  
(908) 321-4200  
EPA CONTRACT 68-C4-0022**

Date: 4/14/99 Samplers: Farrar, Megan Chain of Custody No.: \_\_\_\_\_  
 Site Name: Lackawanna Fanchy REAC Task Leader: Farrar  
 Time: 1100 Sample Location: #250 subsurface south of building EPA WAM: Campaign  
 Work Assignment No.: 7-439

| SITE DESCRIPTION  |            |                   | SOIL TYPE          | SURFACE WATER   | STREAM              | BOTTOM |             |
|-------------------|------------|-------------------|--------------------|-----------------|---------------------|--------|-------------|
| landfill          | old field  | upland palustrine | rock clay          | color _____     | width _____         | rock   | silt        |
| <u>industrial</u> | wooded     | lowland riverine  | gravel muck        | odor _____      | depth _____         | rubble | clay        |
| commercial        | farmland   | lacustrine        | sand loam          | flow _____      | velocity _____ cm/s | gravel | organic     |
| residential       | gully      |                   | silt peat          | direction _____ | pools _____ %       | shell  | other _____ |
| hedgerows         | floodplain |                   | color <u>Black</u> |                 | riffles _____ %     | sand   |             |

| SAMPLE TYPE   |             | DEVICE       | SAMPLE INFORMATION |             | WEATHER PARAMETERS |                           |
|---------------|-------------|--------------|--------------------|-------------|--------------------|---------------------------|
| surface water | effluent    | kemmerer     | ponar              | color _____ | pH _____           | ambient temp _____        |
| groundwater   | sludge      | trowel       | other _____        | odor _____  | ORP _____          | barometric pressure _____ |
| potable water | leachate    | bucket       |                    | temp _____  | salinity _____     | relative humidity _____   |
| sediment      | waste       | <u>auger</u> |                    | DO _____    | sample depth _____ | weather conditions _____  |
| <u>soil</u>   | other _____ | ekman        |                    | cond _____  | tide stage _____   |                           |

**ANALYSES TO BE PERFORMED**

**ORGANICS**

- A. halogenated & aromatic volatiles
- B. volatiles
- C. trihalomethanes
- D. pesticides/PCB
- E. PCB
- F. base neutral/acid extractables
- G. pesticides, drinking water
- H. herbicides, drinking water
- I. other \_\_\_\_\_

**INORGANICS**

- A. metals, priority pollutant
- B. metals, TAL
- C. metals scan (ICP)
- D. metals, other \_\_\_\_\_

**RCRA**

- A. TCLP
- B. ignitability
- C. corrosivity \_\_\_\_\_ pH \_\_\_\_\_
- D. reactivity
- E. other \_\_\_\_\_

**OTHER ANALYSES**

- A. total cyanide
- B. total phenol
- C. petroleum hydrocarbons
- D. pH
- E. alkalinity
- F. hardness
- G. total dissolved solids
- H. total suspended solids
- I. sulfate
- J. TOC
- K. grain size
- L. percent moisture
- M. other \_\_\_\_\_

**SAMPLE PREPARATION**

**CONTAINER**

- glass jar
- plastic jar
- acetate core
- plastic bag
- plastic bucket
- other \_\_\_\_\_

NaOH

**PRESERVATIVES**

- HNO<sub>3</sub>
- Zn Acetate
- HCl
- Na<sub>2</sub>SO<sub>4</sub>
- other \_\_\_\_\_

**STORAGE**

- wet ice
- dry ice
- ambient

**COMMENTS**

12' south of building on south west side

**FIELD DATA SHEET**

16577

**REAC, EDISON, NJ  
(908) 321-4200  
EPA CONTRACT 68-C4-0022**

Date: 4/14/95      Samplers: Fanni, Magan      Chain of Custody No.: \_\_\_\_\_  
 Time: 1100      Site Name: Lackawanna      REAC Task Leader: Fanni  
 Sample Location: #26 Surface 12' South of building      EPA WAM: Compton  
 Work Assignment No.: \_\_\_\_\_

| SITE DESCRIPTION  |            |                   | SOIL TYPE |       | SURFACE WATER |       | STREAM   |            | BOTTOM |         |
|-------------------|------------|-------------------|-----------|-------|---------------|-------|----------|------------|--------|---------|
| landfill          | old field  | upland palustrine | rock      | clay  | color         | _____ | width    | _____      | rock   | silt    |
| <u>industrial</u> | wooded     | lowland riverine  | gravel    | muck  | odor          | _____ | depth    | _____      | rubble | clay    |
| commercial        | farmland   | lacustrine        | sand      | loam  | flow          | _____ | velocity | _____ cm/s | gravel | organic |
| residential       | gully      |                   | silt      | peat  | direction     | _____ | pools    | _____ %    | shell  | other   |
| hedgerows         | floodplain |                   | color     | _____ |               |       | riffles  | _____ %    | sand   |         |

| SAMPLE TYPE     |          | DEVICE   |                   | SAMPLE INFORMATION |       | WEATHER PARAMETERS |       |
|-----------------|----------|----------|-------------------|--------------------|-------|--------------------|-------|
| surface water   | effluent | kemmerer | ponar             | color              | _____ | pH                 | _____ |
| groundwater     | sludge   | trowel   | other <u>span</u> | odor               | _____ | ORP                | _____ |
| potable water   | leachate | bucket   |                   | temp               | _____ | salinity           | _____ |
| <u>sediment</u> | waste    | auger    |                   | DO                 | _____ | sample depth       | _____ |
| soil            | other    | ekman    |                   | cond               | _____ | tide stage         | _____ |

**ANALYSES TO BE PERFORMED**

- ORGANICS**
- A. halogenated & aromatic volatiles
  - B. volatiles
  - C. trihalomethanes
  - D. pesticides/PCB
  - E. PCB
  - F. base neutral/acid extractables
  - G. pesticides, drinking water
  - H. herbicides, drinking water
  - I. other \_\_\_\_\_

- INORGANICS**
- A. metals, priority pollutant
  - B. metals, TAL
  - C. metals scan (ICP)
  - D. metals, other \_\_\_\_\_

- RCRA**
- A. TCLP
  - B. ignitability
  - C. corrosivity \_\_\_\_\_ pH \_\_\_\_\_
  - D. reactivity
  - E. other \_\_\_\_\_

- OTHER ANALYSES**
- A. total cyanide
  - B. total phenol
  - C. petroleum hydrocarbons
  - D. pH
  - E. alkalinity
  - F. hardness
  - G. total dissolved solids
  - H. total suspended solids
  - I. sulfate
  - J. TOC
  - K. grain size
  - L. percent moisture
  - M. other \_\_\_\_\_

**SAMPLE PREPARATION**

- |                  |                                 |
|------------------|---------------------------------|
| <b>CONTAINER</b> | <b>PRESERVATIVES</b>            |
| glass jar        | HNO <sub>3</sub>                |
| plastic jar      | NaOH                            |
| acetate core     | Zn Acetate                      |
| plastic bag      | HCl                             |
| plastic bucket   | Na <sub>2</sub> SO <sub>4</sub> |
| other _____      | other _____                     |

**STORAGE**

- wet ice
- dry ice
- ambient

**COMMENTS**

12' south of southeast section of building

**FIELD DATA SHEET**

16678

**REAC, EDISON, NJ  
(908) 321-4200  
EPA CONTRACT 68-C4-0022**

Date: 9/14/99 Samplers: Fame, Megan Chain of Custody No.: \_\_\_\_\_  
 Site Name: Lackawanna Foundry REAC Task Leader: Fame  
 Time: 1:00 Sample Location: #2615 subsurface EPA WAM: Campaign  
 Work Assignment No.: 3-439

| SITE DESCRIPTION  |            |                   | SOIL TYPE          |      | SURFACE WATER   |                     | STREAM |        | BOTTOM      |  |
|-------------------|------------|-------------------|--------------------|------|-----------------|---------------------|--------|--------|-------------|--|
| landfill          | old field  | upland palustrine | rock               | clay | color _____     | width _____         |        | rock   | silt        |  |
| <u>industrial</u> | wooded     | lowland riverine  | gravel             | muck | odor _____      | depth _____         |        | rubble | clay        |  |
| commercial        | farmland   | lacustrine        | sand               | loam | flow _____      | velocity _____ cm/s |        | gravel | organic     |  |
| residential       | gully      |                   | silt               | peat | direction _____ | pools _____ %       |        | shell  | other _____ |  |
| hedgerows         | floodplain |                   | color <u>black</u> |      |                 | riffles _____ %     |        | sand   |             |  |

| SAMPLE TYPE   |             | DEVICE       |             | SAMPLE INFORMATION |                    | WEATHER PARAMETERS        |  |
|---------------|-------------|--------------|-------------|--------------------|--------------------|---------------------------|--|
| surface water | effluent    | kemmerer     | ponar       | color _____        | pH _____           | ambient temp _____        |  |
| groundwater   | sludge      | trowel       | other _____ | odor _____         | ORP _____          | barometric pressure _____ |  |
| potable water | leachate    | bucket       |             | temp _____         | salinity _____     | relative humidity _____   |  |
| sediment      | waste       | <u>auger</u> |             | DO _____           | sample depth _____ | weather conditions _____  |  |
| <u>soil</u>   | other _____ | ekman        |             | cond _____         | tide stage _____   |                           |  |

**ANALYSES TO BE PERFORMED**

**ORGANICS**

- A. halogenated & aromatic volatiles
- B. volatiles
- C. trihalomethanes
- D. pesticides/PCB
- E. PCB
- F. base neutral/acid extractables
- G. pesticides, drinking water
- H. herbicides, drinking water
- I. other \_\_\_\_\_

**INORGANICS**

- A. metals, priority pollutant
- B. metals, TAL
- C. metals scan (ICP)
- D. metals other \_\_\_\_\_

**RCRA**

- A. TCLP
- B. ignitability
- C. corrosivity \_\_\_\_\_ pH \_\_\_\_\_
- D. reactivity
- E. other \_\_\_\_\_

**OTHER ANALYSES**

- A. total cyanide
- B. total phenol
- C. petroleum hydrocarbons
- D. pH
- E. alkalinity
- F. hardness
- G. total dissolved solids
- H. total suspended solids
- I. sulfate
- J. TOC
- K. grain size
- L. percent moisture
- M. other \_\_\_\_\_

**SAMPLE PREPARATION**

**CONTAINER**

- glass jar
- plastic jar
- acetate core
- plastic bag
- plastic bucket
- other \_\_\_\_\_

NaOH

**PRESERVATIVES**

- HNO<sub>3</sub>
- Zn Acetate
- HCl
- Na<sub>2</sub>SO<sub>4</sub>
- other \_\_\_\_\_

**STORAGE**

- wet ice
- dry ice
- ambient

**COMMENTS**

*12' south of south east section of building*



# FIELD DATA SHEET

16679

**REAC, EDISON, NJ  
(908) 321-4200  
EPA CONTRACT 68-C4-0022**

Date: 4/14/60 Samplers: Fama Metz Chain of Custody No.: \_\_\_\_\_  
 Time: 1100 Site Name: Lackawanna Foundry REAC Task Leader: Fama  
 Sample Location: #27 Composite east wing inside EPA WAM: Campaign  
 Work Assignment No.: 3-439

| SITE DESCRIPTION  |            |                   | SOIL TYPE |       | SURFACE WATER |       | STREAM   |            | BOTTOM |         |
|-------------------|------------|-------------------|-----------|-------|---------------|-------|----------|------------|--------|---------|
| landfill          | old field  | upland palustrine | rock      | clay  | color         | _____ | width    | _____      | rock   | silt    |
| <u>industrial</u> | wooded     | lowland riverine  | gravel    | muck  | odor          | _____ | depth    | _____      | rubble | clay    |
| commercial        | farmland   | lacustrine        | sand      | loam  | flow          | _____ | velocity | _____ cm/s | gravel | organic |
| residential       | gully      |                   | silt      | peat  | direction     | _____ | pools    | _____ %    | shell  | other   |
| hedgerows         | floodplain |                   | color     | _____ |               |       | riffles  | _____ %    | sand   |         |

| SAMPLE TYPE     |          | DEVICE        |       | SAMPLE INFORMATION |       | WEATHER PARAMETERS |       |
|-----------------|----------|---------------|-------|--------------------|-------|--------------------|-------|
| surface water   | effluent | kemmerer      | ponar | color              | _____ | pH                 | _____ |
| groundwater     | sludge   | <u>trowel</u> | other | odor               | _____ | ORP                | _____ |
| potable water   | leachate | bucket        |       | temp               | _____ | salinity           | _____ |
| <u>sediment</u> | waste    | auger         |       | DO                 | _____ | sample depth       | _____ |
| soil            | other    | ekman         |       | cond               | _____ | tide stage         | _____ |

**ANALYSES TO BE PERFORMED**

**ORGANICS**

- A. halogenated & aromatic volatiles
- B. volatiles
- C. trihalomethanes
- D. pesticides/PCB
- E. PCB
- F. base neutral/acid extractables
- G. pesticides, drinking water
- H. herbicides, drinking water
- I. other \_\_\_\_\_

**INORGANICS**

- A. metals, priority pollutant
- B. metals, TAL
- C. metals scan (ICP)
- D. metals, other \_\_\_\_\_

**RCRA**

- A. TCLP
- B. ignitability
- C. corrosivity \_\_\_\_\_ pH \_\_\_\_\_
- D. reactivity
- E. other \_\_\_\_\_

**OTHER ANALYSES**

- A. total cyanide
- B. total phenol
- C. petroleum hydrocarbons
- D. pH
- E. alkalinity
- F. hardness
- G. total dissolved solids
- H. total suspended solids
- I. sulfate
- J. TOC
- K. grain size
- L. percent moisture
- M. other \_\_\_\_\_

**SAMPLE PREPARATION**

**CONTAINER**

- glass jar
- plastic jar
- acetate core
- plastic bag
- plastic bucket
- other \_\_\_\_\_

NaOH

**PRESERVATIVES**

- HNO<sub>3</sub>
- Zn Acetate
- HCl
- Na<sub>2</sub>SO<sub>4</sub>
- other \_\_\_\_\_

**STORAGE**

- wet ice
- dry ice
- ambient

**COMMENTS**

**FIELD DATA SHEET**

16680

**REAC, EDISON, NJ  
(908) 321-4200  
EPA CONTRACT 68-C4-0022**

Date: 4/14/99 Samplers: \_\_\_\_\_ Chain of Custody No.: \_\_\_\_\_  
 Site Name: Lackhanna Fuvdry REAC Task Leader: Famc  
 Time: \_\_\_\_\_ Sample Location: Field/Blank/Trip Blank (VOC) EPA WAM: Compagné  
 Work Assignment No.: 3-439

| SITE DESCRIPTION  |            |                   | SOIL TYPE |       | SURFACE WATER |       | STREAM   |            | BOTTOM |         |
|-------------------|------------|-------------------|-----------|-------|---------------|-------|----------|------------|--------|---------|
| landfill          | old field  | upland palustrine | rock      | clay  | color         | _____ | width    | _____      | rock   | silt    |
| <u>industrial</u> | wooded     | lowland riverine  | gravel    | muck  | odor          | _____ | depth    | _____      | rubble | clay    |
| commercial        | farmland   | lacustrine        | sand      | loam  | flow          | _____ | velocity | _____ cm/s | gravel | organic |
| residential       | gully      |                   | silt      | peat  | direction     | _____ | pools    | _____ %    | shell  | other   |
| hedgerows         | floodplain |                   | color     | _____ |               |       | riffles  | _____ %    | sand   |         |

| SAMPLE TYPE   |                   | DEVICE   |       | SAMPLE INFORMATION |       | WEATHER PARAMETERS |       |
|---------------|-------------------|----------|-------|--------------------|-------|--------------------|-------|
| surface water | effluent          | kemmerer | ponar | color              | _____ | pH                 | _____ |
| groundwater   | sludge            | trowel   | other | odor               | _____ | ORP                | _____ |
| potable water | leachate          | bucket   |       | temp               | _____ | salinity           | _____ |
| sediment      | waste             | auger    |       | DO                 | _____ | sample depth       | _____ |
| soil          | other <u>Sand</u> | ekman    |       | cond               | _____ | tide stage         | _____ |

**ANALYSES TO BE PERFORMED**

- ORGANICS**
- A. halogenated & aromatic volatiles
  - B. volatiles
  - C. trihalomethanes
  - D. pesticides/PCB
  - E. PCB
  - F. base neutral/acid extractables
  - G. pesticides, drinking water
  - H. herbicides, drinking water
  - I. other \_\_\_\_\_

- INORGANICS**
- A. metals, priority pollutant
  - B. metals, TAL
  - C. metals scan (ICP)
  - D. metals, other \_\_\_\_\_

- RCRA**
- A. TCLP
  - B. ignitability
  - C. corrosivity \_\_\_\_\_ pH \_\_\_\_\_
  - D. reactivity
  - E. other \_\_\_\_\_

- OTHER ANALYSES**
- A. total cyanide
  - B. total phenol
  - C. petroleum hydrocarbons
  - D. pH
  - E. alkalinity
  - F. hardness
  - G. total dissolved solids
  - H. total suspended solids
  - I. sulfate
  - J. TOC
  - K. grain size
  - L. percent moisture
  - M. other \_\_\_\_\_

**SAMPLE PREPARATION**

- CONTAINER**
- glass jar
  - plastic jar
  - acetate core
  - plastic bag
  - plastic bucket
  - other \_\_\_\_\_
- PRESERVATIVES**
- HNO<sub>3</sub>
  - NaOH
  - Zn Acetate
  - HCl
  - Na<sub>2</sub>SO<sub>4</sub>
  - other \_\_\_\_\_

- STORAGE**
- wet ice
  - dry ice
  - ambient

**COMMENTS**

Field Blank (VOC) A 16680  
 Trip Blank (VOC) B 16680

## Appendix C

APPENDIX C  
CHAIN OF CUSTODY RECORDS  
LACKAWANNA FOUNDRY SITE  
LACKAWANNA, NEW YORK  
MAY 1999

**Sample Identification**

**Analyses Requested**

| REAC # | Sample No. | Sampling Location | Matrix | Date Collected | # of Bottles | Container/Preservative | PVA | AS | AR | VOC | Trace Metals |
|--------|------------|-------------------|--------|----------------|--------------|------------------------|-----|----|----|-----|--------------|
|        | C16656     | 644R EN100        | S      | 4/14/96        | 1            | 4oz/40C                |     |    |    |     |              |
|        | A16657     | 7                 |        |                |              | 8oz/40C                |     |    |    |     |              |
|        | 316657     | 7                 |        |                |              | 40ml/40C               |     |    |    |     |              |
|        | C16657     | 7                 |        |                |              | 4oz/40C                |     |    |    |     |              |
|        | A16658     | 8                 |        |                |              | 8oz/40C                |     |    |    |     |              |
|        | 316658     | 8                 |        |                |              | 40ml/40C               |     |    |    |     |              |
|        | C16658     | 8                 |        |                |              | 4oz/40C                |     |    |    |     |              |
|        | A16659     | 9                 |        |                |              | 8oz/40C                |     |    |    |     |              |
|        | 316659     | 9                 |        |                |              | 40ml/40C               |     |    |    |     |              |
|        | C16659     | 9                 |        |                |              | 4oz/40C                |     |    |    |     |              |
|        | A16660     | 10                |        |                |              | 8oz/40C                |     |    |    |     |              |
|        | 316660     | 10                |        |                |              | 40ml/40C               |     |    |    |     |              |
|        | C16660     | 10                |        |                |              | 4oz/40C                |     |    |    |     |              |
|        | A16661     | 11                |        |                |              | 8oz/40C                |     |    |    |     |              |
|        | 316661     | 11                |        |                |              | 40ml/40C               |     |    |    |     |              |
|        | C16661     | 11                |        |                |              | 4oz/40C                |     |    |    |     |              |
|        | A16662     | 12                |        |                |              | 8oz/40C                |     |    |    |     |              |
|        | 316662     | 12                |        |                |              | 40ml/40C               |     |    |    |     |              |
|        | C16662     | 12                |        |                |              | 4oz/40C                |     |    |    |     |              |
|        | A16663     | 13                |        |                |              | 8oz/40C                |     |    |    |     |              |

Matrix:  
 SD - Sediment  
 DS - Drum Solids  
 DL - Drum Liquids  
 X - Other

PW - Potable Water  
 GW - Groundwater  
 SW - Surface Water  
 SL - Sludge

S - Soil  
 W - Water  
 O - Oil  
 A - Air

Special Instructions:  
 Lab to select m/s/m/s/D  
 Water BVA - Kion new method available  
 PEST - pesticides  
 DCM - Polychlorinated biphenyls  
 VOC - volatile organic compounds

**FOR SUBCONTRACTING USE ONLY**  
**FROM CHAIN OF CUSTODY #**

| Items/Reason    | Relinquished By   | Date          | Received By | Date | Time | Items/Reason | Relinquished By | Date | Received By | Date | Time |
|-----------------|-------------------|---------------|-------------|------|------|--------------|-----------------|------|-------------|------|------|
| <i>Hand 511</i> | <i>John Jones</i> | <i>6/1/95</i> |             |      |      |              |                 |      |             |      |      |
|                 |                   |               |             |      |      |              |                 |      |             |      |      |
|                 |                   |               |             |      |      |              |                 |      |             |      |      |
|                 |                   |               |             |      |      |              |                 |      |             |      |      |
|                 |                   |               |             |      |      |              |                 |      |             |      |      |

Project Name: Leak & Seep Tank Trucking Site  
 Project Number: 05-97-143-01-3439-01  
 RFW Contact: Jackson/Fane Phone: 732-321-4800

**Sample Identification**

**Analyses Requested**

| REAC # | Sample No. | Sampling Location | Matrix | Date Collected | # of Bottles | Container/Preservative | BVA, RES, RB | VOC | TA, METALS |
|--------|------------|-------------------|--------|----------------|--------------|------------------------|--------------|-----|------------|
|        | A16650     | 1 West Fence      | S      | 4/14/99        | 1            | 8oz/40c                | ✓            |     |            |
|        | B16650     | 1 West Fence      | S      | 4/14/99        | 1            | 40ml/40c               |              | ✓   |            |
|        | C16650     | 1 West Fence      | S      | 4/14/99        | 1            | 4oz/40c                |              |     | ✓          |
|        | A16651     | 2 Surface Middle  | S      | 4/14/99        | 1            | 8oz/40c                | ✓            |     |            |
|        | B16651     | 2 Surface Middle  | S      | 4/14/99        | 1            | 40ml/40c               |              | ✓   |            |
|        | C16651     | 2 Surface Middle  | S      | 4/14/99        | 1            | 4oz/40c                |              |     | ✓          |
|        | A16652     | 2D Middle         | S      | 4/14/99        | 1            | 8oz/40c                | ✓            |     |            |
|        | B16652     | 2D Middle         | S      | 4/14/99        | 1            | 40ml/40c               |              | ✓   |            |
|        | C16652     | 2D Middle         | S      | 4/14/99        | 1            | 4oz/40c                |              |     | ✓          |
|        | A16653     | 3 West Fence      | S      | 4/14/99        | 1            | 8oz/40c                | ✓            |     |            |
|        | B16653     | 3 West Fence      | S      | 4/14/99        | 1            | 40ml/40c               |              | ✓   |            |
|        | C16653     | 3 West Fence      | S      | 4/14/99        | 1            | 4oz/40c                |              |     | ✓          |
|        | A16654     | 4 West Fence      | S      | 4/14/99        | 1            | 8oz/40c                | ✓            |     |            |
|        | B16654     | 4 West Fence      | S      | 4/14/99        | 1            | 40ml/40c               |              | ✓   |            |
|        | C16654     | 4 West Fence      | S      | 4/14/99        | 1            | 4oz/40c                |              |     | ✓          |
|        | A16655     | 5 West Fence      | S      | 4/14/99        | 1            | 8oz/40c                | ✓            |     |            |
|        | B16655     | 5 West Fence      | S      | 4/14/99        | 1            | 40ml/40c               |              | ✓   |            |
|        | C16655     | 5 West Fence      | S      | 4/14/99        | 1            | 4oz/40c                |              |     | ✓          |
|        | A16656     | 6 West Entrance   | S      | 4/14/99        | 1            | 8oz/40c                | ✓            |     |            |
|        | B16656     | 6 West Entrance   | S      | 4/14/99        | 1            | 40ml/40c               |              | ✓   |            |
|        | C16656     | 6 West Entrance   | S      | 4/14/99        | 1            | 4oz/40c                |              |     | ✓          |

Special Instructions: Lab to check MS/MSD

**Matrix:**  
 SD - Sediment  
 DS - Drum Solids  
 DL - Drum Liquids  
 X - Other

PW - Potable Water  
 GW - Groundwater  
 SW - Surface Water  
 SL - Sludge

S - Soil  
 W - Water  
 O - Oil  
 A - Air

VOC - Volatile organic compounds  
 M - Millifilter

Post-pesticides  
 PCP - Polychlorinated biphenyls

FOR SUBCONTRACTING USE ONLY  
 FROM CHAIN OF CUSTODY #

| Item/Reason | Relinquished By | Date    | Received By | Date | Time | Items/Reason | Relinquished By | Date | Received By | Date | Time |
|-------------|-----------------|---------|-------------|------|------|--------------|-----------------|------|-------------|------|------|
| Hand 1      | J. Fane         | 4/14/99 |             |      |      |              |                 |      |             |      |      |
|             |                 |         |             |      |      |              |                 |      |             |      |      |
|             |                 |         |             |      |      |              |                 |      |             |      |      |
|             |                 |         |             |      |      |              |                 |      |             |      |      |
|             |                 |         |             |      |      |              |                 |      |             |      |      |
|             |                 |         |             |      |      |              |                 |      |             |      |      |
|             |                 |         |             |      |      |              |                 |      |             |      |      |
|             |                 |         |             |      |      |              |                 |      |             |      |      |
|             |                 |         |             |      |      |              |                 |      |             |      |      |
|             |                 |         |             |      |      |              |                 |      |             |      |      |







Sample Identification

Analyses Requested

| REAC # | Sample No. | Sampling Location | Matrix | Date Collected | # of Bottles | Container/Preservative | SW/RES/RES | VOCS | WETALS |
|--------|------------|-------------------|--------|----------------|--------------|------------------------|------------|------|--------|
|        | A16672     | LOCATION 22       | S      | 14-APR-99      | 1            | 402 glass / ICE        | ✓          |      | ✓      |
|        | A16673     | LOCATION 23       |        |                | 1            | 402 glass /            | ✓          |      | ✓      |
|        | B16673     | ↓                 |        |                | 1            | 402 glass /            |            |      | ✓      |
|        | C16673     | ↓                 |        |                | 1            | 402 glass /            |            |      | ✓      |
|        | A16674     | LOCATION 24       |        |                | 1            | 802 /                  | ✓          |      | ✓      |
|        | B16674     | ↓                 |        |                | 1            | 402 ml /               |            |      | ✓      |
|        | C16674     | ↓                 |        |                | 1            | 402 /                  |            |      | ✓      |
|        | A16675     | LOCATION 25       |        |                | 1            | 802 /                  | ✓          |      | ✓      |
|        | B16675     | ↓                 |        |                | 1            | 402 ml /               |            |      | ✓      |
|        | C16675     | ↓                 |        |                | 1            | 402 /                  |            |      | ✓      |
|        | A16676     | LOCATION 25-D     |        |                | 1            | 802 /                  | ✓          |      | ✓      |
|        | B16676     | ↓                 |        |                | 1            | 402 ml /               |            |      | ✓      |
|        | C16676     | ↓                 |        |                | 1            | 402 /                  |            |      | ✓      |
|        | A16677     | LOCATION 26       |        |                | 1            | 802 /                  | ✓          |      | ✓      |
|        | B16677     | ↓                 |        |                | 1            | 402 ml /               |            |      | ✓      |
|        | C16677     | ↓                 |        |                | 1            | 402 /                  |            |      | ✓      |
|        | A16678     | LOCATION 26-D     |        |                | 1            | 802 /                  | ✓          |      | ✓      |
|        | B16678     | ↓                 |        |                | 1            | 402 ml /               |            |      | ✓      |
|        | C16678     | ↓                 |        |                | 1            | 402 /                  |            |      | ✓      |
|        | A16679     | LOCATION 27       |        |                | 1            | 802                    | ✓          |      | ✓      |

Matrix:  
SD - Sediment  
DS - Drum Solids  
DL - Drum Liquids  
X - Other

PW - Potable Water  
GW - Groundwater  
SL - Surface Water  
St - Sludge

S - Soil  
W - Water  
O - Oil  
A - Air

Special Instructions: Lab to send m/s/msd

GC-MS analysis of organic compounds  
ACR Polychlorinated biphenyls

FOR SUBCONTRACTING USE ONLY  
FROM CHAIN OF CUSTODY #

| Item/Reason                         | Relinquished By | Date    | Received By | Date | Time | Item/Reason | Relinquished By | Date | Received By | Date | Time |
|-------------------------------------|-----------------|---------|-------------|------|------|-------------|-----------------|------|-------------|------|------|
| GC/MS analysis of organic compounds | [Signature]     | 4/14/99 |             |      |      |             |                 |      |             |      |      |

REAC, L John, NJ  
 (908) 321-4200  
 EPA Contract 68-C4-0022

Project Name: LACALVAWA FOUNDRY  
 Project Number: 03517 H3 001 3439 01  
 RFW Contact: Sal FAWA Phone: 732 321 4200

No: 06913

**CH. OF CUSTODY RECORD**

**Sample Identification**

**Analyses Requested**

| REAC # | Sample No. | Sampling Location | Matrix | Date Collected | # of Bottles | Container/Preservative | PAHs/Res/KB | VOCs | Metals |
|--------|------------|-------------------|--------|----------------|--------------|------------------------|-------------|------|--------|
|        | 316679     | Location 27       | S      | 14-APR-99      | 1            | 40 ml glass / TCE      |             | ✓    |        |
|        | 316679     | ↓                 |        |                | 1            | 40 ml glass / TCE      |             |      | ✓      |
|        | 316680     | Field Pump (VOC)  |        |                | 1            | 40 ml /                |             | ✓    |        |
|        | 316680     | TRAP Pump (VOC)   |        |                | 1            | 40 ml /                |             | ✓    |        |

Matrix:  
 SD - Sediment  
 DS - Drum Solids  
 DL - Drum Liquids  
 X - Other

PW - Potable Water  
 GW - Groundwater  
 SW - Surface Water  
 SL - Sludge

S - Soil  
 W - Water  
 O - Oil  
 A - Air

Special Instructions: Lab to select m/s/m/s/D  
RNA - bear reagent placed extractable  
PEST - Pesticides  
PCB - poly chlorinated biphenyl  
OZ - ounces  
ml - milliliters

**FOR SUBCONTRACTING USE ONLY**  
**FROM CHAIN OF CUSTODY #**

| Items/Reason | Relinquished By | Date    | Received By | Date | Time | Items/Reason | Relinquished By | Date | Received By | Date | Time |
|--------------|-----------------|---------|-------------|------|------|--------------|-----------------|------|-------------|------|------|
| all analysis | Sal FAWA        | 4/14/99 |             |      |      |              |                 |      |             |      |      |
|              |                 |         |             |      |      |              |                 |      |             |      |      |
|              |                 |         |             |      |      |              |                 |      |             |      |      |
|              |                 |         |             |      |      |              |                 |      |             |      |      |
|              |                 |         |             |      |      |              |                 |      |             |      |      |

**Sample Identification**

**Analyses Requested**

| REAC # | Sample No. | Sampling Location | Matrix | Date Collected | # of Bottles | Container/Preservative | Analysis Requested |
|--------|------------|-------------------|--------|----------------|--------------|------------------------|--------------------|
|        | D16670     | Location 20       | S      | 14-APR-99      | 1            | 2 (R) Glass Ice        | CYANIDE            |
|        | D16671     | 21                |        |                | 1            |                        |                    |
|        | D16672     | 22                |        |                | 1            |                        |                    |
|        | D16673     | 23                |        |                | 1            |                        |                    |
|        | D16674     | 24                |        |                | 1            |                        |                    |
|        | D16675     | 25                |        |                | 1            |                        |                    |
|        | D16676     | 25-D              |        |                | 1            |                        |                    |
|        | D16677     | 26                |        |                | 1            |                        |                    |
|        | D16678     | 26-D              |        |                | 1            |                        |                    |
|        | D16679     | 27                |        |                | 1            |                        |                    |

Matrix:  
 SD - Sediment  
 DS - Drum Solids  
 DL - Drum Liquids  
 X - Other

PW - Potable Water  
 GW - Groundwater  
 SW - Surface Water  
 SL - Sludge

S - Soil  
 W - Water  
 O - Oil  
 A - Air

Special Instructions:

**FOR SUBCONTRACTING USE ONLY**  
**FROM CHAIN OF CUSTODY #**

| Items/Reason        | Relinquished By | Date           | Received By | Date | Time | Items/Reason | Relinquished By | Date | Received By | Date | Time |
|---------------------|-----------------|----------------|-------------|------|------|--------------|-----------------|------|-------------|------|------|
| <i>all analysis</i> | <i>Sal Fawa</i> | <i>4/29/99</i> |             |      |      |              |                 |      |             |      |      |
|                     |                 |                |             |      |      |              |                 |      |             |      |      |
|                     |                 |                |             |      |      |              |                 |      |             |      |      |

Project Name: LAKAWANNA FOUNDRY  
 Project Number: 03341 143 061 3439 01  
 RFW Contact: SALVATORE FAJVA Phone: 732 321 4200

**CHAIN OF CUSTODY RECORD**

**Sample Identification**

**Analyses Requested**

| REAC # | Sample No. | Sampling Location | Matrix | Date Collected | # of Bottles | Container/Preservative | Analysis Requested |
|--------|------------|-------------------|--------|----------------|--------------|------------------------|--------------------|
|        | D16650     | Location 1        | S      | 14-APR-99      | 1            | Roz Glass / Ice        | ✓                  |
|        | D16651     |                   |        |                | 1            |                        | ✓                  |
|        | D16652     |                   |        |                | 1            |                        | ✓                  |
|        | D16653     |                   |        |                | 1            |                        | ✓                  |
|        | D16654     |                   |        |                | 1            |                        | ✓                  |
|        | D16655     |                   |        |                | 1            |                        | ✓                  |
|        | D16656     |                   |        |                | 1            |                        | ✓                  |
|        | D16657     |                   |        |                | 1            |                        | ✓                  |
|        | D16658     |                   |        |                | 1            |                        | ✓                  |
|        | D16659     |                   |        |                | 1            |                        | ✓                  |
|        | D16660     |                   |        |                | 1            |                        | ✓                  |
|        | D16661     |                   |        |                | 1            |                        | ✓                  |
|        | D16662     |                   |        |                | 1            |                        | ✓                  |
|        | D16663     |                   |        |                | 1            |                        | ✓                  |
|        | D16664     |                   |        |                | 1            |                        | ✓                  |
|        | D16665     |                   |        |                | 1            |                        | ✓                  |
|        | D16666     |                   |        |                | 1            |                        | ✓                  |
|        | D16667     |                   |        |                | 1            |                        | ✓                  |
|        | D16668     |                   |        |                | 1            |                        | ✓                  |
|        | D16669     |                   |        |                | 1            |                        | ✓                  |
|        | D16670     |                   |        |                | 1            |                        | ✓                  |

Matrix:  
 SD - Sediment  
 DS - Drum Solids  
 DL - Drum Liquids  
 X - Other

PW - Potable Water  
 GW - Groundwater  
 SW - Surface Water  
 SL - Sludge

S - Soil  
 W - Water  
 O - Oil  
 A - Air

Special Instructions:

**FOR SUBCONTRACTING USE ONLY**  
**FROM CHAIN OF CUSTODY #**

| Items/Reason                  | Relinquished By  | Date           | Received By | Date | Time | Items/Reason | Relinquished By | Date | Received By | Date | Time |
|-------------------------------|------------------|----------------|-------------|------|------|--------------|-----------------|------|-------------|------|------|
| <i>all analysis performed</i> | <i>Salvatore</i> | <i>4/14/99</i> |             |      |      |              |                 |      |             |      |      |
|                               |                  |                |             |      |      |              |                 |      |             |      |      |
|                               |                  |                |             |      |      |              |                 |      |             |      |      |
|                               |                  |                |             |      |      |              |                 |      |             |      |      |

## **Appendix D**

APPENDIX D  
FINAL ANALYTICAL REPORT  
LACKAWANNA FOUNDRY SITE  
LACKAWANNA, NEW YORK  
MAY 1999

APPENDIX D  
FINAL ANALYTICAL REPORT  
LACKAWANNA FOUNDRY SITE  
LACKAWANNA, NEW YORK  
MAY 1999

ANALYTICAL REPORT

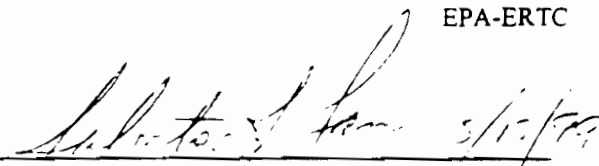
Prepared by  
Roy F. Weston, Inc.

Lackawanna Foundry Site  
Lackawanna, NY

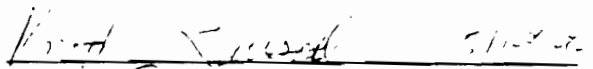
May 1999

EPA Work Assignment No. 3-439  
WESTON Work Order No. 03347-143-001-3439-01  
EPA Contract No. 68-C4-0022

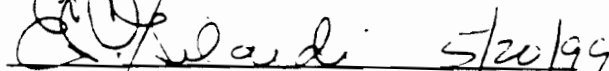
Submitted to  
P. Campagna  
EPA-ERTC

  
\_\_\_\_\_  
S. Eama Date

Task Leader

  
\_\_\_\_\_  
V. Kansal Date

Analytical Section Leader

  
\_\_\_\_\_  
E. Gilardi Date

Program Manager

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REAC  
Accutest

Prepared by:  
G. Karustis

Reviewed by:  
M. Barkley



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Introduction

REAC, in response to WA 3-439, provided analytical support for environmental samples collected from Lackawanna Foundry Site located in Lackawanna, NY as described in the following table. The support also included QA/QC, data review, and preparation of an analytical report containing a summary of the analytical methods, the results, and the QA/QC results.

The samples were treated with procedures consistent with those described in SOP # 1008 and are summarized in the following table:

| COC # | Number of Samples | Sampling Date | Date Received | Matrix | Analysis | Laboratory |
|-------|-------------------|---------------|---------------|--------|----------|------------|
| 06908 | 7                 | 4/14/99       | 4/15/99       | Soil   | BNA      | REAC       |
| 06908 | 6                 | 4/14/99       | 4/15/99       | Soil   | Metals   | REAC       |
| 06909 | 7                 | 4/14/99       | 4/15/99       | Soil   | BNA      | REAC       |
| 06909 | 7                 | 4/14/99       | 4/15/99       | Soil   | Metals   | REAC       |
| 06910 | 2                 | 4/14/99       | 4/15/99       | Soil   | BNA      | REAC       |
| 06910 | 3                 | 4/14/99       | 4/15/99       | Soil   | Metals   | REAC       |
| 06911 | 7                 | 4/14/99       | 4/15/99       | Soil   | BNA      | REAC       |
| 06911 | 6                 | 4/14/99       | 4/15/99       | Soil   | Metals   | REAC       |
| 06912 | 7                 | 4/14/99       | 4/15/99       | Soil   | BNA      | REAC       |
| 06912 | 7                 | 4/14/99       | 4/15/99       | Soil   | Metals   | REAC       |
| 06913 | 1                 | 4/14/99       | 4/15/99       | Soil   | Metals   | REAC       |
| 06914 | 20                | 4/14/99       | 4/15/99       | Soil   | Cyanide  | Accutest   |
| 06915 | 10                | 4/14/99       | 4/15/99       | Soil   | Cyanide  | Accutest   |

Analyses requested on these chains of custody, but not presented in this report, will be given in a separate report.

## Case Narrative

The data in this report have been validated to two significant figures. Any other representation of the data is the responsibility of the user.

### Metals in Soil Package I 171

The percent recoveries for C 16655 MS/MSD exceeded the acceptable QC limits for antimony (0%, 0%), copper (20%, 63%), lead (12%, 38%), selenium (42%, 54%) and arsenic C 16655 MSD (126%). The antimony results for samples C 16650, C 16651, C 16652, C 16653, C 16654, C 16655, C 16657, C 16658, C 16659, C 16660 should be regarded as unusable. The result for antimony in sample C 16656 should be regarded as estimated. The results for copper, lead, selenium and arsenic for samples C 16650, C 16651, C 16652, C 16653, C 16654, C 16655, C 16656, C 16657, C 16658, C 16659 and C 16660 should be regarded as estimated.

The percent recoveries for C 16661 MS/MSD exceeded the acceptable QC limits for antimony (64%, 58%), manganese (10%, 0%), zinc (44%, 26%) and thallium C 16661 MS (69%). The antimony, manganese, zinc and thallium results for samples C 16661, C 16662, C 16663, C 16664, C 16665, C 16666, C 16667 and C 16668 should be regarded as estimated.

The percent recoveries for C 16669 MS/MSD exceeded the acceptable QC limits for antimony (46%, 39%), chromium (219%, 139%), nickel (173%, 136%), selenium (43%, 45%), thallium (20%, 21%) and C 16669 MS for arsenic (145%) and vanadium (126%). The thallium results for samples C 16669, C 16670, C 16671, C 16672, C 16673, C 16674, C 16675, C 16676, C 16677, C 16678 and C 16679 should be regarded as unusable. The results for antimony, chromium, nickel, selenium, arsenic and vanadium for samples C 16669, C 16670, C 16671, C 16672, C 16673, C 16674, C 16675, C 16676, C 16677, C 16678 and C 16679 should be regarded as estimated.

### Cyanide in Soil Package I 182

The data were examined and were found to be acceptable.

### BNA in Soil Package I 190

In the continuing calibration check standard of 4/28/99, the acceptable QC limits were exceeded for bis (2-chloroisopropyl) ether (29%). The data are not affected because this analyte was not detected in the associated samples.

The acceptable QC limits were exceeded for the percent recoveries of one base/neutral surrogate for A 16652 MS. The data are not affected.

### Summary of Abbreviations

|                |   |    |           |    |           |
|----------------|---|----|-----------|----|-----------|
| AA             | Atomic Absorption   |    |           |    |           |
| B              | The analyte was found in the blank  |    |           |    |           |
| BFB            | Bromofluorobenzene  |    |           |    |           |
| C              | Centigrade  |    |           |    |           |
| D              | (Surrogate Table) this value is from a diluted sample and was not calculated<br>(Result Table) this result was obtained from a diluted sample |    |           |    |           |
| Dioxin         | denotes Polychlorinated Dibenzo-p-dioxins and Polychlorinated Dibenzofurans and/or  |    |           |    |           |
| PCDD           | and PCDF  |    |           |    |           |
| CLP            | Contract Laboratory Protocol  |    |           |    |           |
| COC            | Chain of Custody  |    |           |    |           |
| CONC           | Concentration   |    |           |    |           |
| CRDL           | Contract Required Detection Limit   |    |           |    |           |
| CRQL           | Contract Required Quantitation Limit  |    |           |    |           |
| DFTPP          | Decafluorotriphenylphosphine  |    |           |    |           |
| DL             | Detection Limit   |    |           |    |           |
| E              | The value is greater than the highest linear standard and is estimated  |    |           |    |           |
| EMPC           | Estimated maximum possible concentration  |    |           |    |           |
| ICAP           | Inductively Coupled Argon Plasma  |    |           |    |           |
| ISTD           | Internal Standard   |    |           |    |           |
| J              | The value is below the method detection limit and is estimated  |    |           |    |           |
| LCS            | Laboratory Control Sample   |    |           |    |           |
| LCSD           | Laboratory Control Sample Duplicate   |    |           |    |           |
| MDL            | Method Detection Limit  |    |           |    |           |
| MI             | Matrix Interference   |    |           |    |           |
| MS             | Matrix Spike  |    |           |    |           |
| MSD            | Matrix Spike Duplicate  |    |           |    |           |
| MW             | Molecular Weight  |    |           |    |           |
| NA             | either Not Applicable or Not Available  |    |           |    |           |
| NC             | Not Calculated  |    |           |    |           |
| NR             | Not Requested   |    |           |    |           |
| NS             | Not Spiked  |    |           |    |           |
| % D            | Percent Difference  |    |           |    |           |
| % REC          | Percent Recovery  |    |           |    |           |
| PPB            | Parts per billion   |    |           |    |           |
| PPBV           | Parts per billion by volume   |    |           |    |           |
| PQL            | Practical Quantitation Limit  |    |           |    |           |
| QL             | Quantitation Limit  |    |           |    |           |
| RPD            | Relative Percent Difference   |    |           |    |           |
| RSD            | Relative Standard Deviation   |    |           |    |           |
| SIM            | Selected Ion Monitoring   |    |           |    |           |
| TCLP           | Toxic Characteristics Leaching Procedure  |    |           |    |           |
| U              | Denotes not detected  |    |           |    |           |
| W              | Weathered sample; the results should be regarded as estimated   |    |           |    |           |
| m <sup>3</sup> | cubic meter   | kg | kilogram  | μg | microgram |
| L              | liter   | g  | gram      | pg | picogram  |
| mL             | milliliter  | mg | milligram | ng | nanogram  |
| μL             | microliter  |    |           |    |           |
| *              | denotes a value that exceeds the acceptable QC limit  |    |           |    |           |
|                | Abbreviations that are specific to a particular table are explained in footnotes on that table  |    |           |    |           |

Revision 7/23/98

## Analytical Procedure for Metals in Soil

### Sample Preparation

A representative 1-2 g (wet weight) sample, weighed to 0.01 g accuracy, was mixed with 10-mL 1:1 nitric acid, placed in a clean beaker and digested in nitric acid and hydrogen peroxide according to SW-846, Method 3050. The final reflux was either nitric acid or hydrochloric acid depending on the metals to be determined. After digestion, the samples were allowed to cool to room temperature and transferred to 100 mL volumetric flasks and diluted to volume with ASTM Type II water. The samples were analyzed for all metals, except mercury, by USEPA SW-846, Method 7000 (Atomic absorption) or Method 6010 (Inductively Coupled Argon Plasma-ICAP) procedures.

A representative 0.5-0.6 g (wet weight) sample, weighed to 0.01-g accuracy, was prepared and analyzed separately for mercury on a Varian SpectrAA-300 Atomic Absorption Spectrophotometer equipped with a Varian VGA-76 vapor gas analyzer according to SW-846, Method 7471.

A separate sample was used to determine total solids.

A reagent blank and a blank spike sample were carried through the sample preparation procedure for each batch of samples processed. One matrix spike (MS) and one matrix spike duplicate (MSD) were analyzed for each batch or for every ten samples.

### Analysis and Calculations

The instruments were calibrated and operated according to SW-846, Method 7000/7471/6010 and the manufacturers operating instructions. After calibration, initial calibration verification (ICV), initial calibration blank (ICB) and quality control check standards were run to verify proper calibration. The continuing calibration verification (CCV) and continuing calibration blank (CCB) were run after every ten samples to assure proper operation during sample analysis.

The metal concentrations in solution, in micrograms per liter ( $\mu\text{g/L}$ ) were taken from the read-out systems of the Atomic Absorption instruments. The results were converted to milligrams per kilogram ( $\text{mg/kg}$ ) by correcting the reading for the sample weight and percent solids. The ICAP results ( $\text{mg/kg}$ ) were corrected for sample weight prior to instrument read-out; the instrument read-out was then corrected for percent solids.

Final concentrations, based on wet weight are given by:

$$\text{mg metal/kg sample} = [(A \times V) / W] \times DF \times CF$$

where:

- A = Instrument read-out ( $\mu\text{g/L}$ , AA;  $\text{mg/kg}$ , ICAP)
- V = final volume of processed sample (mL, AA: 100, ICAP)
- W = weight of sample (g, AA: 1.00, ICAP)
- DF = Dilution Factor (1.00 for no dilution)
- CF = conversion factor (0.001, AA; 1.00, ICAP)

For samples that required dilution to be within the instrument calibration range, DF is given by:

$$DF = (C+B)/C$$

where:

B = acid blank matrix used for dilution (mL)

C = sample blank aliquot (mL)

Final concentrations, based on dry weight, are given by:

$$\text{mg/kg(dry)} = [\text{mg/kg (wet)} \times 100] / S$$

where

S = percent solids

The results of the analysis are listed in Table 1.1.

### Analytical Procedure for Cyanide in Soil

The subcontract laboratory determined the concentration of cyanide in the samples by analyzing them for cyanide by USEPA Method 9012 M found in SW-846. The results of both analyses are listed in Table 1.2.

## Analytical Procedure for BNA in Soil

### Extraction Procedure

Prior to extraction each sample was spiked with a six component surrogate mixture consisting of nitrobenzene-d<sub>5</sub>, 2-fluorobiphenyl, terphenyl-d<sub>14</sub>, phenol-d<sub>5</sub>, 2-fluorophenol, and 2,4,6-tribromophenol. Thirty grams of sample was mixed with 30 g anhydrous sodium sulfate, and Soxhlet extracted for 16 hours with 300 mL of methylene chloride. The extract was concentrated to 1.0 mL, an internal standard mixture consisting of 1,4-dichlorobenzene-d<sub>4</sub>, naphthalene-d<sub>8</sub>, acenaphthene-d<sub>10</sub>, phenanthrene-d<sub>10</sub>, chrysene-d<sub>12</sub>, and perylene-d<sub>12</sub> was added, and analyzed.

### Analysis Procedure

An HP 6890/5972 Gas Chromatograph/Mass Spectrometer (GC/MS), equipped with a 6890 autosampler and controlled by a PC computer equipped with Enviroquant software was used to analyze the samples.

The instrument conditions were:

|  |  |
|--|--|
| Column   | Restek Rtx-5 (crossbonded SE-54)<br>30 meter x 0.25mm ID, 0.50 µm<br>film thickness  |
| Injection Temperature                          | 280° C   |
| Transfer Temperature                           | 280° C   |
| Source Temperature and<br>Analyzer Temperature | Controlled by thermal transfer of heat from transfer<br>line   |
| Temperature Program                            | 50°C for 0.5 min<br>20° C/min to 295° C<br>hold for 8.5 min<br>25° C/min to 310° C<br>hold for 15 min                      |
| Pulsed Split Injection                         | Split time = 2.00 min @ 8:1 split ratio<br>Pressure Pulse = 16 psi for 0.5 min, then normal                                |
| Injection Volume                               | 1 µL<br><br>Must use 4 mm ID single gooseneck liners packed<br>with 10 mm plug of silanized and conditioned glass<br>wool. |

The GC MS system was calibrated using 5 BNA standard mixtures at 20, 50, 80, 120, and 160 µg/mL. Before analysis each day, the system was tuned with 50-ng decafluorotriphenylphosphine (DFTPP) and passed a continuing calibration check when analyzing a 50-µg/mL standard mixture in which the responses were evaluated by comparison to the average response of the calibration curve.



The BNA results, based on dry weight, are listed in Table 1.3; the tentatively identified compounds are listed in Table 1.4. The concentration of the detected compounds was calculated using the following equation:

$$C_u = \frac{DF \times A_u \times I_{is} \times V_i}{A_{is} \times RF \text{ ( or } RF_{ave} \text{ )} \times V_i \times W \times D}$$

where

- $C_u$  = Concentration of target analyte ( $\mu\text{g/Kg}$ )
- DF = Dilution Factor
- $A_u$  = Area of target analyte
- $I_{is}$  = Mass of specific internal standard (ng)
- $V_i$  = Volume of extract ( $\mu\text{L}$ )
- $A_{is}$  = Area of specific internal standard
- RF = Response Factor (unitless)
- $RF_{ave}$  = average Response Factor
- $V_i$  = Volume of extract injected ( $\mu\text{L}$ )
- W = Weight of sample (g)
- D = Decimal per cent solids

The  $RF_{ave}$  is used when a sample is associated with an initial calibration curve. The RF is used when a sample is associated with a continuing calibration.

Response Factor calculation:

The RF for each specific analyte is quantitated based on the area response from the continuing calibration check as follows:

$$RF = \frac{A_c \times I_{is}}{A_{is} \times I_c}$$

where

- RF = Response factor for a specific analyte
- $A_c$  = Area of the analyte in the standard
- $I_{is}$  = Mass of the specific internal standard
- $A_{is}$  = Area of the specific internal standard
- $I_c$  = Mass of the analyte in the standard

$$RF_{ave} = \frac{RF_1 + \dots + RF_n}{n}$$

and

n = number of Samples

Revision of 7/08/94

**Table 1.1 Results of the Analysis for Metals in Soil  
WA # 3-439 Lackawanna Foundry Site  
Results Based on Dry Weight**

| Client ID | Method Blank    | C16650       |           | C16651           |           | C16652     |           | C16653       |           | C16654     |           |            |           |
|-----------|-----------------|--------------|-----------|------------------|-----------|------------|-----------|--------------|-----------|------------|-----------|------------|-----------|
| Location  | Lab             | 1 West Fence |           | 2 Surface Middle |           | 2D middle  |           | 3 West Fence |           | 4 West     |           |            |           |
| % Solids  | NA              | 64.02        |           | 94.99            |           | 94.41      |           | 79.41        |           | 80.00      |           |            |           |
| Parameter | Analysis Method | Conc mg/kg   | MDL mg/kg | Conc mg/kg       | MDL mg/kg | Conc mg/kg | MDL mg/kg | Conc mg/kg   | MDL mg/kg | Conc mg/kg | MDL mg/kg | Conc mg/kg | MDL mg/kg |
| Aluminum  | ICAP            | U            | 18        | 11000            | 24        | 1300       | 16        | 1200         | 16        | 7800       | 21        | 13000      |           |
| Antimony  | ICAP            | U            | 6.0       | U                | 8.0       | U          | 5.2       | U            | 5.4       | U          | 7.1       | U          |           |
| Arsenic   | AA-Fur          | U            | 0.50      | 14               | 1.5       | 1.6        | 0.50      | 2.5          | 0.48      | 11         | 0.62      | 13         |           |
| Barium    | ICAP            | U            | 1.0       | 160              | 1.3       | 18         | 0.87      | 17           | 0.90      | 140        | 1.2       | 150        |           |
| Beryllium | ICAP            | U            | 0.50      | 0.93             | 0.67      | U          | 0.44      | U            | 0.45      | 0.81       | 0.59      | 2.2        |           |
| Cadmium   | ICAP            | U            | 0.50      | 3.2              | 0.67      | 0.48       | 0.44      | 0.66         | 0.45      | 3.7        | 0.59      | 2.4        |           |
| Calcium   | ICAP            | U            | 50        | 21000            | 67        | 1900       | 44        | 3300         | 45        | 94000      | 59        | 120000     |           |
| Chromium  | ICAP            | U            | 0.50      | 45               | 0.67      | 25         | 0.44      | 15           | 0.45      | 33         | 0.59      | 28         |           |
| Cobalt    | ICAP            | U            | 1.0       | 12               | 1.3       | 2.1        | 0.87      | 1.2          | 0.90      | 7.7        | 1.2       | 6.2        |           |
| Copper    | ICAP            | U            | 1.0       | 74               | 1.3       | 57         | 0.87      | 35           | 0.90      | 75         | 1.2       | 67         |           |
| Iron      | ICAP            | U            | 10        | 43000            | 13        | 24000      | 8.7       | 13000        | 9.0       | 46000      | 12        | 30000      |           |
| Lead      | ICAP            | U            | 4.0       | 300              | 5.3       | 44         | 3.5       | 38           | 3.6       | 200        | 4.7       | 260        |           |
| Magnesium | ICAP            | U            | 50        | 4800             | 67        | 600        | 44        | 560          | 45        | 7200       | 59        | 11000      |           |
| Manganese | ICAP            | U            | 1.0       | 1200             | 1.3       | 250        | 0.87      | 140          | 0.90      | 800        | 1.2       | 1300       |           |
| Mercury   | Cold Vapor      | U            | 0.04      | 0.21             | 0.06      | 0.06       | 0.03      | 0.1          | 0.04      | 0.11       | 0.05      | 0.16       |           |
| Nickel    | ICAP            | U            | 1.0       | 42               | 1.3       | 15         | 0.87      | 9.2          | 0.90      | 28         | 1.2       | 32         |           |
| Potassium | ICAP            | U            | 200       | 1500             | 270       | U          | 170       | U            | 180       | 920        | 240       | 1500       |           |
| Selenium  | AA-Fur          | U            | 0.50      | 1.0              | 0.75      | U          | 0.50      | U            | 0.48      | 0.79       | 0.62      | 0.98       |           |
| Silver    | ICAP            | U            | 0.50      | U                | 0.67      | U          | 0.44      | U            | 0.45      | U          | 0.59      | U          |           |
| Sodium    | ICAP            | U            | 50        | 110              | 67        | U          | 44        | 64           | 45        | 140        | 59        | 520        |           |
| Thallium  | AA-Fur          | U            | 0.50      | U                | 0.75      | U          | 0.50      | U            | 0.48      | U          | 0.62      | U          |           |
| Vanadium  | ICAP            | U            | 2.0       | 27               | 2.7       | 5.2        | 1.7       | 3.7          | 1.8       | 17         | 2.4       | 18         |           |
| Zinc      | ICAP            | U            | 2.0       | 760              | 2.7       | 140        | 1.7       | 79           | 1.8       | 640        | 2.4       | 540        |           |

00009

Table 1.1 (Cont) Results of the Analysis for Metals in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 Results Based on Dry Weight

| Client ID |                 | C16655       |           | C16656          |           | C16657     |           | C16658     |           | C16659     |           | C16660     |           |
|-----------|-----------------|--------------|-----------|-----------------|-----------|------------|-----------|------------|-----------|------------|-----------|------------|-----------|
| Location  |                 | 5 west Fence |           | 6 Near Entrance |           | 7          |           | 8          |           | 9          |           | 10         |           |
| % Solids  |                 | 81.63        |           | 82.59           |           | 75.99      |           | 67.39      |           | 88.49      |           | 99.00      |           |
| Parameter | Analysis Method | Conc mg/kg   | MDL mg/kg | Conc mg/kg      | MDL mg/kg | Conc mg/kg | MDL mg/kg | Conc mg/kg | MDL mg/kg | Conc mg/kg | MDL mg/kg | Conc mg/kg | MDL mg/kg |
| Aluminum  | ICAP            | 12000        | 20        | 4600            | 15        | 5300       | 22        | 3900       | 23        | 5800       | 18        | 2100       | 16        |
| Antimony  | ICAP            | U            | 6.6       | 6.4             | 4.9       | U          | 7.4       | U          | 7.5       | U          | 6.0       | U          | 5.4       |
| Arsenic   | AA-Fur          | 15           | 0.55      | 20              | 0.50      | 6.3        | 0.63      | 6.5        | 0.69      | 16         | 0.51      | 1.3        | 0.49      |
| Barium    | ICAP            | 130          | 1.1       | 270             | 0.81      | 120        | 1.2       | 83         | 1.3       | 170        | 1.0       | 22         | 0.90      |
| Beryllium | ICAP            | 0.85         | 0.55      | U               | 0.41      | U          | 0.62      | U          | 0.63      | U          | 0.50      | U          | 0.45      |
| Cadmium   | ICAP            | 2.4          | 0.55      | 20              | 0.41      | 3.6        | 0.62      | 2.4        | 0.63      | 4.4        | 0.50      | 0.59       | 0.45      |
| Calcium   | ICAP            | 14000        | 55        | 21000           | 41        | 18000      | 62        | 8300       | 63        | 14000      | 50        | 2200       | 45        |
| Chromium  | ICAP            | 29           | 0.55      | 190             | 0.41      | 55         | 0.62      | 36         | 0.63      | 63         | 0.50      | 14         | 0.45      |
| Cobalt    | ICAP            | 14           | 1.1       | 20              | 0.81      | 6.7        | 1.2       | 8.5        | 1.3       | 8.4        | 1.0       | 2.0        | 0.90      |
| Copper    | ICAP            | 200          | 1.1       | 440             | 0.81      | 110        | 1.2       | 81         | 1.3       | 200        | 1.0       | 41         | 0.90      |
| Iron      | ICAP            | 35000        | 11        | 200000          | 41        | 47000      | 12        | 73000      | 13        | 76000      | 10        | 19000      | 9.0       |
| Lead      | ICAP            | 180          | 4.4       | 610             | 3.3       | 2100       | 5.0       | 230        | 5.0       | 340        | 4.0       | 34         | 3.6       |
| Magnesium | ICAP            | 4800         | 55        | 3100            | 41        | 3300       | 62        | 1800       | 63        | 4300       | 50        | 610        | 45        |
| Manganese | ICAP            | 700          | 1.1       | 1800            | 0.81      | 1100       | 1.2       | 790        | 1.3       | 1100       | 1.0       | 180        | 0.90      |
| Mercury   | Cold Vapor      | 0.11         | 0.04      | 1.1             | 0.04      | 0.22       | 0.05      | 0.13       | 0.05      | 0.95       | 0.04      | 0.27       | 0.04      |
| Nickel    | ICAP            | 41           | 1.1       | 160             | 0.81      | 110        | 1.2       | 41         | 1.3       | 160        | 1.0       | 15         | 0.90      |
| Potassium | ICAP            | 1200         | 220       | 630             | 160       | 570        | 250       | 540        | 250       | 390        | 200       | 500        | 180       |
| Selenium  | AA-Fur          | 0.58         | 0.55      | 0.77            | 0.50      | U          | 0.63      | U          | 0.69      | U          | 0.51      | U          | 0.49      |
| Silver    | ICAP            | U            | 0.55      | U               | 2.1       | U          | 0.62      | U          | 0.63      | U          | 0.50      | U          | 0.45      |
| Sodium    | ICAP            | 79           | 55        | 410             | 41        | 230        | 62        | 230        | 63        | 270        | 50        | 1500       | 45        |
| Thallium  | AA-Fur          | U            | 0.55      | U               | 0.50      | U          | 0.63      | U          | 0.69      | U          | 0.51      | U          | 0.49      |
| Vanadium  | ICAP            | 23           | 2.2       | 20              | 1.6       | 20         | 2.5       | 14         | 2.5       | 16         | 2.0       | 4.2        | 1.8       |
| Zinc      | ICAP            | 340          | 2.2       | 1100            | 1.6       | 1000       | 2.5       | 450        | 2.5       | 550        | 2.0       | 42         | 1.8       |

00010

Table 1.1 (Cont) Results of the Analysis for Metals in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 Results Based on Dry Weight

| Client ID |                 | C16661     |           | C16662     |           | C16663     |           | C16664     |           | C16665     |           | C16666     |           |
|-----------|-----------------|------------|-----------|------------|-----------|------------|-----------|------------|-----------|------------|-----------|------------|-----------|
| Location  |                 | 11         |           | 12         |           | 13         |           | 14         |           | 15         |           | Loc 16     |           |
| % Solids  |                 | 85.56      |           | 81.01      |           | 89.86      |           | 78.52      |           | 94.98      |           | 9          |           |
| Parameter | Analysis Method | Conc mg/kg | MDL mg/kg | Conc mg/kg | MDL mg/kg | Conc mg/kg | MDL mg/kg | Conc mg/kg | MDL mg/kg | Conc mg/kg | MDL mg/kg | Conc mg/kg | MDL mg/kg |
| Aluminum  | ICAP            | 740        | 18        | 4200       | 12        | 4600       | 19        | 2300       | 15        | 2200       | 17        | 1300       |           |
| Antimony  | ICAP            | U          | 5.9       | U          | 4.2       | 13         | 6.3       | U          | 4.8       | U          | 5.6       | U          |           |
| Arsenic   | AA-Fur          | 1.8        | 0.40      | 7.0        | 0.35      | 16         | 0.39      | 4.4        | 0.41      | 2.7        | 0.48      | 1.5        |           |
| Barium    | ICAP            | 14         | 0.98      | 84         | 0.69      | 100        | 1.0       | 39         | 0.81      | 50         | 0.93      | 15         |           |
| Beryllium | ICAP            | U          | 0.49      | U          | 0.35      | U          | 0.52      | U          | 0.40      | U          | 0.47      | U          |           |
| Cadmium   | ICAP            | 4.2        | 0.49      | 2.0        | 0.35      | 3.4        | 0.52      | 0.92       | 0.40      | 0.48       | 0.47      | U          |           |
| Calcium   | ICAP            | 1100       | 49        | 21000      | 35        | 1200       | 52        | 24000      | 40        | 1600       | 47        | 1200       |           |
| Chromium  | ICAP            | 13         | 0.49      | 100        | 0.35      | 79         | 0.52      | 86         | 0.40      | 14         | 0.47      | 6.1        |           |
| Cobalt    | ICAP            | 1.7        | 0.98      | 7.0        | 0.69      | 11         | 1.0       | 3.6        | 0.81      | 0.98       | 0.93      | 1.4        |           |
| Copper    | ICAP            | 33         | 0.98      | 270        | 0.69      | 370        | 1.0       | 93         | 0.81      | 100        | 0.93      | 44         |           |
| Iron      | ICAP            | 22000      | 9.8       | 56000      | 6.9       | 140000     | 53        | 39000      | 8.1       | 11000      | 9.3       | 16000      |           |
| Lead      | ICAP            | 48         | 3.9       | 150        | 2.8       | 540        | 4.2       | 110        | 3.2       | 58         | 3.7       | 58         |           |
| Magnesium | ICAP            | 240        | 49        | 2400       | 35        | 480        | 52        | 3400       | 40        | 460        | 47        | 410        |           |
| Manganese | ICAP            | 160        | 0.98      | 810        | 0.69      | 770        | 1.0       | 2500       | 0.81      | 120        | 0.93      | 210        |           |
| Mercury   | Cold Vapor      | 0.04       | 0.04      | 0.17       | 0.03      | 0.07       | 0.03      | 0.12       | 0.03      | U          | 0.04      | U          |           |
| Nickel    | ICAP            | 13         | 0.98      | 75         | 0.69      | 72         | 1.0       | 25         | 0.81      | 31         | 0.93      | 9.5        |           |
| Potassium | ICAP            | U          | 200       | 430        | 140       | 500        | 210       | 260        | 160       | U          | 190       | U          |           |
| Selenium  | AA-Fur          | U          | 0.40      | U          | 0.35      | 0.39       | 0.39      | U          | 0.41      | U          | 0.48      | U          |           |
| Silver    | ICAP            | U          | 0.49      | U          | 0.35      | U          | 0.52      | U          | 0.40      | U          | 0.47      | U          |           |
| Sodium    | ICAP            | U          | 49        | 210        | 35        | 200        | 52        | 92         | 40        | 72         | 47        | U          |           |
| Thallium  | AA-Fur          | U          | 0.40      | U          | 0.35      | U          | 0.39      | U          | 0.41      | U          | 0.48      | U          |           |
| Vanadium  | ICAP            | 2.7        | 2.0       | 14         | 1.4       | 16         | 2.1       | 43         | 1.6       | 5.4        | 1.9       | 3.1        |           |
| Zinc      | ICAP            | 81         | 2.0       | 280        | 1.4       | 660        | 2.1       | 190        | 1.6       | 110        | 1.9       | 98         |           |

00011

**Table 1.1 (Cont) Results of the Analysis for Metals in Soil  
WA # 3-439 Lackawanna Foundry Site  
Results Based on Dry Weight**

| Client ID |                 | C16667      |           | C16668      |           | C16669      |           | C16670      |           | C16671      |           | C16672      |           |
|-----------|-----------------|-------------|-----------|-------------|-----------|-------------|-----------|-------------|-----------|-------------|-----------|-------------|-----------|
| Location  |                 | Location 17 |           | Location 18 |           | Location 19 |           | Location 20 |           | Location 21 |           | Location 22 |           |
| % Solids  |                 | 90.83       |           | 92.93       |           | 95.49       |           | 88.29       |           | 51.93       |           | 78.77       |           |
| Parameter | Analysis Method | Conc mg/kg  | MDL mg/kg | Conc mg/kg  | MDL mg/kg | Conc mg/kg  | MDL mg/kg | Conc mg/kg  | MDL mg/kg | Conc mg/kg  | MDL mg/kg | Conc mg/kg  | MDL mg/kg |
| Aluminum  | ICAP            | 5900        | 17        | 790         | 17        | 1400        | 15        | 2400        | 17        | 12000       | 23        | 15000       | 20        |
| Antimony  | ICAP            | 6.5         | 5.8       | U           | 5.6       | U           | 5.0       | U           | 5.7       | U           | 7.8       | U           | 6.7       |
| Arsenic   | AA-Fur          | 7.1         | 0.95      | 4.4         | 0.45      | 6.7         | 0.51      | 6.5         | 0.46      | 20          | 0.54      | 77          | 0.58      |
| Barium    | ICAP            | 180         | 0.97      | 9.0         | 0.94      | 29          | 0.83      | 130         | 0.94      | 200         | 1.3       | 170         | 1.1       |
| Beryllium | ICAP            | U           | 0.48      | U           | 0.47      | U           | 0.42      | U           | 0.47      | 0.96        | 0.65      | 1.2         | 0.56      |
| Cadmium   | ICAP            | U           | 0.48      | U           | 0.47      | 0.59        | 0.42      | 2.3         | 0.47      | 13          | 0.65      | 2.3         | 0.56      |
| Calcium   | ICAP            | 130000      | 48        | 1200        | 47        | 850         | 42        | 3800        | 47        | 23000       | 65        | 34000       | 56        |
| Chromium  | ICAP            | 33          | 0.48      | 8.9         | 0.47      | 97          | 0.42      | 130         | 0.47      | 95          | 0.65      | 32          | 0.56      |
| Cobalt    | ICAP            | 47          | 0.97      | U           | 0.94      | 4.4         | 0.83      | 12          | 0.94      | 15          | 1.3       | 11          | 1.1       |
| Copper    | ICAP            | 240         | 0.97      | 28          | 0.94      | 300         | 0.83      | 720         | 0.94      | 270         | 1.3       | 47          | 1.1       |
| Iron      | ICAP            | 25000       | 9.7       | 5800        | 9.4       | 58000       | 8.3       | 93000       | 9.4       | 110000      | 13        | 32000       | 11        |
| Lead      | ICAP            | 300         | 3.9       | 21          | 3.7       | 140         | 3.3       | 280         | 3.8       | 660         | 5.2       | 180         | 4.5       |
| Magnesium | ICAP            | 19000       | 48        | 340         | 47        | 410         | 42        | 1000        | 47        | 4700        | 65        | 5900        | 56        |
| Manganese | ICAP            | 840         | 0.97      | 75          | 0.94      | 760         | 0.83      | 690         | 0.94      | 1600        | 1.3       | 1200        | 1.1       |
| Mercury   | Cold Vapor      | 0.06        | 0.04      | 0.04        | 0.04      | U           | 0.03      | 1.3         | 0.04      | 0.77        | 0.05      | 0.73        | 0.04      |
| Nickel    | ICAP            | 38          | 0.97      | 2.7         | 0.94      | 57          | 0.83      | 150         | 0.94      | 100         | 1.3       | 36          | 1.1       |
| Potassium | ICAP            | 470         | 190       | U           | 190       | U           | 170       | 280         | 190       | 900         | 260       | 1500        | 220       |
| Selenium  | AA-Fur          | U           | 0.47      | U           | 0.45      | U           | 0.51      | U           | 0.46      | 1.0         | 0.54      | U           | 0.58      |
| Silver    | ICAP            | U           | 0.48      | U           | 0.47      | U           | 0.42      | U           | 0.47      | U           | 0.65      | U           | 0.56      |
| Sodium    | ICAP            | 330         | 48        | U           | 47        | 68          | 42        | 68          | 47        | 360         | 65        | 120         | 56        |
| Thallium  | AA-Fur          | U           | 0.47      | U           | 0.45      | U           | 0.51      | U           | 0.46      | U           | 0.54      | U           | 0.58      |
| Vanadium  | ICAP            | 11          | 1.9       | 2.6         | 1.9       | 14          | 1.7       | 21          | 1.9       | 36          | 2.6       | 27          | 2.2       |
| Zinc      | ICAP            | 1300        | 1.9       | 55          | 1.9       | 310         | 1.7       | 290         | 1.9       | 1000        | 2.6       | 1100        | 2.2       |

00012

Table 1.1 (Cont) Results of the Analysis for Metals in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 Results Based on Dry Weight

| Client ID |                 | C16673      |           | C16674      |           | C16675      |           | C16676        |           | C16677      |           | C16678      |           |
|-----------|-----------------|-------------|-----------|-------------|-----------|-------------|-----------|---------------|-----------|-------------|-----------|-------------|-----------|
| Location  |                 | Location 23 |           | Location 24 |           | Location 25 |           | Location 25-D |           | Location 26 |           | Location 27 |           |
| % Solids  |                 | 81.54       |           | 77.01       |           | 80.54       |           | 78.92         |           | 74.19       |           | 63.00       |           |
| Parameter | Analysis Method | Conc mg/kg  | MDL mg/kg | Conc mg/kg  | MDL mg/kg | Conc mg/kg  | MDL mg/kg | Conc mg/kg    | MDL mg/kg | Conc mg/kg  | MDL mg/kg | Conc mg/kg  | MDL mg/kg |
| Aluminum  | ICAP            | 9300        | 19        | 7300        | 19        | 12000       | 18        | 17000         | 21        | 14000       | 23        | 14000       | 23        |
| Antimony  | ICAP            | U           | 6.5       | U           | 6.5       | U           | 6.1       | U             | 7.0       | U           | 7.6       | U           | U         |
| Arsenic   | AA-Fur          | 8.4         | 0.46      | 5           | 0.98      | 21          | 0.49      | 25            | 0.50      | 66          | 0.53      | 48          | U         |
| Barium    | ICAP            | 94          | 1.1       | 64          | 1.1       | 130         | 1.0       | 140           | 1.2       | 250         | 1.3       | 280         | U         |
| Beryllium | ICAP            | 0.66        | 0.54      | U           | 0.54      | 0.73        | 0.50      | 0.97          | 0.59      | 1.3         | 0.64      | 0.67        | U         |
| Cadmium   | ICAP            | 1.5         | 0.54      | 1.2         | 0.54      | 2.3         | 0.50      | U             | 0.59      | 5.2         | 0.64      | 1.3         | U         |
| Calcium   | ICAP            | 12000       | 54        | 10000       | 54        | 39000       | 50        | 11000         | 59        | 53000       | 64        | 120000      | U         |
| Chromium  | ICAP            | 24          | 0.54      | 18          | 0.54      | 31          | 0.50      | 25            | 0.59      | 51          | 0.64      | 23          | U         |
| Cobalt    | ICAP            | 8.0         | 1.1       | 4.8         | 1.1       | 9.7         | 1.0       | 15            | 1.2       | 9.7         | 1.3       | 7.3         | U         |
| Copper    | ICAP            | 52          | 1.1       | 49          | 1.1       | 65          | 1.0       | 27            | 1.2       | 110         | 1.3       | 34          | U         |
| Iron      | ICAP            | 30000       | 11        | 25000       | 11        | 29000       | 10        | 32000         | 12        | 45000       | 13        | 20000       | U         |
| Lead      | ICAP            | 160         | 4.3       | 130         | 4.3       | 230         | 4.0       | 40            | 4.7       | 740         | 5.1       | 100         | U         |
| Magnesium | ICAP            | 3600        | 54        | 3500        | 54        | 5400        | 50        | 5400          | 59        | 6300        | 64        | 7300        | U         |
| Manganese | ICAP            | 560         | 1.1       | 390         | 1.1       | 710         | 1.0       | 610           | 1.2       | 1500        | 1.3       | 570         | U         |
| Mercury   | Cold Vapor      | 0.13        | 0.04      | 0.13        | 0.04      | 0.17        | 0.04      | 0.08          | 0.04      | 1.6         | 0.04      | 0.17        | U         |
| Nickel    | ICAP            | 28          | 1.1       | 18          | 1.1       | 38          | 1.0       | 41            | 1.2       | 39          | 1.3       | 22          | U         |
| Potassium | ICAP            | 1100        | 220       | 1100        | 220       | 1400        | 200       | 1500          | 230       | 1600        | 250       | 1400        | U         |
| Selenium  | AA-Fur          | 0.57        | 0.46      | U           | 0.49      | U           | 0.49      | U             | 0.50      | 1.6         | 0.53      | U           | U         |
| Silver    | ICAP            | U           | 0.54      | U           | 0.54      | U           | 0.50      | U             | 0.59      | U           | 0.64      | U           | U         |
| Sodium    | ICAP            | 130         | 54        | 250         | 54        | 170         | 50        | U             | 59        | 190         | 64        | 190         | U         |
| Thallium  | AA-Fur          | U           | 0.46      | U           | 0.49      | U           | 0.49      | U             | 0.50      | U           | 0.53      | U           | U         |
| Vanadium  | ICAP            | 19          | 2.2       | 14          | 2.2       | 23          | 2.0       | 27            | 2.3       | 29          | 2.5       | 20          | U         |
| Zinc      | ICAP            | 660         | 2.2       | 350         | 2.2       | 760         | 2.0       | 150           | 2.3       | 1600        | 2.5       | 330         | U         |

00013

Table 1.1 (Cont) Results of the Analysis for Metals in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 Results Based on Dry Weight

| Client ID | C16679          |            |           |
|-----------|-----------------|------------|-----------|
| Location  | Location 27     |            |           |
| % Solids  | 83.78           |            |           |
| Parameter | Analysis Method | Conc mg/kg | MDL mg/kg |
| Aluminum  | ICAP            | 2900       | 17        |
| Antimony  | ICAP            | U          | 5.6       |
| Arsenic   | AA-Fur          | 19         | 0.34      |
| Barium    | ICAP            | 59         | 0.94      |
| Beryllium | ICAP            | U          | 0.47      |
| Cadmium   | ICAP            | 0.84       | 0.47      |
| Calcium   | ICAP            | 15000      | 47        |
| Chromium  | ICAP            | 95         | 0.47      |
| Cobalt    | ICAP            | 8.7        | 0.94      |
| Copper    | ICAP            | 280        | 0.94      |
| Iron      | ICAP            | 78000      | 9.4       |
| Lead      | ICAP            | 140        | 3.8       |
| Magnesium | ICAP            | 2500       | 47        |
| Manganese | ICAP            | 770        | 0.94      |
| Mercury   | Cold Vapor      | 0.09       | 0.04      |
| Nickel    | ICAP            | 89         | 0.94      |
| Potassium | ICAP            | 320        | 190       |
| Selenium  | AA-Fur          | U          | 0.34      |
| Silver    | ICAP            | U          | 0.47      |
| Sodium    | ICAP            | 120        | 47        |
| Thallium  | AA-Fur          | U          | 0.34      |
| Vanadium  | ICAP            | 15         | 1.9       |
| Zinc      | ICAP            | 250        | 1.9       |

03014

Table 1.2 Results of the Analysis for Cyanide in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 Results Based on Dry Weight

| Sample ID    | Location    | % Solids | Cyanide       |              |
|--------------|-------------|----------|---------------|--------------|
|              |             |          | Conc<br>mg/kg | MDL<br>mg/kg |
| Method Blank | -           | 100      | U             | 1.0          |
| Method Blank | -           | 100      | U             | 1.0          |
| D16650       | 1           | 56.3     | U             | 1.8          |
| D16651       | 2           | 94.1     | U             | 1.1          |
| D16652       | 2-D         | 95.2     | U             | 1.0          |
| D16653       | 3           | 81.4     | U             | 1.2          |
| D16654       | 4           | 86.6     | U             | 1.2          |
| D16655       | 5           | 84.0     | U             | 1.2          |
| D16656       | 6           | 86.6     | U             | 1.2          |
| D16657       | 7           | 79.9     | U             | 1.2          |
| D16658       | 8           | 79.1     | U             | 1.3          |
| D16659       | 9           | 89.7     | U             | 1.1          |
| D16660       | 10          | 99.2     | U             | 1.0          |
| D16661       | 11          | 85.8     | U             | 1.2          |
| D16662       | 12          | 77.3     | U             | 1.3          |
| D16663       | 13          | 92.0     | U             | 1.1          |
| D16664       | 14          | 75.9     | U             | 1.3          |
| D16665       | 15          | 94.7     | U             | 1.0          |
| D16666       | 16          | 90.3     | U             | 1.1          |
| D16667       | 17          | 91.8     | U             | 1.1          |
| D16668       | 18          | 93.1     | U             | 1.1          |
| D16669       | 19          | 93.2     | U             | 1.1          |
| D16670       | Location 20 | 90.8     | U             | 1.1          |
| D16671       | 21          | 43.8     | U             | 2.3          |
| D16672       | 22          | 81.8     | U             | 1.2          |
| D16673       | 23          | 84.0     | U             | 1.2          |
| D16674       | 24          | 81.4     | U             | 1.2          |
| D16675       | 25          | 87.5     | U             | 1.1          |
| D16676       | 25-D        | 78.2     | U             | 1.3          |
| D16677       | 26          | 76.7     | U             | 1.3          |
| D16678       | 26-D        | 75.2     | U             | 1.3          |
| D16679       | 27          | 84.7     | U             | 1.2          |

00015



Table 1.3 Results of the Analysis for BNA in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 (Results are Based on Dry Weight)

| Sample No.                  | SBLK041599     |              | A16650         |              | A16651           |              | A16652         |              | A16653         |              |
|-----------------------------|----------------|--------------|----------------|--------------|------------------|--------------|----------------|--------------|----------------|--------------|
| Sample Location             | Lab Blank      |              | 1 West Fence   |              | 2 Surface Middle |              | 2D Middle      |              | 3 West Fence   |              |
| GC/MS File Name             | LFS002         |              | LFS003         |              | LFS004           |              | LFS005         |              | LFS008         |              |
| Dilution Factor             | 1              |              | 10             |              | 1                |              | 5              |              | 10             |              |
| % Solid                     | 100            |              | 57             |              | 95               |              | 91             |              | 73             |              |
| Compound Name               | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg   | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg |
| Phenol                      | U              | 330          | U              | 5800         | 16000            | 350          | 23000          | 1800         | U              | 4600         |
| bis(-2-Chloroethyl)Ether    | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| 2-Chlorophenol              | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| 1,3-Dichlorobenzene         | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| 1,4-Dichlorobenzene         | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| Benzyl alcohol              | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| 1,2-Dichlorobenzene         | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| 2-Methylphenol              | U              | 330          | U              | 5800         | 370              | 350          | 410 J          | 1800         | U              | 4600         |
| bis(2-Chloroisopropyl)ether | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| 4-Methylphenol              | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| N-Nitroso-Di-n-propylamine  | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| Hexachloroethane            | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| Nitrobenzene                | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| Isophorone                  | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| 2-Nitrophenol               | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| 2,4-Dimethylphenol          | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| bis(2-Chloroethoxy)methane  | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| 2,4-Dichlorophenol          | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| 1,2,4-Trichlorobenzene      | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| Naphthalene                 | U              | 330          | U              | 5800         | 970              | 350          | 960 J          | 1800         | U              | 4600         |
| 4-Chloroaniline             | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| Hexachlorobutadiene         | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| 4-Chloro-3-methylphenol     | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| 2-Methylnaphthalene         | U              | 330          | U              | 5800         | 360              | 350          | 640 J          | 1800         | U              | 4600         |
| Hexachlorocyclopentadiene   | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| 2,4,6-Trichlorophenol       | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| 2,4,5-Trichlorophenol       | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| 2-Chloronaphthalene         | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| 2-Nitroaniline              | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| Dimethylphthalate           | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| Acenaphthylene              | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| 2,6-Dinitrotoluene          | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| 3-Nitroaniline              | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| Acenaphthene                | U              | 330          | U              | 5800         | 95 J             | 350          | U              | 1800         | U              | 4600         |
| 2,4-Dinitrophenol           | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| 4-Nitrophenol               | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| Dibenzofuran                | U              | 330          | U              | 5800         | 100 J            | 350          | U              | 1800         | U              | 4600         |
| 2,4-Dinitrotoluene          | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| Diethylphthalate            | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| 4-Chlorophenyl-phenylether  | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| Fluorene                    | U              | 330          | U              | 5800         | 87 J             | 350          | U              | 1800         | U              | 4600         |
| 4-Nitroaniline              | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| 4,6-Dinitro-2-methylphenol  | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| N-Nitrosodiphenylamine      | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| 4-Bromophenyl-phenylether   | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| Hexachlorobenzene           | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| Pentachlorophenol           | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| Phenanthrene                | U              | 330          | U              | 5800         | 1400             | 350          | U              | 1800         | U              | 4600         |
| Anthracene                  | U              | 330          | U              | 5800         | 220 J            | 350          | U              | 1800         | U              | 4600         |
| Carbazole                   | U              | 330          | U              | 5800         | 190 J            | 350          | U              | 1800         | U              | 4600         |
| Di-n-butylphthalate         | U              | 330          | U              | 5800         | 150 J            | 350          | U              | 1800         | U              | 4600         |
| Fluoranthene                | U              | 330          | U              | 5800         | 2000             | 350          | 380 J          | 1800         | 1000 J         | 4600         |
| Pyrene                      | U              | 330          | U              | 5800         | 1500             | 350          | U              | 1800         | U              | 4600         |
| Butylbenzylphthalate        | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| Benzo(a)anthracene          | U              | 330          | U              | 5800         | 800              | 350          | U              | 1800         | U              | 4600         |
| 3,3'-Dichlorobenzidine      | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| Chrysene                    | U              | 330          | U              | 5800         | 920              | 350          | U              | 1800         | U              | 4600         |
| Bis(2-Ethylhexyl)phthalate  | U              | 330          | U              | 5800         | 640              | 350          | 1300 J         | 1800         | 1000 J         | 4600         |
| Di-n-octylphthalate         | U              | 330          | U              | 5800         | U                | 350          | U              | 1800         | U              | 4600         |
| Benzo(b)fluoranthene        | U              | 330          | U              | 5800         | 1100             | 350          | U              | 1800         | U              | 4600         |
| Benzo(k)fluoranthene        | U              | 330          | U              | 5800         | 880              | 350          | U              | 1800         | U              | 4600         |
| Benzo(a)pyrene              | U              | 330          | U              | 5800         | 1100             | 350          | U              | 1800         | U              | 4600         |
| Indeno(1,2,3-cd)pyrene      | U              | 330          | U              | 5800         | 750              | 350          | U              | 1800         | U              | 4600         |
| Dibenzo(a,h)anthracene      | U              | 330          | U              | 5800         | 300 J            | 350          | U              | 1800         | U              | 4600         |
| Benzo(g,h,i)perylene        | U              | 330          | U              | 5800         | 890              | 350          | U              | 1800         | U              | 4600         |

Table 1.3 (cont.) Results of the Analysis for BNA in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 (Results are Based on Dry Weight)

| Sample No                   | A16654         |              | A16655         |              | A16656          |              |
|-----------------------------|----------------|--------------|----------------|--------------|-----------------|--------------|
| Sample Location             | 4 West Fence   |              | 5 West Fence   |              | 6 Near Entrance |              |
| GC/MS File Name             | LFS009         |              | LFS010         |              | LFS012          |              |
| Dilution Factor             | 10             |              | 1              |              | 10              |              |
| % Solid                     | 82             |              | 81             |              | 85              |              |
| Compound Name               | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg  | MDL<br>µg/kg |
| Phenol                      | U              | 4100         | U              | 410          | U               | 3900         |
| bis(-2-Chloroethyl)Ether    | U              | 4100         | U              | 410          | U               | 3900         |
| 2-Chlorophenol              | U              | 4100         | U              | 410          | U               | 3900         |
| 1,3-Dichlorobenzene         | U              | 4100         | U              | 410          | U               | 3900         |
| 1,4-Dichlorobenzene         | U              | 4100         | U              | 410          | U               | 3900         |
| Benzyl alcohol              | U              | 4100         | U              | 410          | U               | 3900         |
| 1,2-Dichlorobenzene         | U              | 4100         | U              | 410          | U               | 3900         |
| 2-Methylphenol              | U              | 4100         | U              | 410          | U               | 3900         |
| bis(2-Chloroisopropyl)ether | U              | 4100         | U              | 410          | U               | 3900         |
| 4-Methylphenol              | U              | 4100         | U              | 410          | U               | 3900         |
| N-Nitroso-Di-n-propylamine  | U              | 4100         | U              | 410          | U               | 3900         |
| Hexachloroethane            | U              | 4100         | U              | 410          | U               | 3900         |
| Nitrobenzene                | U              | 4100         | U              | 410          | U               | 3900         |
| Isophorone                  | U              | 4100         | U              | 410          | U               | 3900         |
| 2-Nitrophenol               | U              | 4100         | U              | 410          | U               | 3900         |
| 2,4-Dimethylphenol          | U              | 4100         | U              | 410          | U               | 3900         |
| bis(2-Chloroethoxy)methane  | U              | 4100         | U              | 410          | U               | 3900         |
| 2,4-Dichlorophenol          | U              | 4100         | U              | 410          | U               | 3900         |
| 1,2,4-Trichlorobenzene      | U              | 4100         | U              | 410          | U               | 3900         |
| Naphthalene                 | U              | 4100         | 310 J          | 410          | U               | 3900         |
| 4-Chloroaniline             | U              | 4100         | U              | 410          | U               | 3900         |
| Hexachlorobutadiene         | U              | 4100         | U              | 410          | U               | 3900         |
| 4-Chloro-3-methylphenol     | U              | 4100         | U              | 410          | U               | 3900         |
| 2-Methylnaphthalene         | U              | 4100         | 270 J          | 410          | U               | 3900         |
| Hexachlorocyclopentadiene   | U              | 4100         | U              | 410          | U               | 3900         |
| 2,4,6-Trichlorophenol       | U              | 4100         | U              | 410          | U               | 3900         |
| 2,4,5-Trichlorophenol       | U              | 4100         | U              | 410          | U               | 3900         |
| 2-Chloronaphthalene         | U              | 4100         | U              | 410          | U               | 3900         |
| 2-Nitroaniline              | U              | 4100         | U              | 410          | U               | 3900         |
| Dimethylphthalate           | U              | 4100         | U              | 410          | U               | 3900         |
| Acenaphthylene              | U              | 4100         | 200 J          | 410          | U               | 3900         |
| 2,6-Dinitrotoluene          | U              | 4100         | U              | 410          | U               | 3900         |
| 3-Nitroaniline              | U              | 4100         | U              | 410          | U               | 3900         |
| Acenaphthene                | U              | 4100         | 740            | 410          | U               | 3900         |
| 2,4-Dinitrophenol           | U              | 4100         | U              | 410          | U               | 3900         |
| 4-Nitrophenol               | U              | 4100         | U              | 410          | U               | 3900         |
| Dibenzofuran                | U              | 4100         | 490            | 410          | U               | 3900         |
| 2,4-Dinitrotoluene          | U              | 4100         | U              | 410          | U               | 3900         |
| Diethylphthalate            | U              | 4100         | U              | 410          | U               | 3900         |
| 4-Chlorophenyl-phenylether  | U              | 4100         | U              | 410          | U               | 3900         |
| Fluorene                    | U              | 4100         | 880            | 410          | U               | 3900         |
| 4-Nitroaniline              | U              | 4100         | U              | 410          | U               | 3900         |
| 4,6-Dinitro-2-methylphenol  | U              | 4100         | U              | 410          | U               | 3900         |
| N-Nitrosodiphenylamine      | U              | 4100         | U              | 410          | U               | 3900         |
| 4-Bromophenyl-phenylether   | U              | 4100         | U              | 410          | U               | 3900         |
| Hexachlorobenzene           | U              | 4100         | U              | 410          | U               | 3900         |
| Pentachlorophenol           | U              | 4100         | U              | 410          | U               | 3900         |
| Phenanthrene                | U              | 4100         | 7900           | 410          | U               | 3900         |
| Anthracene                  | U              | 4100         | 1800           | 410          | U               | 3900         |
| Carbazole                   | U              | 4100         | 1200           | 410          | U               | 3900         |
| Di-n-butylphthalate         | U              | 4100         | U              | 410          | U               | 3900         |
| Fluoranthene                | 2300 J         | 4100         | 9900           | 410          | 2000 J          | 3900         |
| Pyrene                      | 1900 J         | 4100         | 7800           | 410          | 1700 J          | 3900         |
| Butylbenzylphthalate        | U              | 4100         | U              | 410          | U               | 3900         |
| Benzo(a)anthracene          | 1100 J         | 4100         | 5000           | 410          | 1000 J          | 3900         |
| 3,3'-Dichlorobenzidine      | U              | 4100         | U              | 410          | U               | 3900         |
| Chrysene                    | 1400 J         | 4100         | 5300           | 410          | 1100 J          | 3900         |
| Bis(2-Ethylhexyl)phthalate  | 1000 J         | 4100         | 260 J          | 410          | 810 J           | 3900         |
| Di-n-octylphthalate         | U              | 4100         | U              | 410          | U               | 3900         |
| Benzo(b)fluoranthene        | 1700 J         | 4100         | 5800           | 410          | 1400 J          | 3900         |
| Benzo(k)fluoranthene        | 1400 J         | 4100         | 4500           | 410          | 1200 J          | 3900         |
| Benzo(a)pyrene              | 1700 J         | 4100         | 5800           | 410          | 1300 J          | 3900         |
| Indeno(1,2,3-cd)pyrene      | 1300 J         | 4100         | 3400           | 410          | 990 J           | 3900         |
| Dibenzo(a,h)anthracene      | U              | 4100         | 1300           | 410          | U               | 3900         |
| Benzo(g,h,i)perylene        | 1600 J         | 4100         | 3700           | 410          | 1100 J          | 3900         |

Table 1.3 (cont.) Results of the Analysis for BNA in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 (Results are Based on Dry Weight)

| Sample No.                  | SBLK041999     | A16657       | -A16658        | A16659       | A16660         |              |                |              |                |              |      |      |      |      |
|-----------------------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|------|------|------|------|
| Sample Location             | Lab Blank      | 7            | 8              | 9            | 10             |              |                |              |                |              |      |      |      |      |
| GC/MS File Name             | LFS023         | LFS024       | LFS025         | LFS026       | LFS027         |              |                |              |                |              |      |      |      |      |
| Dilution Factor             | 1              | 10           | 10             | 10           | 5              |              |                |              |                |              |      |      |      |      |
| % Solid                     | 100            | 76           | 87             | 91           | 99             |              |                |              |                |              |      |      |      |      |
| Compound Name               | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg |      |      |      |      |
| Phenol                      | U              | 330          | 1900           | J            | 4400           | 1500         | J              | 3800         | 4900           | 3700         | 1800 | 1700 |      |      |
| bis(-2-Chloroethyl)Ether    | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| 2-Chlorophenol              | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| 1,3-Dichlorobenzene         | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| 1,4-Dichlorobenzene         | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| Benzyl alcohol              | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| 1,2-Dichlorobenzene         | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| 2-Methylphenol              | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | 570  | J    | 1700 |      |
| bis(2-Chloroisopropyl)ether | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| 4-Methylphenol              | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| N-Nitroso-Di-n-propylamine  | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| Hexachloroethane            | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| Nitrobenzene                | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| Isophorone                  | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| 2-Nitrophenol               | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| 2,4-Dimethylphenol          | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| bis(2-Chloroethoxy)methane  | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| 2,4-Dichlorophenol          | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| 1,2,4-Trichlorobenzene      | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| Naphthalene                 | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | 3100 | 1700 |      |      |
| 4-Chloroaniline             | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| Hexachlorobutadiene         | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| 4-Chloro-3-methylphenol     | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| 2-Methylnaphthalene         | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | 1500 | J    | 1700 |      |
| Hexachlorocyclopentadiene   | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| 2,4,6-Trichlorophenol       | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| 2,4,5-Trichlorophenol       | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| 2-Chloronaphthalene         | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| 2-Nitroaniline              | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| Dimethylphthalate           | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| Acenaphthylene              | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| 2,6-Dinitrotoluene          | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| 3-Nitroaniline              | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| Acenaphthene                | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| 2,4-Dinitrophenol           | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| 4-Nitrophenol               | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| Dibenzofuran                | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| 2,4-Dinitrotoluene          | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| Diethylphthalate            | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| 4-Chlorophenyl-phenylether  | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| Fluorene                    | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| 4-Nitroaniline              | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| 4,6-Dinitro-2-methylphenol  | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| N-Nitrosodiphenylamine      | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| 4-Bromophenyl-phenylether   | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| Hexachlorobenzene           | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| Pentachlorophenol           | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| Phenanthrene                | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | 680  | J    | 1700 |      |
| Anthracene                  | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| Carbazole                   | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| Di-n-butylphthalate         | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| Fluoranthene                | U              | 330          | 2400           | J            | 4400           | 1300         | J              | 3800         | 1000           | J            | 3700 | U    | 1700 |      |
| Pyrene                      | U              | 330          | 2200           | J            | 4400           | 1200         | J              | 3800         | 880            | J            | 3700 | U    | 1700 |      |
| Butylbenzylphthalate        | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | 6000           | 3700         | U    | 1700 |      |      |
| Benzo(a)anthracene          | U              | 330          | 1500           | J            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| 3,3'-Dichlorobenzidine      | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| Chrysene                    | U              | 330          | 1900           | J            | 4400           | 1000         | J              | 3800         | U              | 3700         | U    | 1700 |      |      |
| Bis(2-Ethylhexyl)phthalate  | U              | 330          | U              | U            | 4400           | 8700         | U              | 3800         | 1200           | J            | 3700 | 1300 | J    | 1700 |
| Di-n-octylphthalate         | U              | 330          | U              | U            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| Benzo(b)fluoranthene        | U              | 330          | 3100           | J            | 4400           | 1600         | J              | 3800         | 940            | J            | 3700 | U    | 1700 |      |
| Benzo(k)fluoranthene        | U              | 330          | 2400           | J            | 4400           | 1400         | J              | 3800         | 830            | J            | 3700 | U    | 1700 |      |
| Benzo(a)pyrene              | U              | 330          | 3300           | J            | 4400           | 1400         | J              | 3800         | 920            | J            | 3700 | U    | 1700 |      |
| Indeno(1,2,3-cd)pyrene      | U              | 330          | 3200           | J            | 4400           | 1700         | J              | 3800         | 770            | J            | 3700 | U    | 1700 |      |
| Dibenzo(a,h)anthracene      | U              | 330          | 1200           | J            | 4400           | U            | U              | 3800         | U              | 3700         | U    | 1700 |      |      |
| Benzo(g,h,i)perylene        | U              | 330          | 3600           | J            | 4400           | 1900         | J              | 3800         | 920            | J            | 3700 | U    | 1700 |      |

Table 1.3 (cont.) Results of the Analysis for BNA in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 (Results are Based on Dry Weight)

| Sample No.                  | A16661         |              | A16662         |              | A16663         |              | A16664         |              | A16665         |              |
|-----------------------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|
| Sample Location             | 11             |              | 12             |              | 13             |              | 14             |              | 15             |              |
| GC/MS File Name             | LS030          |              | LS031          |              | LS032          |              | LS033          |              | LS034          |              |
| Dilution Factor             | 10             |              | 5              |              | 5              |              | 5              |              | 5              |              |
| % Solid                     | 85             |              | 80             |              | 91             |              | 74             |              | 62             |              |
| Compound Name               | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg |
| Phenol                      | 9700           | 3900         | 4700           | 2100         | 2200           | 1800         | 7000           | 2300         | 1600           | J 2700       |
| bis(-2-Chloroethyl)Ether    | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 2-Chlorophenol              | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 1,3-Dichlorobenzene         | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 1,4-Dichlorobenzene         | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Benzyl alcohol              | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 1,2-Dichlorobenzene         | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 2-Methylphenol              | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| bis(2-Chloroisopropyl)ether | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 4-Methylphenol              | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| N-Nitroso-Di-n-propylamine  | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Hexachloroethane            | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Nitrobenzene                | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Isophorone                  | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 2-Nitrophenol               | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 2,4-Dimethylphenol          | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| bis(2-Chloroethoxy)methane  | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 2,4-Dichlorophenol          | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 1,2,4-Trichlorobenzene      | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Naphthalene                 | 2100           | J 3900       | 960            | J 2100       | 670            | J 1800       | 1400           | J 2300       | U              | 2700         |
| 4-Chloroaniline             | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Hexachlorobutadiene         | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 4-Chloro-3-methylphenol     | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 2-Methylnaphthalene         | U              | 3900         | 500            | J 2100       | 400            | J 1800       | U              | 2300         | U              | 2700         |
| Hexachlorocyclopentadiene   | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 2,4,6-Trichlorophenol       | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 2,4,5-Trichlorophenol       | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 2-Chloronaphthalene         | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 2-Nitroaniline              | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Dimethylphthalate           | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Acenaphthylene              | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 2,6-Dinitrotoluene          | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 3-Nitroaniline              | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Acenaphthene                | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 2,4-Dinitrophenol           | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 4-Nitrophenol               | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Dibenzofuran                | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 2,4-Dinitrotoluene          | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Diethylphthalate            | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 4-Chlorophenyl-phenylether  | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Fluorene                    | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 4-Nitroaniline              | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 4,6-Dinitro-2-methylphenol  | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| N-Nitrosodiphenylamine      | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 4-Bromophenyl-phenylether   | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Hexachlorobenzene           | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Pentachlorophenol           | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Phenanthrene                | U              | 3900         | 530            | J 2100       | U              | 1800         | U              | 2300         | U              | 2700         |
| Anthracene                  | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Carbazole                   | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Di-n-butylphthalate         | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Fluoranthene                | U              | 3900         | 600            | J 2100       | U              | 1800         | U              | 2300         | U              | 2700         |
| Pyrene                      | U              | 3900         | 900            | J 2100       | U              | 1800         | U              | 2300         | U              | 2700         |
| Butylbenzylphthalate        | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Benzo(a)anthracene          | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 3,3'-Dichlorobenzidine      | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Chrysene                    | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Bis(2-Ethylhexyl)phthalate  | 16000          | 3900         | 1300           | J 2100       | 540            | J 1800       | U              | 2300         | U              | 2700         |
| Di-n-octylphthalate         | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Benzo(b)fluoranthene        | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Benzo(k)fluoranthene        | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Benzo(a)pyrene              | U              | 3900         | U              | 2100         | U              | 1800         | 460            | J 2300       | U              | 2700         |
| Indeno(1,2,3-cd)pyrene      | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Dibenzo(a,h)anthracene      | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Benzo(g,h,i)perylene        | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |

Table 1.3 (cont.) Results of the Analysis for BNA in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 (Results are Based on Dry Weight)

| Sample No.                  | A16666         |              | A16667         |              | A16668         |              |
|-----------------------------|----------------|--------------|----------------|--------------|----------------|--------------|
| Sample Location             | Location 16    |              | Location 17    |              | Location 18    |              |
| GC/MS File Name             | LFS035         |              | LFS036         |              | LFS037         |              |
| Dilution Factor             | 5              |              | 5              |              | 5              |              |
| % Solid                     | 90             |              | 88             |              | 93             |              |
| Compound Name               | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg |
| Phenol                      | 840            | J 1900       | 910            | J 1900       | 11000          | 1800         |
| bis(-2-Chloroethyl)Ether    | U              | 1900         | U              | 1900         | U              | 1800         |
| 2-Chlorophenol              | U              | 1900         | U              | 1900         | U              | 1800         |
| 1,3-Dichlorobenzene         | U              | 1900         | U              | 1900         | U              | 1800         |
| 1,4-Dichlorobenzene         | U              | 1900         | U              | 1900         | U              | 1800         |
| Benzyl alcohol              | U              | 1900         | U              | 1900         | U              | 1800         |
| 1,2-Dichlorobenzene         | U              | 1900         | U              | 1900         | U              | 1800         |
| 2-Methylphenol              | U              | 1900         | U              | 1900         | U              | 1800         |
| bis(2-Chloroisopropyl)ether | U              | 1900         | U              | 1900         | U              | 1800         |
| 4-Methylphenol              | U              | 1900         | U              | 1900         | U              | 1800         |
| N-Nitroso-Di-n-propylamine  | U              | 1900         | U              | 1900         | U              | 1800         |
| Hexachloroethane            | U              | 1900         | U              | 1900         | U              | 1800         |
| Nitrobenzene                | U              | 1900         | U              | 1900         | U              | 1800         |
| Isophorone                  | U              | 1900         | U              | 1900         | U              | 1800         |
| 2-Nitrophenol               | U              | 1900         | U              | 1900         | U              | 1800         |
| 2,4-Dimethylphenol          | U              | 1900         | U              | 1900         | U              | 1800         |
| bis(2-Chloroethoxy)methane  | U              | 1900         | U              | 1900         | U              | 1800         |
| 2,4-Dichlorophenol          | U              | 1900         | U              | 1900         | U              | 1800         |
| 1,2,4-Trichlorobenzene      | U              | 1900         | U              | 1900         | U              | 1800         |
| Naphthalene                 | U              | 1900         | U              | 1900         | 850            | J 1800       |
| 4-Chloroaniline             | U              | 1900         | U              | 1900         | U              | 1800         |
| Hexachlorobutadiene         | U              | 1900         | U              | 1900         | U              | 1800         |
| 4-Chloro-3-methylphenol     | U              | 1900         | U              | 1900         | U              | 1800         |
| 2-Methylnaphthalene         | U              | 1900         | U              | 1900         | U              | 1800         |
| Hexachlorocyclopentadiene   | U              | 1900         | U              | 1900         | U              | 1800         |
| 2,4,6-Trichlorophenol       | U              | 1900         | U              | 1900         | U              | 1800         |
| 2,4,5-Trichlorophenol       | U              | 1900         | U              | 1900         | U              | 1800         |
| 2-Chloronaphthalene         | U              | 1900         | U              | 1900         | U              | 1800         |
| 2-Nitroaniline              | U              | 1900         | U              | 1900         | U              | 1800         |
| Dimethylphthalate           | U              | 1900         | U              | 1900         | U              | 1800         |
| Acenaphthylene              | U              | 1900         | U              | 1900         | U              | 1800         |
| 2,6-Dinitrotoluene          | U              | 1900         | U              | 1900         | U              | 1800         |
| 3-Nitroaniline              | U              | 1900         | U              | 1900         | U              | 1800         |
| Acenaphthene                | U              | 1900         | U              | 1900         | U              | 1800         |
| 2,4-Dinitrophenol           | U              | 1900         | U              | 1900         | U              | 1800         |
| 4-Nitrophenol               | U              | 1900         | U              | 1900         | U              | 1800         |
| Dibenzofuran                | U              | 1900         | U              | 1900         | U              | 1800         |
| 2,4-Dinitrotoluene          | U              | 1900         | U              | 1900         | U              | 1800         |
| Diethylphthalate            | U              | 1900         | U              | 1900         | U              | 1800         |
| 4-Chlorophenyl-phenylether  | U              | 1900         | U              | 1900         | U              | 1800         |
| Fluorene                    | U              | 1900         | U              | 1900         | U              | 1800         |
| 4-Nitroaniline              | U              | 1900         | U              | 1900         | U              | 1800         |
| 4,6-Dinitro-2-methylphenol  | U              | 1900         | U              | 1900         | U              | 1800         |
| N-Nitrosodiphenylamine      | U              | 1900         | U              | 1900         | U              | 1800         |
| 4-Bromophenyl-phenylether   | U              | 1900         | U              | 1900         | U              | 1800         |
| Hexachlorobenzene           | U              | 1900         | U              | 1900         | U              | 1800         |
| Pentachlorophenol           | U              | 1900         | U              | 1900         | U              | 1800         |
| Phenanthrene                | U              | 1900         | U              | 1900         | U              | 1800         |
| Anthracene                  | U              | 1900         | U              | 1900         | U              | 1800         |
| Carbazole                   | U              | 1900         | U              | 1900         | U              | 1800         |
| Di-n-butylphthalate         | U              | 1900         | U              | 1900         | U              | 1800         |
| Fluoranthene                | U              | 1900         | 860            | J 1900       | U              | 1800         |
| Pyrene                      | U              | 1900         | 740            | J 1900       | U              | 1800         |
| Butylbenzylphthalate        | U              | 1900         | U              | 1900         | U              | 1800         |
| Benzo(a)anthracene          | U              | 1900         | 410            | J 1900       | U              | 1800         |
| 3,3'-Dichlorobenzidine      | U              | 1900         | U              | 1900         | U              | 1800         |
| Chrysene                    | U              | 1900         | 510            | J 1900       | U              | 1800         |
| Bis(2-Ethylhexyl)phthalate  | U              | 1900         | U              | 1900         | U              | 1800         |
| Di-n-octylphthalate         | U              | 1900         | U              | 1900         | U              | 1800         |
| Benzo(b)fluoranthene        | U              | 1900         | 580            | J 1900       | U              | 1800         |
| Benzo(k)fluoranthene        | U              | 1900         | 520            | J 1900       | U              | 1800         |
| Benzo(a)pyrene              | U              | 1900         | 570            | J 1900       | U              | 1800         |
| Indeno(1,2,3-cd)pyrene      | U              | 1900         | 420            | J 1900       | U              | 1800         |
| Dibenzo(a,h)anthracene      | U              | 1900         | U              | 1900         | U              | 1800         |
| Benzo(g,h,i)perylene        | U              | 1900         | 470            | J 1900       | U              | 1800         |

Table 1.3 (cont.) Results of the Analysis for BNA in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 (Results are Based on Dry Weight)

| Sample No.                  | SBLK042099     | A16669       | A16670         | A16671       | A16672         |              |                |              |                |              |      |      |      |      |
|-----------------------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|------|------|------|------|
| Sample Location             | Lab Blank      | Location 19  | Location 20    | Location 21  | Location 22    |              |                |              |                |              |      |      |      |      |
| GC/MS File Name             | LFS041         | LFS042       | LFS043         | LFS044       | LFS045         |              |                |              |                |              |      |      |      |      |
| Dilution Factor             | 1              | 1            | 5              | 10           | 5              |              |                |              |                |              |      |      |      |      |
| % Solid                     | 100            | 93           | 91             | 48           | 81             |              |                |              |                |              |      |      |      |      |
| Compound Name               | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg |      |      |      |      |
| Phenol                      | U              | 330          | 4000           | 360          | 18000          | 1800         | 5600           | J            | 6900           | U            | 2000 |      |      |      |
| bis(-2-Chloroethyl)Ether    | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| 2-Chlorophenol              | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| 1,3-Dichlorobenzene         | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| 1,4-Dichlorobenzene         | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| Benzyl alcohol              | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| 1,2-Dichlorobenzene         | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| 2-Methylphenol              | U              | 330          | U              | 360          | 400            | J            | 1800           | U            | 6900           | U            | 2000 |      |      |      |
| bis(2-Chloroisopropyl)ether | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| 4-Methylphenol              | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| N-Nitroso-Di-n-propylamine  | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| Hexachloroethane            | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| Nitrobenzene                | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| Isophorone                  | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| 2-Nitrophenol               | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| 2,4-Dimethylphenol          | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| bis(2-Chloroethoxy)methane  | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| 2,4-Dichlorophenol          | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| 1,2,4-Trichlorobenzene      | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| Naphthalene                 | U              | 330          | 580            | 360          | 4400           | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| 4-Chloroaniline             | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| Hexachlorobutadiene         | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| 4-Chloro-3-methylphenol     | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| 2-Methylnaphthalene         | U              | 330          | 100            | J            | 360            | 1100         | J              | 1800         | U              | 6900         | U    | 2000 |      |      |
| Hexachlorocyclopentadiene   | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| 2,4,6-Trichlorophenol       | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| 2,4,5-Trichlorophenol       | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| 2-Chloronaphthalene         | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| 2-Nitroaniline              | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| Dimethylphthalate           | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| Acenaphthylene              | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | 420          | J    | 2000 |      |      |
| 2,6-Dinitrotoluene          | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| 3-Nitroaniline              | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| Acenaphthene                | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| 2,4-Dinitrophenol           | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| 4-Nitrophenol               | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| Dibenzofuran                | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| 2,4-Dinitrotoluene          | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| Diethylphthalate            | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| 4-Chlorophenyl-phenylether  | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| Fluorene                    | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| 4-Nitroaniline              | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| 4,6-Dinitro-2-methylphenol  | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| N-Nitrosodiphenylamine      | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| 4-Bromophenyl-phenylether   | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| Hexachlorobenzene           | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| Pentachlorophenol           | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| Phenanthrene                | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| Anthracene                  | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| Carbazole                   | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| Di-n-butylphthalate         | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| Fluoranthene                | U              | 330          | 74             | J            | 360            | 1600         | J              | 1800         | 3400           | J            | 6900 | 4700 | 2000 |      |
| Pyrene                      | U              | 330          | U              | 360          | 1400           | J            | 1800           | 3400         | J              | 6900         | 4100 | 2000 |      |      |
| Butylbenzylphthalate        | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| Benzo(a)anthracene          | U              | 330          | U              | 360          | 840            | J            | 1800           | 2100         | J              | 6900         | 2600 | 2000 |      |      |
| 3,3'-Dichlorobenzidine      | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| Chrysene                    | U              | 330          | U              | 360          | 900            | J            | 1800           | 2600         | J              | 6900         | 2400 | 2000 |      |      |
| Bis(2-Ethylhexyl)phthalate  | U              | 330          | 180            | J            | 360            | 760          | J              | 1800         | 1700           | J            | 6900 | 430  | J    | 2000 |
| Di-n-octylphthalate         | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| Benzo(b)fluoranthene        | U              | 330          | U              | 360          | 910            | J            | 1800           | 3500         | J              | 6900         | 2700 | 2000 |      |      |
| Benzo(k)fluoranthene        | U              | 330          | U              | 360          | 890            | J            | 1800           | 3400         | J              | 6900         | 2500 | 2000 |      |      |
| Benzo(a)pyrene              | U              | 330          | U              | 360          | 980            | J            | 1800           | 3700         | J              | 6900         | 3100 | 2000 |      |      |
| Indeno(1,2,3-cd)pyrene      | U              | 330          | U              | 360          | 660            | J            | 1800           | 3100         | J              | 6900         | 1600 | J    | 2000 |      |
| Dibenzo(a,h)anthracene      | U              | 330          | U              | 360          | U              | 1800         | U              | U            | 6900           | U            | 2000 |      |      |      |
| Benzo(g,h,i)perylene        | U              | 330          | U              | 360          | 770            | J            | 1800           | 3600         | J              | 6900         | 1700 | J    | 2000 |      |

Table 1.3 (cont.) Results of the Analysis for BNA in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 (Results are Based on Dry Weight)

| Sample No.                  | A16673      |        | A16674      |        | A16675      |        | A16676        |       | A16677      |        |
|-----------------------------|-------------|--------|-------------|--------|-------------|--------|---------------|-------|-------------|--------|
| Sample Location             | Location 23 |        | Location 24 |        | Location 25 |        | Location 25-D |       | Location 26 |        |
| GC/MS File Name             | LFS048      |        | LFS049      |        | LFS050      |        | LFS051        |       | LFS052      |        |
| Dilution Factor             | 10          |        | 5           |        | 10          |        | 1             |       | 10          |        |
| % Solid                     | 82          |        | 75          |        | 83          |        | 77            |       | 70          |        |
| Compound Name               | Conc.       | MDL    | Conc.       | MDL    | Conc.       | MDL    | Conc.         | MDL   | Conc.       | MDL    |
|                             | µg/kg       | µg/kg  | µg/kg       | µg/kg  | µg/kg       | µg/kg  | µg/kg         | µg/kg | µg/kg       | µg/kg  |
| Phenol                      | 2400        | J 4100 | 4000        | 2200   | 990         | J 4000 | U             | 430   | U           | 4800   |
| bis(-2-Chloroethyl)Ether    | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| 2-Chlorophenol              | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| 1,3-Dichlorobenzene         | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| 1,4-Dichlorobenzene         | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| Benzyl alcohol              | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| 1,2-Dichlorobenzene         | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| 2-Methylphenol              | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| bis(2-Chloroisopropyl)ether | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| 4-Methylphenol              | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| N-Nitroso-Di-n-propylamine  | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| Hexachloroethane            | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| Nitrobenzene                | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| Isophorone                  | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| 2-Nitrophenol               | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| 2,4-Dimethylphenol          | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| bis(2-Chloroethoxy)methane  | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| 2,4-Dichlorophenol          | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| 1,2,4-Trichlorobenzene      | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| Naphthalene                 | 1100        | J 4100 | 1400        | J 2200 | U           | 4000   | U             | 430   | U           | 4800   |
| 4-Chloroaniline             | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| Hexachlorobutadiene         | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| 4-Chloro-3-methylphenol     | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| 2-Methylnaphthalene         | 850         | J 4100 | 790         | J 2200 | U           | 4000   | U             | 430   | U           | 4800   |
| Hexachlorocyclopentadiene   | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| 2,4,6-Trichlorophenol       | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| 2,4,5-Trichlorophenol       | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| 2-Chloronaphthalene         | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| 2-Nitroaniline              | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| Dimethylphthalate           | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| Acenaphthylene              | 1300        | J 4100 | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| 2,6-Dinitrotoluene          | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| 3-Nitroaniline              | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| Acenaphthene                | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| 2,4-Dinitrophenol           | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| 4-Nitrophenol               | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| Dibenzofuran                | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| 2,4-Dinitrotoluene          | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| Diethylphthalate            | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| 4-Chlorophenyl-phenylether  | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| Fluorene                    | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| 4-Nitroaniline              | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| 4,6-Dinitro-2-methylphenol  | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| N-Nitrosodiphenylamine      | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| 4-Bromophenyl-phenylether   | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| Hexachlorobenzene           | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| Pentachlorophenol           | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| Phenanthrene                | 2200        | J 4100 | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| Anthracene                  | 1400        | J 4100 | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| Carbazole                   | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| Di-n-butylphthalate         | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| Fluoranthene                | 8500        | 4100   | 1200        | J 2200 | 5800        | 4000   | 100           | J 430 | 3300        | J 4800 |
| Pyrene                      | 8800        | 4100   | 940         | J 2200 | 4800        | 4000   | U             | 430   | 2900        | J 4800 |
| Butylbenzylphthalate        | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| Benzo(a)anthracene          | 5200        | 4100   | 650         | J 2200 | 2700        | J 4000 | U             | 430   | 1600        | J 4800 |
| 3,3'-Dichlorobenzidine      | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| Chrysene                    | 5400        | 4100   | 910         | J 2200 | 3100        | J 4000 | U             | 430   | 2000        | J 4800 |
| Bis(2-Ethylhexyl)phthalate  | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | 1100        | J 4800 |
| Di-n-octylphthalate         | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| Benzo(b)fluoranthene        | 5700        | 4100   | 1200        | J 2200 | 3600        | J 4000 | U             | 430   | 2100        | J 4800 |
| Benzo(k)fluoranthene        | 5700        | 4100   | 1000        | J 2200 | 3200        | J 4000 | U             | 430   | 1900        | J 4800 |
| Benzo(a)pyrene              | 5300        | 4100   | 1100        | J 2200 | 3700        | J 4000 | U             | 430   | 2200        | J 4800 |
| Indeno(1,2,3-cd)pyrene      | 3100        | J 4100 | 860         | J 2200 | 2400        | J 4000 | U             | 430   | 1400        | J 4800 |
| Dibenzo(a,h)anthracene      | U           | 4100   | U           | 2200   | U           | 4000   | U             | 430   | U           | 4800   |
| Benzo(g,h,i)perylene        | 3200        | J 4100 | 970         | J 2200 | 2700        | J 4000 | U             | 430   | U           | 4800   |

Table 1.3 (cont.) Results of the Analysis for BNA in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 (Results are Based on Dry Weight)

| Sample No.                  | A16678         | A16679       |                |              |
|-----------------------------|----------------|--------------|----------------|--------------|
| Sample Location             | Location 26-D  | Location 27  |                |              |
| GC/MS File Name             | LFS053         | LFS054       |                |              |
| Dilution Factor             | 1              | 5            |                |              |
| % Solid                     | 77             | 83           |                |              |
| Compound Name               | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg |
| Phenol                      | U              | 440          | 16000          | 2000         |
| bis(-2-Chloroethyl)Ether    | U              | 440          | U              | 2000         |
| 2-Chlorophenol              | U              | 440          | U              | 2000         |
| 1,3-Dichlorobenzene         | U              | 440          | U              | 2000         |
| 1,4-Dichlorobenzene         | U              | 440          | U              | 2000         |
| Benzyl alcohol              | U              | 440          | U              | 2000         |
| 1,2-Dichlorobenzene         | U              | 440          | U              | 2000         |
| 2-Methylphenol              | U              | 440          | 820            | J 2000       |
| bis(2-Chloroisopropyl)ether | U              | 440          | U              | 2000         |
| 4-Methylphenol              | U              | 440          | U              | 2000         |
| N-Nitroso-Di-n-propylamine  | U              | 440          | U              | 2000         |
| Hexachloroethane            | U              | 440          | U              | 2000         |
| Nitrobenzene                | U              | 440          | U              | 2000         |
| Isophorone                  | U              | 440          | U              | 2000         |
| 2-Nitrophenol               | U              | 440          | U              | 2000         |
| 2,4-Dimethylphenol          | U              | 440          | U              | 2000         |
| bis(2-Chloroethoxy)methane  | U              | 440          | U              | 2000         |
| 2,4-Dichlorophenol          | U              | 440          | U              | 2000         |
| 1,2,4-Trichlorobenzene      | U              | 440          | 1700           | J 2000       |
| Naphthalene                 | U              | 440          | 2900           | J 2000       |
| 4-Chloroaniline             | U              | 440          | U              | 2000         |
| Hexachlorobutadiene         | U              | 440          | U              | 2000         |
| 4-Chloro-3-methylphenol     | U              | 440          | U              | 2000         |
| 2-Methylnaphthalene         | U              | 440          | 810            | J 2000       |
| Hexachlorocyclopentadiene   | U              | 440          | U              | 2000         |
| 2,4,6-Trichlorophenol       | U              | 440          | U              | 2000         |
| 2,4,5-Trichlorophenol       | U              | 440          | U              | 2000         |
| 2-Chloronaphthalene         | U              | 440          | U              | 2000         |
| 2-Nitroaniline              | U              | 440          | U              | 2000         |
| Dimethylphthalate           | U              | 440          | U              | 2000         |
| Acenaphthylene              | 96             | J 440        | U              | 2000         |
| 2,6-Dinitrotoluene          | U              | 440          | U              | 2000         |
| 3-Nitroaniline              | U              | 440          | U              | 2000         |
| Acenaphthene                | U              | 440          | U              | 2000         |
| 2,4-Dinitrophenol           | U              | 440          | U              | 2000         |
| 4-Nitrophenol               | U              | 440          | U              | 2000         |
| Dibenzofuran                | U              | 440          | U              | 2000         |
| 2,4-Dinitrotoluene          | U              | 440          | U              | 2000         |
| Diethylphthalate            | U              | 440          | U              | 2000         |
| 4-Chlorophenyl-phenylether  | U              | 440          | U              | 2000         |
| Fluorene                    | U              | 440          | U              | 2000         |
| 4-Nitroaniline              | U              | 440          | U              | 2000         |
| 4,6-Dinitro-2-methylphenol  | U              | 440          | U              | 2000         |
| N-Nitrosodiphenylamine      | U              | 440          | U              | 2000         |
| 4-Bromophenyl-phenylether   | U              | 440          | U              | 2000         |
| Hexachlorobenzene           | U              | 440          | U              | 2000         |
| Pentachlorophenol           | U              | 440          | U              | 2000         |
| Phenanthrene                | 440            | U            | U              | 2000         |
| Anthracene                  | 92             | J 440        | U              | 2000         |
| Carbazole                   | U              | 440          | U              | 2000         |
| Di-n-butylphthalate         | U              | 440          | U              | 2000         |
| Fluoranthene                | 930            | 440          | 420            | J 2000       |
| Pyrene                      | 810            | 440          | 500            | J 2000       |
| Butylbenzylphthalate        | U              | 440          | U              | 2000         |
| Benzo(a)anthracene          | 470            | 440          | U              | 2000         |
| 3,3'-Dichlorobenzidine      | U              | 440          | U              | 2000         |
| Chrysene                    | 500            | 440          | U              | 2000         |
| Bis(2-Ethylhexyl)phthalate  | U              | 440          | 1000           | J 2000       |
| Di-n-octylphthalate         | U              | 440          | U              | 2000         |
| Benzo(b)fluoranthene        | 540            | 440          | U              | 2000         |
| Benzo(k)fluoranthene        | 520            | 440          | U              | 2000         |
| Benzo(a)pyrene              | 600            | 440          | U              | 2000         |
| Indeno(1,2,3-cd)pyrene      | 360            | J 440        | U              | 2000         |
| Dibenzo(a,h)anthracene      | U              | 440          | U              | 2000         |
| Benzo(g,h,i)perylene        | 400            | J 440        | U              | 2000         |



Table 1.4 Results of TIC for BNA in Soil  
WA # 3-439 Lackawanna Foundry Site

| Sample ID   | LabFile# | CAS #       | Compound   | Conc**<br>ug/kg |
|-------------|----------|-------------|--|-----------------|
| SBLK 041599 | LFS 002  |             | No TICs were detected                              |                 |
| A 16650     | LFS 003  |             | No TICs were detected                              |                 |
| A 16653     | LFS 008  |             | Unknown  | 2500            |
| A 16653     | LFS 008  |             | Unknown Alkane                                     | 3800            |
| A 16653     | LFS 008  |             | Unknown Alkane                                     | 5900            |
| A 16654     | LFS 009  |             | Unknown Aldehyde + Unknown                         | 1700            |
| A 16654     | LFS 009  |             | Unknown Alkane                                     | 2600            |
| A 16656     | LFS 012  |             | No TICs were detected                              |                 |
| SBLK 041999 | LFS 023  |             | Unknown  | 210             |
| SBLK 041999 | LFS 023  |             | Unknown  | 140             |
| A 16659     | LFS 026  |             | Benzene, methylenebis[isocyanato-] + alkyl benzene | 1600            |
| A 16665     | LFS 034  |             | No TICs were detected                              |                 |
| A 16666     | LFS 035  |             | Unknown  | 890             |
| A 16666     | LFS 035  |             | Unknown Phthalate                                  | 760             |
| A 16666     | LFS 035  | 000127-63-9 | Diphenyl sulfone                                   | 1200            |
| A 16667     | LFS 036  |             | No TICs were detected                              |                 |
| SBLK 042099 | LFS 041  |             | Unknown  | 160             |
| A 16671     | LFS 044  |             | Phenylpropyl-pyridine isomer                       | 4000            |
| A 16671     | LFS 044  |             | Hexadecenoic acid + unknown                        | 5600            |
| A 16675     | LFS 050  |             | Unknown PAH isomer                                 | 2700            |
| A 16676     | LFS 051  |             | Unknown  | 340             |
| A 16676     | LFS 051  |             | Unknown  | 210             |

\*\* the concentration is estimated - the response factor was assumed to be 1

00024

Table 1.4 (cont.) Results of TIC for BNA in Soil  
WA # 3-439 Lackawanna Foundry Site

Sample #  
LabFile#

A16651  
LFS004

Con. Factor

25

|    | CAS#        | Compound   | Q  | RT    | Cor.<br>µg/kg |
|----|-------------|--|----|-------|---------------|
| 1  |             | Trimethyl-benzene isomer                           |    | 4.92  | 1600          |
| 2  |             | Trimethyl-benzene isomer                           |    | 5.09  | 1500          |
| 3  |             | Trimethyl-benzene isomer                           |    | 5.21  | 4900          |
| 4  |             | Trimethyl-benzene isomer                           |    | 5.48  | 1900          |
| 5  |             | Ethyl-dimethyl-benzene isomer                      |    | 5.73  | 2200          |
| 6  |             | 2-butoxyethyl acetate                              |    | 5.87  | 4300          |
| 7  |             | Ethyl-dimethyl-benzene isomer                      |    | 5.91  | 860           |
| 8  |             | Ethyl-dimethyl-benzene isomer                      |    | 5.93  | 650           |
| 9  |             | Ethyl-dimethyl-benzene isomer                      |    | 5.98  | 1200          |
| 10 |             | Ethyl-dimethyl-benzene isomer                      |    | 6.25  | 870           |
| 11 |             | Ethyl-dimethyl-benzene isomer                      |    | 6.29  | 710           |
| 12 |             | Alkyl benzene isomer                               |    | 6.43  | 670           |
| 13 |             | Alkylbenzene isomer + dihydro-methyl-indene isomer |    | 6.57  | 1100          |
| 14 | 000627-93-0 | Hexanedioic acid, dimethyl ester                   | 91 | 7.08  | 670           |
| 15 | 000090-01-7 | Salicyl Alcohol                                    | 90 | 7.39  | 2600          |
| 16 |             | Unknown acid                                       |    | 8.36  | 560           |
| 17 |             | Phenylpropyl-pyridine + unknown                    |    | 10.69 | 4200          |
| 18 |             | Phthalate isomer                                   |    | 11.08 | 5100          |
| 19 | 000101-68-8 | Benzene, 1,1'-methylenebis(4-isocyanato-)          | 99 | 12.37 | 3400          |
| 20 |             | Unknown PAH isomer                                 |    | 18.15 | 820           |

\* Estimated Concentration (Response Factor = 1)

00025

Table 1.4 (cont.) Results of TIC for BNA in Soil  
 WA # 3-439 Lackawanna Foundry Site

| Sample # | LabFile#    | A16652  | LFS005 | Con. Factor | 180             |
|----------|-------------|---|--------|-------------|-----------------|
|          | CAS#        | Compound  | Q      | RT          | Conc.*<br>µg/kg |
| 1        |             | Trimethyl-benzene isomer                            |        | 4.92        | 1600            |
| 2        |             | Trimethyl-benzene isomer                            |        | 5.09        | 1600            |
| 3        |             | Trimethyl-benzene isomer                            |        | 5.21        | 5400            |
| 4        |             | Trimethyl-benzene isomer                            |        | 5.49        | 2000            |
| 5        |             | Ethyl-dimethyl-benzene isomer                       |        | 5.73        | 2100            |
| 6        |             | Butoxyethyl acetate                                 |        | 5.87        | 6500            |
| 7        |             | Ethyl-dimethyl-benzene isomer                       |        | 5.91        | 810             |
| 8        |             | Ethyl-dimethyl-benzene isomer                       |        | 5.98        | 940             |
| 9        |             | Alkyl benzene isomer + dihydro-methyl-indene isomer |        | 6.57        | 990             |
| 10       | 000090-1-7  | Salicyl Alcohol                                     | 90     | 7.39        | 2500            |
| 11       |             | Phenylpropyl-pyridine + unknown                     |        | 10.69       | 5900            |
| 12       |             | Phthalate isomer                                    |        | 11.08       | 6700            |
| 13       |             | Unknown   |        | 12.02       | 810             |
| 14       | 000101-68-8 | Benzene, 1,1'-methylenebis(4-isocyanato)-           | 99     | 12.37       | 9400            |
| 15       |             |   |        |             |                 |
| 16       |             |   |        |             |                 |
| 17       |             |   |        |             |                 |
| 18       |             |   |        |             |                 |
| 19       |             |   |        |             |                 |
| 20       |             |   |        |             |                 |

\* Estimated Concentration (Response Factor = 1)

00026

Table 1.4 (cont.) Results of TIC for BNA in Soil  
WA # 3-439 Lackawanna Foundry Site

Sample # A16655  
LabFile# LFS010

Con. Factor 41

|    | CAS#        | Compound                           | Q  | RT    | Cor<br>µg/kg |
|----|-------------|------------------------------------|----|-------|--------------|
| 1  | 000132-65-0 | Dibenzothiophene                   | 97 | 10.78 | 500          |
| 2  |             | Methyl-phenanthrene isomer         |    | 11.53 | 780          |
| 3  |             | Methyl-phenanthrene isomer         |    | 11.57 | 920          |
| 4  |             | Methyl-phenanthrene isomer         |    | 11.61 | 380          |
| 5  |             | Cyclopenta[xxx]phenanthrene isomer |    | 11.70 | 2000         |
| 6  |             | Unknown PAH isomer                 |    | 11.89 | 640          |
| 7  |             | 9,10-Anthracenedione + unknown     |    | 11.95 | 690          |
| 8  |             | Unknown + unknown PAH isomer       |    | 12.31 | 420          |
| 9  |             | Unknown                            |    | 16.80 | 480          |
| 10 |             | Unknown + unknown alkane           |    | 17.08 | 1400         |
| 11 |             | Unknown PAH isomer                 |    | 17.59 | 1100         |
| 12 |             | Unknown PAH isomer                 |    | 18.18 | 4900         |
| 13 |             | Unknown PAH isomer                 |    | 18.65 | 1500         |
| 14 |             | Unknown alkane                     |    | 19.69 | 1600         |
| 15 |             | Unknown aldehyde + unknown         |    | 22.22 | 1200         |
| 16 |             | Unknown PAH isomer                 |    | 22.45 | 550          |
| 17 |             | Unknown PAH isomer                 |    | 22.54 | 460          |
| 18 |             | Unknown PAH isomer                 |    | 23.33 | 420          |
| 19 |             | Unknown PAH isomer                 |    | 23.49 | 860          |
| 20 |             | Unknown PAH isomer                 |    | 29.50 | 1300         |

\* Estimated Concentration (Response Factor = 1)

00027

Table 1.4 (cont.) Results of TIC for BNA in Soil  
 WA # 3-439 Lackawanna Foundry Site

| Sample # | A16657 |                             |   | Con. Factor | 440             |
|----------|--------|-----------------------------|---|-------------|-----------------|
| LabFile# | LFS024 |                             |   |             |                 |
|          | CAS#   | Compound                    | Q | RT          | Conc.*<br>µg/kg |
| 1        |        | Unknown                     |   | 12.31       | 2300            |
| 2        |        | Unknown aldehyde            |   | 16.53       | 2500            |
| 3        |        | Unknown alkane              |   | 17.05       | 2900            |
| 4        |        | Unknown benzo pyrene isomer |   | 18.14       | 2400            |
| 5        |        | Unknown alkane              |   | 19.66       | 6900            |
| 6        |        |                             |   |             |                 |
| 7        |        |                             |   |             |                 |
| 8        |        |                             |   |             |                 |
| 9        |        |                             |   |             |                 |
| 10       |        |                             |   |             |                 |
| 11       |        |                             |   |             |                 |
| 12       |        |                             |   |             |                 |
| 13       |        |                             |   |             |                 |
| 14       |        |                             |   |             |                 |
| 15       |        |                             |   |             |                 |
| 16       |        |                             |   |             |                 |
| 17       |        |                             |   |             |                 |
| 18       |        |                             |   |             |                 |
| 19       |        |                             |   |             |                 |
| 20       |        |                             |   |             |                 |

\* Estimated Concentration (Response Factor = 1)

Table 1.4 (cont.) Results of TIC for BNA in Soil  
 WA # 3-439 Lackawanna Foundry Site

Sample #  
 LabFile#

A16658  
 LFS025

Con. Factor

|    | CAS# | Compound                                       | Q | RT    | Con.<br>µg/kg |
|----|------|--|---|-------|---------------|
| 1  |      | Unknown  |   | 3.90  | 2300          |
| 2  |      | Benzoic acid, [[[(acetylamino)sulfonyl]phenyl] |   | 7.86  | 1900          |
| 3  |      | Unknown ketone                                 |   | 8.37  | 2300          |
| 4  |      | Unknown  |   | 11.54 | 1700          |
| 5  |      | Unknown  |   | 12.31 | 4100          |
| 6  |      | Unknown alkane                                 |   | 19.67 | 1800          |
| 7  |      |  |   |       |               |
| 8  |      |  |   |       |               |
| 9  |      |  |   |       |               |
| 10 |      |  |   |       |               |
| 11 |      |  |   |       |               |
| 12 |      |  |   |       |               |
| 13 |      |  |   |       |               |
| 14 |      |  |   |       |               |
| 15 |      |  |   |       |               |
| 16 |      |  |   |       |               |
| 17 |      |  |   |       |               |
| 18 |      |  |   |       |               |
| 19 |      |  |   |       |               |
| 20 |      |  |   |       |               |

\* Estimated Concentration (Response Factor = 1)

00029

Table 1.4 (cont.) Results of TIC for BNA in Soil  
WA # 3-439 Lackawanna Foundry Site

Sample #  
LabFile#

A16660  
LFS027

Con. Factor 170

|    | CAS# | Compound                                       | Q | RT    | Conc.*<br>µg/kg |
|----|------|--|---|-------|-----------------|
| 1  |      | Trimethyl-benzene isomer                       |   | 5.21  | 1300            |
| 2  |      | Trimethyl-benzene isomer                       |   | 5.48  | 830             |
| 3  |      | Unknown acid                                   |   | 6.26  | 1400            |
| 4  |      | Phenyl-butene                                  |   | 6.57  | 740             |
| 5  |      | Methyl-naphthalene isomer                      |   | 7.87  | 1100            |
| 6  |      | Dimethyl-naphthalene isomer                    |   | 8.61  | 980             |
| 7  |      | Trimethyl-naphthalene isomer + unknown         |   | 9.28  | 720             |
| 8  |      | Unknown alkane + unknown                       |   | 9.44  | 1400            |
| 9  |      | Trimethyl-naphthalene isomer + unknown alkane  |   | 9.72  | 1400            |
| 10 |      | Unknown alkane                                 |   | 10.58 | 1300            |
| 11 |      | Unknown alkane                                 |   | 10.63 | 910             |
| 12 |      | Unknown alkane                                 |   | 11.10 | 930             |
| 13 |      | Trichloro-biphenyl isomer + unknown            |   | 11.31 | 710             |
| 14 |      | Trichloro-biphenyl isomer + unknown            |   | 11.33 | 710             |
| 15 |      | Trichloro-biphenyl isomer + unknown            |   | 11.53 | 690             |
| 16 |      | Unknown phthalate + methyl-phenanthrene isomer |   | 11.56 | 690             |
| 17 |      | Unknown alkane                                 |   | 11.60 | 1100            |
| 18 |      | Unknown alkane                                 |   | 12.08 | 980             |
| 19 |      | Unknown alkane                                 |   | 12.54 | 740             |
| 20 |      | Unknown alkane                                 |   | 12.71 | 1500            |

\* Estimated Concentration (Response Factor = 1)

00000

Table 1.4 (cont.) Results of TIC for BNA in Soil  
 WA # 3-439 Lackawanna Foundry Site

Sample #  
 LabFile#

A16661  
 LFS030

Con. Factor

|    | CAS#        | Compound                                   | Q  | RT    | Cor.<br>µg/kg |
|----|-------------|--|----|-------|---------------|
| 1  |             | Unknown                                    |    | 3.90  | 1900          |
| 2  |             | Trimethyl-benzene isomer                   |    | 5.21  | 3400          |
| 3  |             | Hydroxy benzaldehyde + alkyl benzene       |    | 5.68  | 2300          |
| 4  |             | Tetramethyl-benzene isomer                 |    | 5.73  | 2500          |
| 5  |             | Tetramethyl-benzene isomer                 |    | 5.98  | 2100          |
| 6  |             | Tetramethyl-benzene isomer + unknown       |    | 6.25  | 2400          |
| 7  |             | Tetramethyl-benzene isomer                 |    | 6.29  | 2400          |
| 8  |             | Tetramethyl-benzene isomer + alkyl benzene |    | 6.57  | 2600          |
| 9  |             | Pyridine, (phenylpropyl-) + unknown        |    | 10.69 | 2100          |
| 10 |             | Unknown phthalate                          |    | 11.08 | 3600          |
| 11 | 000127-63-9 | Diphenyl sulfone                           | 96 | 11.62 | 7700          |
| 12 | 000101-68-8 | Benzene, 1,1'-methylenebis[4-isocyanato-   | 98 | 12.37 | 5700          |
| 13 |             |  |    |       |               |
| 14 |             |  |    |       |               |
| 15 |             |  |    |       |               |
| 16 |             |  |    |       |               |
| 17 |             |  |    |       |               |
| 18 |             |  |    |       |               |
| 19 |             |  |    |       |               |
| 20 |             |  |    |       |               |

\* Estimated Concentration (Response Factor = 1)

00031



Table 1.4 (cont.) Results of TIC for BNA in Soil  
WA # 3-439 Lackawanna Foundry Site

| Sample # | LabFile#    | A16662                                       | LFS031 | Con. Factor | 210             |
|----------|-------------|--|--------|-------------|-----------------|
|          | CAS#        | Compound                                     | Q      | RT          | Conc.*<br>µg/kg |
| 1        |             | Trimethyl-benzene isomer                     |        | 5.21        | 1400            |
| 2        |             | Benzaldehyde, hydroxy- + alkyl benzene       |        | 5.68        | 960             |
| 3        | 000090-01-7 | Salicyl Alcohol                              | 91     | 7.39        | 1200            |
| 4        |             | Unknown alkane                               |        | 8.82        | 1200            |
| 5        |             | Unknown alkane                               |        | 9.44        | 2600            |
| 6        |             | Unknown alkane                               |        | 9.73        | 2600            |
| 7        |             | Azulene, ethyl dimethyl- + unknown           |        | 10.28       | 1200            |
| 8        |             | Unknown alkane                               |        | 10.58       | 2200            |
| 9        |             | Unknown alkane                               |        | 10.64       | 3100            |
| 10       |             | Pyridine, (phenylpropyl-) + unknown          |        | 10.69       | 1400            |
| 11       |             | Unknown phthalate                            |        | 11.07       | 1000            |
| 12       |             | Unknown alkane                               |        | 11.10       | 1600            |
| 13       |             | Trichloro-biphenyl isomer + unknown          |        | 11.33       | 1300            |
| 14       |             | Trichloro-biphenyl isomer + unknown          |        | 11.52       | 1000            |
| 15       |             | Unknown alkane                               |        | 11.61       | 1600            |
| 16       | 000127-63-9 | Diphenyl sulfone                             | 93     | 11.62       | 2600            |
| 17       |             | Tetra-chloro-biphenyl isomer + unknown       |        | 11.67       | 1100            |
| 18       |             | Unknown alkane                               |        | 12.08       | 1500            |
| 19       |             | Benzene, methylenebis[isocyanato-] + unknown |        | 12.37       | 1700            |
| 20       |             | Trimethyl-phenanthrene isomer                |        | 12.79       | 1300            |

\* Estimated Concentration (Response Factor = 1)

00032

Table 1.4 (cont.) Results of TIC for BNA in Soil  
 WA # 3-439 Lackawanna Foundry Site

Sample #  
 LabFile#

A16663  
 LFS032

Con. Factor

|    | CAS#        | Compound                               | Q  | RT    | Conc<br>µg/kg |
|----|-------------|--|----|-------|---------------|
| 1  |             | Unknown                                |    | 10.63 | 1700          |
| 2  |             | Unknown phthalate                      |    | 11.08 | 730           |
| 3  | 000127-63-9 | Diphenyl sulfone                       | 90 | 11.62 | 750           |
| 4  |             | Unknown                                |    | 12.09 | 2600          |
| 5  |             | Tetra-chloro-biphenyl isomer + unknown |    | 12.52 | 750           |
| 6  |             | Unknown                                |    | 13.41 | 1500          |
| 7  |             |  |    |       |               |
| 8  |             |  |    |       |               |
| 9  |             |  |    |       |               |
| 10 |             |  |    |       |               |
| 11 |             |  |    |       |               |
| 12 |             |  |    |       |               |
| 13 |             |  |    |       |               |
| 14 |             |  |    |       |               |
| 15 |             |  |    |       |               |
| 16 |             |  |    |       |               |
| 17 |             |  |    |       |               |
| 18 |             |  |    |       |               |
| 19 |             |  |    |       |               |
| 20 |             |  |    |       |               |

\* Estimated Concentration (Response Factor = 1)

Table 1.4 (cont.) Results of TIC for BNA in Soil  
WA # 3-439 Lackawanna Foundry Site

| Sample # | A16664      |  | Con. Factor | 230   |                 |
|----------|-------------|--|-------------|-------|-----------------|
| LabFile# | LFS033      |  |             |       |                 |
|          | CAS#        | Compound                                   | Q           | RT    | Conc.*<br>µg/kg |
| 1        |             | Trimethyl-benzene isomer                   |             | 5.21  | 1500            |
| 2        |             | Hydroxy-benzaldehyde + alkyl benzene       |             | 5.68  | 1700            |
| 3        |             | Tetramethyl-benzene isomer                 |             | 5.73  | 1500            |
| 4        |             | Tetramethyl-benzene isomer                 |             | 5.98  | 1200            |
| 5        |             | Tetramethyl-benzene isomer                 |             | 6.25  | 1500            |
| 6        |             | Tetramethyl-benzene isomer                 |             | 6.29  | 1700            |
| 7        |             | Tetramethyl-benzene isomer + alkyl benzene |             | 6.57  | 1600            |
| 8        |             | Phenylpropyl-pyridine                      |             | 10.69 | 1900            |
| 9        |             | Unknown phthalate                          |             | 11.08 | 2500            |
| 10       |             | Trichloro-biphenyl isomer                  |             | 11.31 | 1000            |
| 11       |             | Trichloro-biphenyl isomer                  |             | 11.33 | 1300            |
| 12       | 000127-63-9 | Diphenyl sulfone                           | 95          | 11.62 | 1300            |
| 13       |             | Tetrachloro-biphenyl isomer + unknown      |             | 11.67 | 1200            |
| 14       |             | Tetrachloro-biphenyl isomer + unknown      |             | 11.91 | 1200            |
| 15       |             | Tetrachloro-biphenyl isomer + unknown      |             | 12.29 | 1100            |
| 16       |             | Tetrachloro-biphenyl isomer + unknown      |             | 12.34 | 950             |
| 17       |             | Benzene, methylenebis[isocyanato-]         |             | 12.37 | 3300            |
| 18       |             | Tetrachloro-biphenyl isomer + unknown      |             | 12.52 | 1000            |
| 19       |             |  |             |       |                 |
| 20       |             |  |             |       |                 |

\* Estimated Concentration (Response Factor = 1)

00034

Table 1.4 (cont.) Results of TIC for BNA in Soil  
 WA # 3-439 Lackawanna Foundry Site

Sample #  
 LabFile#

A16668  
 LFS037

Con. Factor

1

|    | CAS# | Compound                                     | Q | RT    | Conc.<br>µg/kg |
|----|------|--|---|-------|----------------|
| 1  |      | Trimethyl-benzene isomer                     |   | 5.21  | 2300           |
| 2  |      | Trimethyl-benzene isomer                     |   | 5.48  | 790            |
| 3  |      | Benzaldehyde, hydroxy + alkyl benzene        |   | 5.68  | 4300           |
| 4  |      | Tetramethyl-benzene isomer                   |   | 5.73  | 880            |
| 5  |      | Butoxyethyl acetate                          |   | 5.86  | 1500           |
| 6  |      | Dimethyl ester hexanedioic acid + unknown    |   | 7.08  | 760            |
| 7  |      | Unknown                                      |   | 7.34  | 970            |
| 8  |      | Unknown                                      |   | 7.97  | 1200           |
| 9  |      | Phenylpropyl-pyridine                        |   | 10.69 | 940            |
| 10 |      | Unknown phthalate                            |   | 11.08 | 4400           |
| 11 |      | Benzene, methylenebis[isocyanato-] + unknown |   | 12.37 | 3800           |
| 12 |      | Unknown alkane                               |   | 17.06 | 1100           |
| 13 |      | Unknown aldehyde + unknown                   |   | 18.95 | 1000           |
| 14 |      | Unknown aldehyde + unknown                   |   | 22.20 | 1000           |
| 15 |      |  |   |       |                |
| 16 |      |  |   |       |                |
| 17 |      |  |   |       |                |
| 18 |      |  |   |       |                |
| 19 |      |  |   |       |                |
| 20 |      |  |   |       |                |

\* Estimated Concentration (Response Factor = 1)

Table 1.4 (cont.) Results of TIC for BNA in Soil  
 WA # 3-439 Lackawanna Foundry Site

| Sample # | A16669 |  |   |             | 36              |
|----------|--------|--|---|-------------|-----------------|
| LabFile# | LFS042 |  |   | Con. Factor |                 |
|          | CAS#   | Compound                                     | Q | RT          | Conc.*<br>µg/kg |
| 1        |        | Trimethyl-benzene isomer                     |   | 4.92        | 460             |
| 2        |        | Trimethyl-benzene isomer                     |   | 5.21        | 1300            |
| 3        |        | Trimethyl-benzene isomer                     |   | 5.48        | 550             |
| 4        |        | Benzaldehyde, hydroxy- + alkyl benzene       |   | 5.67        | 2500            |
| 5        |        | Tetramethyl-benzene isomer                   |   | 5.73        | 960             |
| 6        |        | Butoxyethyl acetate                          |   | 5.86        | 420             |
| 7        |        | Tetramethyl-benzene isomer                   |   | 5.91        | 490             |
| 8        |        | Tetramethyl-benzene isomer                   |   | 5.98        | 570             |
| 9        |        | Tetramethyl-benzene isomer                   |   | 6.25        | 610             |
| 10       |        | Tetramethyl-benzene isomer                   |   | 6.28        | 610             |
| 11       |        | Alkyl benzene                                |   | 6.43        | 370             |
| 12       |        | Tetramethyl-benzene isomer + alkyl benzene   |   | 6.57        | 510             |
| 13       |        | Unknown acid                                 |   | 7.08        | 450             |
| 14       |        | Unknown                                      |   | 7.34        | 370             |
| 15       |        | Unknown alcohol                              |   | 7.39        | 420             |
| 16       |        | Unknown                                      |   | 7.97        | 510             |
| 17       |        | Unknown                                      |   | 8.36        | 460             |
| 18       |        | Unknown phthalate                            |   | 11.08       | 500             |
| 19       |        | Unknown                                      |   | 11.67       | 390             |
| 20       |        | Benzene, methylenebis[isocyanato-] + unknown |   | 12.37       | 1400            |

\* Estimated Concentration (Response Factor = 1)

00036

Table 1.4 (cont.) Results of TIC for BNA in Soil  
WA # 3-439 Lackawanna Foundry Site

| Sample # | LabFile#  | A16670                                   | LFS043 | Con. Factor | 1'                            |
|----------|-----------|--|--------|-------------|-------------------------------|
|          | CAS#      | Compound                                 | Q      | RT          | Conc. $\mu\text{g}/\text{kg}$ |
| 1        |           | Trimethyl-benzene isomer                 |        | 4.92        | 2900                          |
| 2        |           | Trimethyl-benzene isomer                 |        | 5.09        | 2100                          |
| 3        |           | Trimethyl-benzene isomer                 |        | 5.21        | 7500                          |
| 4        |           | Trimethyl-benzene isomer                 |        | 5.48        | 4000                          |
| 5        |           | Tetramethylbenzene isomer                |        | 5.73        | 5600                          |
| 6        |           | Methyl-propyl-benzene isomer             |        | 5.83        | 1200                          |
| 7        |           | Tetramethylbenzene isomer                |        | 5.90        | 3400                          |
| 8        |           | Tetramethylbenzene isomer                |        | 5.93        | 2500                          |
| 9        |           | Tetramethylbenzene isomer                |        | 5.97        | 5200                          |
| 10       |           | Tetramethylbenzene isomer                |        | 6.25        | 4600                          |
| 11       |           | Tetramethylbenzene isomer                |        | 6.28        | 4700                          |
| 12       |           | Alkyl benzene                            |        | 6.42        | 1200                          |
| 13       |           | Tetramethylbenzene isomer + unknown      |        | 6.57        | 4600                          |
| 14       |           | Alkyl benzene                            |        | 6.69        | 1400                          |
| 15       |           | Unknown acid                             |        | 7.08        | 2000                          |
| 16       |           | Unknown alcohol                          |        | 7.38        | 2600                          |
| 17       |           | Unknown                                  |        | 8.36        | 1200                          |
| 18       | 2057-49-0 | Pyridine, 4-(3-phenylpropyl-)            | 93     | 10.69       | 6300                          |
| 19       |           | Unknown phthalate                        |        | 11.08       | 2900                          |
| 20       | 101-68-8  | Benzene, 1,1'-methylenebis[4-isocyanato- | 99     | 12.37       | 3700                          |

\* Estimated Concentration (Response Factor = 1)

Table 1.4 (cont.) Results of TIC for BNA in Soil  
 WA # 3-439 Lackawanna Foundry Site

| Sample # | A16672 |                         |   | Con. Factor | 200             |
|----------|--------|-------------------------|---|-------------|-----------------|
| LabFile# | LFS045 |                         |   |             |                 |
|          | CAS#   | Compound                | Q | RT          | Conc.*<br>µg/kg |
| 1        |        | Unknown PAH isomer      |   | 11.70       | 980             |
| 2        |        | H-Benzo fluorene isomer |   | 13.16       | 920             |
| 3        |        | Unknown aldehyde        |   | 16.54       | 1200            |
| 4        |        | Unknown alkane          |   | 17.06       | 1700            |
| 5        |        | Unknown PAH isomer      |   | 18.15       | 2000            |
| 6        |        | Unknown aldehyde        |   | 18.94       | 1200            |
| 7        |        | Unknown alkane          |   | 19.66       | 1200            |
| 8        |        | Unknown aldehyde        |   | 22.19       | 1400            |
| 9        |        |                         |   |             |                 |
| 10       |        |                         |   |             |                 |
| 11       |        |                         |   |             |                 |
| 12       |        |                         |   |             |                 |
| 13       |        |                         |   |             |                 |
| 14       |        |                         |   |             |                 |
| 15       |        |                         |   |             |                 |
| 16       |        |                         |   |             |                 |
| 17       |        |                         |   |             |                 |
| 18       |        |                         |   |             |                 |
| 19       |        |                         |   |             |                 |
| 20       |        |                         |   |             |                 |

\* Estimated Concentration (Response Factor = 1)

0005

Table 1.4 (cont.) Results of TIC for BNA in Soil  
 WA # 3-439 Lackawanna Foundry Site

| Sample #<br>LabFile# | A16673<br>LFS048 | Con. Factor                                  |   |       | Conc.<br>µg/kg |
|----------------------|------------------|--|---|-------|----------------|
|                      | CAS#             | Compound                                     | Q | RT    |                |
| 1                    |                  | Unknown                                      |   | 12.31 | 1700           |
| 2                    |                  | Benzene, methylenebis[isocyanato]- + unknown |   | 12.37 | 1800           |
| 3                    |                  | Unknown PAH isomer                           |   | 13.15 | 2300           |
| 4                    |                  | Unknown alkane                               |   | 17.05 | 1800           |
| 5                    |                  | Unknown PAH isomer                           |   | 18.15 | 4800           |
| 6                    |                  |  |   |       |                |
| 7                    |                  |  |   |       |                |
| 8                    |                  |  |   |       |                |
| 9                    |                  |  |   |       |                |
| 10                   |                  |  |   |       |                |
| 11                   |                  |  |   |       |                |
| 12                   |                  |  |   |       |                |
| 13                   |                  |  |   |       |                |
| 14                   |                  |  |   |       |                |
| 15                   |                  |  |   |       |                |
| 16                   |                  |  |   |       |                |
| 17                   |                  |  |   |       |                |
| 18                   |                  |  |   |       |                |
| 19                   |                  |  |   |       |                |
| 20                   |                  |  |   |       |                |

\* Estimated Concentration (Response Factor = 1)



Table 1.4 (cont.) Results of TIC for BNA in Soil  
 WA # 3-439 Lackawanna Foundry Site

| Sample # | LabFile# | A16674   | LFS049 | Con. Factor | 220             |
|----------|----------|--|--------|-------------|-----------------|
|          | CAS#     | Compound   | Q      | RT          | Conc.*<br>µg/kg |
| 1        |          | Trimethyl-benzene isomer                           |        | 5.21        | 1100            |
| 2        |          | Benzaldehyde, hydroxy + tetramethyl-benzene isomer |        | 5.68        | 1100            |
| 3        |          | Unknown alkane                                     |        | 9.44        | 910             |
| 4        |          | Unknown  |        | 12.31       | 2000            |
| 5        | 101-68-8 | Benzene, 1,1'-methylenebis[4-isocyanato]-          | 97     | 12.37       | 2100            |
| 6        |          | Unknown  |        | 16.55       | 880             |
| 7        |          | Unknown alkane                                     |        | 17.06       | 2400            |
| 8        |          | Unknown alkane                                     |        | 19.67       | 3300            |
| 9        |          |  |        |             |                 |
| 10       |          |  |        |             |                 |
| 11       |          |  |        |             |                 |
| 12       |          |  |        |             |                 |
| 13       |          |  |        |             |                 |
| 14       |          |  |        |             |                 |
| 15       |          |  |        |             |                 |
| 16       |          |  |        |             |                 |
| 17       |          |  |        |             |                 |
| 18       |          |  |        |             |                 |
| 19       |          |  |        |             |                 |
| 20       |          |  |        |             |                 |

\* Estimated Concentration (Response Factor = 1)

03040

Table 1.4 (cont.) Results of TIC for BNA in Soil  
 WA # 3-439 Lackawanna Foundry Site

Sample #  
 LabFile#

A16677  
 LFS052

Con. Factor

4

|    | CAS# | Compound                            | Q | RT    | Conc.<br>µg/kg |
|----|------|-------------------------------------|---|-------|----------------|
| 1  |      | Unknown aldehyde                    |   | 16.55 | 2600           |
| 2  |      | Unknown alkane                      |   | 17.06 | 3300           |
| 3  |      | Unknown aldehyde                    |   | 18.95 | 2000           |
| 4  |      | Unknown alkane                      |   | 19.67 | 3200           |
| 5  |      | Unknown alkane + alkene/cycloalkene |   | 22.21 | 3100           |
| 6  |      |                                     |   |       |                |
| 7  |      |                                     |   |       |                |
| 8  |      |                                     |   |       |                |
| 9  |      |                                     |   |       |                |
| 10 |      |                                     |   |       |                |
| 11 |      |                                     |   |       |                |
| 12 |      |                                     |   |       |                |
| 13 |      |                                     |   |       |                |
| 14 |      |                                     |   |       |                |
| 15 |      |                                     |   |       |                |
| 16 |      |                                     |   |       |                |
| 17 |      |                                     |   |       |                |
| 18 |      |                                     |   |       |                |
| 19 |      |                                     |   |       |                |
| 20 |      |                                     |   |       |                |

\* Estimated Concentration (Response Factor = 1)

Table 1.4 (cont.) Results of TIC for BNA in Soil  
 WA # 3-439 Lackawanna Foundry Site

| Sample # | LabFile# | A16678                   | LFS053 | Con. Factor | 44              |
|----------|----------|--------------------------|--------|-------------|-----------------|
|          | CAS#     | Compound                 | Q      | RT          | Conc.*<br>µg/kg |
| 1        |          | Unknown                  |        | 3.90        | 280             |
| 2        |          | Unknown                  |        | 3.95        | 180             |
| 3        |          | Unknown alkane           |        | 9.44        | 280             |
| 4        |          | Unknown alkane + unknown |        | 9.73        | 310             |
| 5        |          | Unknown alkane           |        | 10.58       | 280             |
| 6        |          | Unknown alkane           |        | 10.64       | 180             |
| 7        |          | Unknown alkane           |        | 11.11       | 230             |
| 8        |          | Unknown alkane           |        | 11.60       | 240             |
| 9        |          | Unknown PAH isomer       |        | 18.16       | 420             |
| 10       |          |                          |        |             |                 |
| 11       |          |                          |        |             |                 |
| 12       |          |                          |        |             |                 |
| 13       |          |                          |        |             |                 |
| 14       |          |                          |        |             |                 |
| 15       |          |                          |        |             |                 |
| 16       |          |                          |        |             |                 |
| 17       |          |                          |        |             |                 |
| 18       |          |                          |        |             |                 |
| 19       |          |                          |        |             |                 |
| 20       |          |                          |        |             |                 |

\* Estimated Concentration (Response Factor = 1)

00042

Table 1.4 (cont.) Results of TIC for BNA in Soil  
WA # 3-439 Lackawanna Foundry Site

| Sample # | LabFile# | CAS#        | Compound  | Con. Factor | Q  | RT    | Conc.<br>µg/kg |
|----------|----------|-------------|---|-------------|----|-------|----------------|
| 1        |          |             | Trimethyl-benzene isomer                                |             |    | 4.92  | 2100           |
| 2        |          |             | Trimethyl-benzene isomer                                |             |    | 5.21  | 4500           |
| 3        |          |             | Tetramethyl-benzene isomer                              |             |    | 5.98  | 2800           |
| 4        |          |             | Tetramethyl-benzene isomer                              |             |    | 6.25  | 3200           |
| 5        |          |             | Tetramethyl-benzene isomer                              |             |    | 6.28  | 2700           |
| 6        |          |             | Dihydro methylindene + tetramethyl-benzene isomer       |             |    | 6.57  | 3200           |
| 7        |          | 000090-01-7 | Salicyl Alcohol   |             | 94 | 7.39  | 4800           |
| 8        |          |             | Tetrachloro-benzene isomer + unknown                    |             |    | 8.36  | 2100           |
| 9        |          | 2057-49-0   | Pyridine, 4-(3-phenylpropyl)                            |             | 94 | 10.69 | 9300           |
| 10       |          |             | Trichloro-biphenyl isomer + unknown                     |             |    | 10.82 | 2700           |
| 11       |          |             | Trichloro-biphenyl isomer + unknown                     |             |    | 11.05 | 2100           |
| 12       |          |             | Trichloro-biphenyl isomer                               |             |    | 11.31 | 3600           |
| 13       |          |             | Trichloro-biphenyl isomer                               |             |    | 11.33 | 3800           |
| 14       |          |             | Tetrachloro-biphenyl isomer + trichloro-biphenyl isomer |             |    | 11.67 | 2800           |
| 15       |          |             | Tetrachloro-biphenyl isomer + trichloro-biphenyl isomer |             |    | 11.72 | 2100           |
| 16       |          |             | Tetrachloro-biphenyl isomer + trichloro-biphenyl isomer |             |    | 11.91 | 2600           |
| 17       |          |             | Tetrachloro-biphenyl isomer + trichloro-biphenyl isomer |             |    | 11.95 | 2400           |
| 18       |          |             | Tetrachloro-biphenyl isomer                             |             |    | 12.29 | 2200           |
| 19       |          |             | Tetrachloro-biphenyl isomer                             |             |    | 12.34 | 2700           |
| 20       |          |             | Tetrachloro-biphenyl isomer                             |             |    | 12.52 | 2300           |

\* Estimated Concentration (Response Factor = 1)

## QA/QC for Metals

### Results of the QC Standard Analysis for Metals (Soil)

QC standards QC-7x100, QC-21x100, ERA-434, TMAA #1, TMAA #2 and TMWS were used to check the accuracy of the calibration curve. The percent recoveries, listed in Table 2.1, ranged from 90 to 115 and all twenty recovered concentrations for which 95% confidence limits are available were within the 95% confidence limits. The 95% confidence limits for nineteen values are not available.

### Results of the MS/MSD Analysis for Metals in Soil

Samples C 16655, C 16661 and C 16669 were chosen for the matrix spike/matrix spike duplicate analysis (MS/MSD). The percent recoveries, listed in Table 2.2, ranged from 0 (zero) to 219. Sixty-four out of ninety-two reported values were within the acceptable QC limits. Ten values were not calculated because the concentration of analyte in the sample was greater than four times the concentration spiked. The relative percent differences, also listed in Table 2.2, ranged from 0 (zero) to 104. Thirty-five out of forty-four values were within the acceptable QC limits. Seven values were not calculated either because the spike recoveries were not calculated or none of the spiked analyte was recovered.

### Results of the Blank Spike Analysis for Metals in Soil

The results of the blank spike analysis are reported in Table 2.3. The percent recoveries ranged from 92 to 109 and all twenty-three recovered values were within the acceptable QC limits.

Table 2.1 Results of the QC Standard Analysis for Metals (Soil)  
WA # 3-439 Lackawanna Foundry Site

| Metal     | Date Analyzed | Quality Control Standard | Conc. Rec µg/L | Certified Value µg/L | 95 % Confidence Interval µg/L | % Rec |
|-----------|---------------|--------------------------|----------------|----------------------|-------------------------------|-------|
| Aluminum  | 04/20/99      | QC-7 x100                | 998            | 1000                 | NA                            | 100   |
|           | 04/20/99      | ERA-434                  | 703            | 647                  | 531-763                       | 109   |
| Antimony  | 04/20/99      | QC-21 x100               | 981            | 1000                 | NA                            | 98    |
| Arsenic   | 04/19/99      | TMAA #1                  | 49.7           | 50                   | 41.9-55.9                     | 99    |
|           | 04/20/99      | TMAA #1                  | 52.3           | 50                   | 41.9-55.9                     | 105   |
| Barium    | 04/20/99      | QC-7 x100                | 964            | 1000                 | NA                            | 96    |
|           | 04/20/99      | ERA-434                  | 723            | 735                  | 603 - 867                     | 98    |
| Beryllium | 04/20/99      | QC-21 x100               | 1048           | 1000                 | NA                            | 105   |
|           | 04/20/99      | ERA-434                  | 86.2           | 82                   | 68 - 97                       | 105   |
| Cadmium   | 04/20/99      | QC-21 x100               | 1045           | 1000                 | NA                            | 105   |
|           | 04/20/99      | ERA-434                  | 78.3           | 77                   | 63 - 90                       | 102   |
| Calcium   | 04/20/99      | QC-21 x100               | 1061           | 1000                 | NA                            | 106   |
| Chromium  | 04/20/99      | QC-21 x100               | 1061           | 1000                 | NA                            | 106   |
|           | 04/20/99      | ERA-434                  | 113            | 106                  | 87 - 125                      | 107   |
| Cobalt    | 04/20/99      | QC-21 x100               | 1074           | 1000                 | NA                            | 107   |
|           | 04/20/99      | ERA-434                  | 94.1           | 88                   | 72 - 104                      | 107   |
| Copper    | 04/20/99      | QC-21 x100               | 1019           | 1000                 | NA                            | 102   |
|           | 04/20/99      | ERA-434                  | 151            | 147                  | 121 - 173                     | 103   |
| Iron      | 04/20/99      | QC-21 x100               | 1082           | 1000                 | NA                            | 108   |
|           | 04/20/99      | ERA-434                  | 223            | 206                  | 169 - 243                     | 108   |
| Lead      | 04/20/99      | QC-21 x100               | 1067           | 1000                 | NA                            | 107   |
|           | 04/20/99      | ERA-434                  | 108            | 94                   | 77 - 111                      | 115   |
| Magnesium | 04/20/99      | QC-21 x100               | 1005           | 1000                 | NA                            | 100   |
| Manganese | 04/20/99      | QC-21 x100               | 1053           | 1000                 | NA                            | 105   |
|           | 04/20/99      | ERA-434                  | 244            | 235                  | 193 - 277                     | 104   |
| Mercury   | 04/19/99      | TMWS                     | 2.8            | 2.9                  | 2.13 - 3.53                   | 97    |
| Nickel    | 04/20/99      | QC-21 x100               | 1099           | 1000                 | NA                            | 110   |
|           | 04/20/99      | ERA-434                  | 124            | 112                  | 92 - 132                      | 111   |
| Potassium | 04/20/99      | QC-7 x100                | 8988           | 10000                | NA                            | 90    |
| Selenium  | 04/16/99      | TMAA #1                  | 54.76          | 50                   | 39.4-57.4                     | 110   |
|           | 04/26/99      | TMAA #1                  | 49.16          | 50                   | 39.4-57.4                     | 98    |
| Silver    | 04/20/99      | QC-7 x100                | 997            | 1000                 | NA                            | 100   |
|           | 04/20/99      | ERA-434                  | 92.3           | 88                   | 72 - 104                      | 105   |
| Sodium    | 04/20/99      | QC-7 x100                | 952            | 1000                 | NA                            | 95    |
| Thallium  | 04/16/989     | TMAA #2                  | 51.1           | 50                   | 39.9-57.97                    | 102   |
| Vanadium  | 04/20/99      | QC-21 x100               | 1046           | 1000                 | NA                            | 105   |
|           | 04/20/99      | ERA-434                  | 123            | 118                  | 97 - 139                      | 104   |
| Zinc      | 04/20/99      | QC-21 x100               | 1071           | 1000                 | NA                            | 107   |
|           | 04/20/99      | ERA-434                  | 280            | 265                  | 217 - 313                     | 106   |

000-25

Table 2.2 Results of the MS/MSD Analysis for Metals in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 Results Based on Dry Weight

Sample ID: C16655

| Metal     | Sample Conc mg/kg | MS Spike Added mg/kg | MS Conc mg/kg | MS % Rec | MSD Spike Added mg/kg | MSD Conc mg/kg | MSD % Rec | RPD   | Recommended QC Limits %Rec | RPD |
|-----------|-------------------|----------------------|---------------|----------|-----------------------|----------------|-----------|-------|----------------------------|-----|
| Antimony  | U                 | 50.6                 | U             | 0 *      | 50.6                  | U              | 0 *       | NC    | 75-125                     | 20  |
| Arsenic   | 15.1              | 5.89                 | 20.1          | 85       | 5.83                  | 22.5           | 127 *     | 40 *  | 75-125                     | 20  |
| Banum     | 129               | 50.6                 | 171           | 83       | 50.6                  | 172            | 85        | 2     | 75-125                     | 20  |
| Beryllium | 0.853             | 50.6                 | 51.9          | 101      | 50.6                  | 52.9           | 103       | 2     | 75-125                     | 20  |
| Cadmium   | 2.44              | 50.6                 | 51.7          | 97       | 50.6                  | 52.3           | 98        | 1     | 75-125                     | 20  |
| Chromium  | 29                | 50.6                 | 76.5          | 94       | 50.6                  | 78.5           | 98        | 4     | 75-125                     | 20  |
| Cobalt    | 13.9              | 50.6                 | 61.8          | 95       | 50.6                  | 62.9           | 97        | 2     | 75-125                     | 20  |
| Copper    | 195               | 50.6                 | 205           | 20 *     | 50.6                  | 227            | 63 *      | 104 * | 75-125                     | 20  |
| Lead      | 179               | 50.6                 | 185           | 12 *     | 50.6                  | 198            | 38 *      | 104 * | 75-125                     | 20  |
| Manganese | 700               | 50.6                 | 722           | NC       | 50.6                  | 691            | NC        | NC    | 75-125                     | 20  |
| Mercury   | 0.111             | 0.438                | 0.525         | 95       | 0.438                 | 0.525          | 95        | 0     | 75-125                     | 20  |
| Nickel    | 41                | 50.6                 | 87.7          | 92       | 50.6                  | 87.5           | 92        | 0     | 75-125                     | 20  |
| Selenium  | 0.578             | 5.89                 | 3.066         | 42 *     | 5.83                  | 3.74           | 54 *      | 25 *  | 75-125                     | 20  |
| Silver    | U                 | 50.6                 | 48.1          | 95       | 50.6                  | 48.3           | 95        | 0     | 75-125                     | 20  |
| Thallium  | U                 | 5.89                 | 4.84          | 82       | 5.83                  | 4.67           | 80        | 3     | 75-125                     | 20  |
| Vanadium  | 23.4              | 50.6                 | 73.5          | 99       | 50.6                  | 72.4           | 97        | 2     | 75-125                     | 20  |
| Zinc      | 340               | 50.6                 | 301           | NC       | 50.6                  | 319            | NC        | NC    | 75-125                     | 20  |

00046

Table 2.2 (Cont) Results of the MS/MSD Analysis for Metals in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 Results Based on Dry Weight

| Sample ID: C16661 |                   |                   |               |          |                   |                |           |      |                       |     |  |
|-------------------|-------------------|-------------------|---------------|----------|-------------------|----------------|-----------|------|-----------------------|-----|--|
| Metal             | Sample Conc mg/kg | MS                |               |          | MSD               |                |           | RPD  | Recommended QC Limits |     |  |
|                   |                   | Spike Added mg/kg | MS Conc mg/kg | MS % Rec | Spike Added mg/kg | MSD Conc mg/kg | MSD % Rec |      | %Rec                  | RPD |  |
| Antimony          | U                 | 47.9              | 30.6          | 64 *     | 48.3              | 27.8           | 58 *      | 10   | 75-125                | 20  |  |
| Arsenic           | 1.8               | 4.64              | 5.87          | 88       | 4.60              | 5.7            | 85        | 3    | 75-125                | 20  |  |
| Barium            | 14.2              | 47.9              | 58.4          | 92       | 48.3              | 57.7           | 90        | 2    | 75-125                | 20  |  |
| Beryllium         | U                 | 47.9              | 50.1          | 105      | 48.3              | 50.6           | 105       | 0    | 75-125                | 20  |  |
| Cadmium           | 4.15              | 47.9              | 49.2          | 94       | 48.3              | 49.3           | 93        | 1    | 75-125                | 20  |  |
| Chromium          | 12.9              | 47.9              | 58.3          | 95       | 48.3              | 58.1           | 94        | 1    | 75-125                | 20  |  |
| Cobalt            | 1.69              | 47.9              | 50.3          | 101      | 48.3              | 50.7           | 101       | 0    | 75-125                | 20  |  |
| Copper            | 32.8              | 47.9              | 77.7          | 94       | 48.3              | 82.8           | 104       | 10   | 75-125                | 20  |  |
| Lead              | 47.7              | 47.9              | 89.9          | 88       | 48.3              | 87.8           | 83        | 6    | 75-125                | 20  |  |
| Manganese         | 164               | 47.9              | 169           | 10 *     | 48.3              | 148            | 0 *       | NC   | 75-125                | 20  |  |
| Mercury           | 0.038             | 0.354             | 0.39          | 99       | 0.354             | 0.39           | 99        | 0    | 75-125                | 20  |  |
| Nickel            | 13.2              | 47.9              | 59.3          | 96       | 48.3              | 60             | 97        | 1    | 75-125                | 20  |  |
| Selenium          | U                 | 4.64              | 3.88          | 84       | 4.60              | 3.79           | 82        | 2    | 75-125                | 20  |  |
| Silver            | U                 | 47.9              | 46.4          | 97       | 48.3              | 46.6           | 96        | 0    | 75-125                | 20  |  |
| Thallium          | U                 | 4.64              | 3.22          | 69 *     | 4.60              | 3.53           | 77        | 10   | 75-125                | 20  |  |
| Vanadium          | 2.74              | 47.9              | 50.9          | 101      | 48.3              | 51.6           | 101       | 1    | 75-125                | 20  |  |
| Zinc              | 81                | 47.9              | 102           | 44 *     | 48.3              | 93.7           | 26 *      | 50 * | 75-125                | 20  |  |



Table 2.2 (Cont) Results of the MS/MSD Analysis for Metals in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 Results Based on Dry Weight

| Sample ID: C16669 |                   |                      |               |          |                       |                |           |      |                       |     |  |
|-------------------|-------------------|----------------------|---------------|----------|-----------------------|----------------|-----------|------|-----------------------|-----|--|
| Metal             | Sample Conc mg/kg | MS Spike Added mg/kg | MS Conc mg/kg | MS % Rec | MSD Spike Added mg/kg | MSD Conc mg/kg | MSD % Rec | RPD  | Recommended QC Limits |     |  |
|                   |                   |                      |               |          |                       |                |           |      | %Rec                  | RPD |  |
| Antimony          | U                 | 44.4                 | 20.5          | 46 *     | 44.8                  | 17.3           | 39 *      | 18   | 75-125                | 20  |  |
| Arsenic           | 6.74              | 4.72                 | 13.6          | 145 *    | 4.76                  | 10.7           | 85        | 52 * | 75-125                | 20  |  |
| Barium            | 28.7              | 44.4                 | 68.4          | 89       | 44.8                  | 75.8           | 105       | 16   | 75-125                | 20  |  |
| Beryllium         | U                 | 44.4                 | 45.6          | 103      | 44.8                  | 47.4           | 106       | 3    | 75-125                | 20  |  |
| Cadmium           | 0.592             | 44.4                 | 44.8          | 100      | 44.8                  | 46.2           | 102       | 2    | 75-125                | 20  |  |
| Chromium          | 96.9              | 44.4                 | 194           | 219 *    | 44.8                  | 159            | 139 *     | 45 * | 75-125                | 20  |  |
| Cobalt            | 4.42              | 44.4                 | 56.6          | 118      | 44.8                  | 52.9           | 108       | 8    | 75-125                | 20  |  |
| Copper            | 297               | 44.4                 | 402           | NC       | 44.8                  | 284            | NC        | NC   | 75-125                | 20  |  |
| Lead              | 141               | 44.4                 | 185           | 99       | 44.8                  | 177            | 80        | 21 * | 75-125                | 20  |  |
| Manganese         | 763               | 44.4                 | 736           | NC       | 44.8                  | 899            | NC        | NC   | 75-125                | 20  |  |
| Mercury           | U                 | 0.343                | 0.361         | 105      | 0.349                 | 0.367          | 105       | 0    | 75-125                | 20  |  |
| Nickel            | 57.1              | 44.4                 | 134           | 173 *    | 44.8                  | 118            | 136 *     | 24 * | 75-125                | 20  |  |
| Selenium          | U                 | 4.72                 | 2.03          | 43 *     | 4.76                  | 2.16           | 45 *      | 5    | 75-125                | 20  |  |
| Silver            | U                 | 44.4                 | 39            | 88       | 44.8                  | 42.5           | 95        | 8    | 75-125                | 20  |  |
| Thallium          | U                 | 4.72                 | 0.925         | 20 *     | 4.76                  | 0.99           | 21 *      | 6    | 75-125                | 20  |  |
| Vanadium          | 13.7              | 44.4                 | 69.5          | 126 *    | 44.8                  | 63.5           | 111       | 12   | 75-125                | 20  |  |
| Zinc              | 308               | 44.4                 | 381           | NC       | 44.8                  | 413            | NC        | NC   | 75-125                | 20  |  |

00045

**Table 2.3 Results of the Blank Spike Analysis for Metals in Soil  
WA # 3-439 Lackawanna Foundry Site**

| Metal     | Spiked Conc.<br>mg/kg | Sand Blank Conc.<br>mg/kg | Rec Conc.<br>mg/kg | % Rec | Recommended QC Limits<br>%Rec |
|-----------|-----------------------|---------------------------|--------------------|-------|-------------------------------|
| Aluminum  | 400                   | 37.1                      | 470                | 108   | 75-125                        |
| Antimony  | 50.0                  | U                         | 45.9               | 92    | 75-125                        |
| Arsenic   | 5.00                  | U                         | 4.88               | 98    | 75-125                        |
| Barium    | 50.0                  | U                         | 49                 | 98    | 75-125                        |
| Beryllium | 50.0                  | U                         | 53.1               | 106   | 75-125                        |
| Cadmium   | 50.0                  | U                         | 51.2               | 102   | 75-125                        |
| Calcium   | 400                   | U                         | 422                | 106   | 75-125                        |
| Chromium  | 50.0                  | U                         | 52.8               | 106   | 75-125                        |
| Cobalt    | 50.0                  | U                         | 53.2               | 106   | 75-125                        |
| Copper    | 50.0                  | U                         | 50.8               | 102   | 75-125                        |
| Iron      | 400                   | 14.7                      | 439                | 106   | 75-125                        |
| Lead      | 50.0                  | U                         | 53.3               | 107   | 75-125                        |
| Magnesium | 400                   | U                         | 408                | 102   | 75-125                        |
| Manganese | 50.0                  | U                         | 51.3               | 103   | 75-125                        |
| Mercury   | 0.400                 | U                         | 0.4                | 100   | 75-125                        |
| Nickel    | 50.0                  | U                         | 54.5               | 109   | 75-125                        |
| Potassium | 800                   | U                         | 756                | 95    | 75-125                        |
| Selenium  | 5.00                  | U                         | 5.4                | 108   | 75-125                        |
| Silver    | 50.0                  | U                         | 49.5               | 99    | 75-125                        |
| Sodium    | 400                   | U                         | 393                | 98    | 75-125                        |
| Thallium  | 5.00                  | U                         | 4.98               | 100   | 75-125                        |
| Vanadium  | 50.0                  | U                         | 51.9               | 104   | 75-125                        |
| Zinc      | 50.0                  | U                         | 52.9               | 106   | 75-125                        |

## QA/QC for Cyanide

### Results of the Blank Spike Analysis for Cyanide in Soil

Four blanks were spiked with cyanide. The percent recoveries, listed in Table 2.4, ranged from 99 to 104 and all four values were within the acceptable QC limits.

### Results of the Matrix Spike Analysis for Cyanide in Soil

Samples D 16651 and D 16670 were chosen for the matrix spike analysis (MS). The percent recoveries, listed in Table 2.5, were 75 and 83 and both values were within the acceptable QC limits.

### Results of the Duplicate Analysis for Cyanide in Soil

Samples D 16651 and D 16670 were chosen for the duplicate analysis. The relative percent differences, listed in Table 2.6, were not calculated because cyanide was not detected in either analysis.

Table 2.4 Results of the Blank Spike Analysis for Cyanide  
in Soil

WA # 3-439 Lackawanna Foundry Site

| Analyte | Spiked<br>Conc<br>mg/kg | Rec<br>Conc<br>mg/kg | % Rec | Recommended<br>Limit |
|---------|-------------------------|----------------------|-------|----------------------|
| Cyanide | 4.87                    | 4.86                 | 100   | 90-110               |
| Cyanide | 24.4                    | 24.7                 | 101   | 90-110               |
| Cyanide | 4.87                    | 4.84                 | 99    | 90-110               |
| Cyanide | 24.4                    | 25.3                 | 104   | 90-110               |

00051

Table 2.5 Results of the Matrix Spike Analysis for Cyanide in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 (based on dry weight)

| Analyte | Sample ID | Sample<br>Conc<br>mg/kg | Spike<br>Conc<br>mg/kg | Rec<br>Conc<br>mg/kg | %<br>Rec | QC<br>Limits |
|---------|-----------|-------------------------|------------------------|----------------------|----------|--------------|
| Cyanide | D 16651   | U                       | 5.18                   | 4.3                  | 83       | 75-125       |
| Cyanide | D 16670   | U                       | 5.36                   | 4.0                  | 75       | 75-125       |

06052

Table 2.6 Results of the Duplicate Analysis  
for Cyanide in Soil  
WA # 3-439 Lackawanna Foundry Site  
(based on dry weight)

| Analyte | Sample ID | Initial<br>Analysis<br>mg/kg | Duplicate<br>Analysis<br>mg/kg | RPD |
|---------|-----------|------------------------------|--------------------------------|-----|
| Cyanide | D 16651   | U                            | U                              | NC  |
| Cyanide | D 16670   | U                            | U                              | NC  |

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## QA/QC for BNA

### Results of the Internal Standard Areas for BNA in Soil

The internal standard areas (for 1,4-dichlorobenzene-d<sub>4</sub>, naphthalene-d<sub>8</sub>, acenaphthene-d<sub>10</sub>, phenanthrene-d<sub>10</sub>, chrysene-d<sub>12</sub>, perylene-d<sub>12</sub>) are listed in Table 2.7. All two hundred and forty-six areas were within the acceptable QC limits.

### Results of the Surrogate Recoveries for BNA in Soil

Before extraction, each sample was spiked with a six component mixture of CLP surrogate standards consisting of nitrobenzene-d<sub>5</sub>, 2-fluorobiphenyl, terphenyl-d<sub>14</sub>, phenol-d<sub>5</sub>, 2-fluorophenol, and 2,4,6-tribromophenol. The surrogate percent recoveries, listed in Table 2.8, ranged from 51 to 129. Two hundred and forty-five values out of two hundred and forty-six were within the acceptable QC limits.

### Results of the MS/MSD Analysis for BNA in Soil

Samples A 16552, A 16660 and A 16672 were chosen for the matrix spike/matrix spike duplicate (MS/MSD) analyses. The percent recoveries, ranging from 63 to 134, are listed in Table 2.9. Fifty-four calculated values out of sixty-four were within the acceptable QC limits. Two values were not calculated because the concentration spiked was less than one fourth the concentration of analyte in the sample. The relative percent differences, also listed in Table 2.9, ranged from 1 to 26. All thirty-two calculated values were within the acceptable QC limits. One value was not calculated because the concentration spiked was less than one fourth the concentration of analyte in the sample.

Table 2.7 Results of the Internal Standard Areas for BNA in Soil  
WA # 3-439 Lackawanna Foundry Site

| Sample No      | File ID  | IS 1  | IS 2   | IS 3   | IS 4   | IS 5   | IS 6   |
|----------------|----------|-------|--------|--------|--------|--------|--------|
| Cal Check Area | LFS001.D | 48688 | 179730 | 97785  | 167866 | 177152 | 186601 |
| SBLK041599     | LFS002.D | 38510 | 140571 | 78553  | 148212 | 150207 | 125993 |
| A16650         | LFS003.D | 40571 | 153235 | 85090  | 163211 | 173252 | 148749 |
| A16651         | LFS004.D | 41330 | 160993 | 91734  | 175743 | 180938 | 151966 |
| A16652         | LFS005.D | 40603 | 154160 | 85563  | 162650 | 169517 | 144850 |
| A16652 MS      | LFS006.D | 43881 | 164833 | 92390  | 173001 | 175900 | 150056 |
| A16652 MSD     | LFS007.D | 44986 | 165422 | 92428  | 172362 | 178812 | 152838 |
| A16653         | LFS008.D | 43557 | 167418 | 93790  | 175400 | 183577 | 156110 |
| A16654         | LFS009.D | 39551 | 147585 | 82831  | 157607 | 162902 | 138293 |
| A16655         | LFS010.D | 44828 | 164660 | 92090  | 171161 | 177732 | 151656 |
| A16656         | LFS012.D | 45814 | 172544 | 94436  | 174912 | 178680 | 150439 |
| A16651 10x     | LFS013.D | 43165 | 166651 | 89943  | 172543 | 168251 | 138359 |
| Cal Check Area | LFS018.D | 44996 | 167440 | 91487  | 152765 | 171184 | 194601 |
| A16655 5x      | LFS021.D | 47925 | 181511 | 102192 | 199801 | 226495 | 203962 |
| Cal Check Area | LFS022.D | 41707 | 157767 | 85588  | 147737 | 183361 | 230722 |
| SBLK041999     | LFS023.D | 35683 | 129660 | 77135  | 151823 | 185395 | 183296 |
| A16657         | LFS024.D | 40423 | 150208 | 86919  | 176546 | 244298 | 286758 |
| A16658         | LFS025.D | 44152 | 165587 | 96299  | 184119 | 260997 | 288930 |
| A16659         | LFS026.D | 49458 | 187128 | 105200 | 205351 | 272397 | 303064 |
| A16660         | LFS027.D | 49412 | 183395 | 101781 | 200160 | 256189 | 283017 |
| A16660 MS      | LFS028.D | 46718 | 175020 | 98247  | 193740 | 246912 | 262896 |
| A16660 MSD     | LFS029.D | 43718 | 165428 | 90664  | 180347 | 231438 | 246412 |
| A16661         | LFS030.D | 58684 | 223720 | 120367 | 235952 | 286248 | 300780 |
| A16662         | LFS031.D | 52142 | 205487 | 113587 | 228180 | 283717 | 305523 |
| A16663         | LFS032.D | 49650 | 189790 | 105506 | 203526 | 268843 | 293669 |
| A16664         | LFS033.D | 53371 | 195883 | 109444 | 210447 | 257593 | 262743 |
| A16665         | LFS034.D | 48865 | 190182 | 107469 | 206609 | 247928 | 244841 |
| A16666         | LFS035.D | 46744 | 172641 | 95122  | 186850 | 224118 | 215563 |
| A16667         | LFS036.D | 50906 | 187711 | 104538 | 200840 | 250667 | 261343 |
| A16668         | LFS037.D | 45448 | 172680 | 97365  | 187800 | 230510 | 230929 |
| Cal Check Area | LFS040.D | 63317 | 245039 | 131024 | 221397 | 257165 | 289256 |
| SBLK042099     | LFS041.D | 57662 | 210169 | 119222 | 221156 | 250826 | 238932 |
| A16669         | LFS042.D | 61237 | 231646 | 131722 | 250037 | 288580 | 267026 |
| A16670         | LFS043.D | 61347 | 231871 | 132611 | 252844 | 285675 | 269351 |
| A16671         | LFS044.D | 64388 | 240847 | 137526 | 259184 | 292471 | 275487 |
| A16672         | LFS045.D | 68500 | 257101 | 140215 | 269153 | 297219 | 270768 |
| A16672 MS      | LFS046.D | 70715 | 265245 | 149403 | 282760 | 302277 | 277111 |
| A16672 MSD     | LFS047.D | 66152 | 245321 | 136489 | 261503 | 286018 | 256846 |
| A16673         | LFS048.D | 61738 | 231114 | 135262 | 255710 | 281387 | 256022 |
| A16674         | LFS049.D | 64391 | 240974 | 138772 | 261573 | 287460 | 261662 |
| A16675         | LFS050.D | 64370 | 240735 | 135554 | 260538 | 287491 | 256002 |
| A16676         | LFS051.D | 61034 | 230354 | 127683 | 241923 | 264551 | 235349 |
| A16677         | LFS052.D | 64642 | 240201 | 138026 | 262796 | 277436 | 249518 |
| A16678         | LFS053.D | 66386 | 250243 | 141199 | 270854 | 286069 | 255431 |
| A16679         | LFS054.D | 67843 | 260200 | 147376 | 279196 | 297023 | 266294 |

IS 1 = d4-Dichlorobenzene  
 IS 2 = d8-Naphthalene  
 IS 3 = d10-Acenaphthene  
 IS 4 = d10-Phenanthrene  
 IS 5 = d12-Chrysene  
 IS 6 = d12-Perylene

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Table 2.8 Results of the Surrogate Recoveries for BNA in Soil  
WA # 3-439 Lackawanna Foundry Site

| Sample No  | File ID  | Surr. 1 | Surr. 2 | Surr. 3 | Surr. 4 | Surr. 5 | Surr. 6 |
|------------|----------|---------|---------|---------|---------|---------|---------|
| SBLK041599 | LFS002.D | 81      | 90      | 89      | 94      | 81      | 108     |
| A16650     | LFS003.D | 70      | 78      | 67      | 82      | 75      | 90      |
| A16651     | LFS004.D | 63      | 74      | 70      | 78      | 89      | 103     |
| A16652     | LFS005.D | 51      | 67      | 63      | 74      | 85      | 114     |
| A16652 MS  | LFS006.D | 97      | 112     | 105     | 116     | 105     | 129     |
| A16652 MSD | LFS007.D | 83      | 95      | 89      | 101     | 104     | 118     |
| A16653     | LFS008.D | 81      | 93      | 90      | 94      | 77      | 100     |
| A16654     | LFS009.D | 91      | 98      | 98      | 108     | 88      | 111     |
| A16655     | LFS010.D | 66      | 77      | 73      | 86      | 95      | 105     |
| A16656     | LFS012.D | 59      | 72      | 67      | 80      | 64      | 99      |
| A16651 10x | LFS013.D | 59      | 69      | 65      | 76      | 72      | 97      |
| A16655 5x  | LFS021.D | 52      | 60      | 63      | 69      | 63      | 78      |
| SBLK041999 | LFS023.D | 91      | 101     | 102     | 102     | 90      | 112     |
| A16657     | LFS024.D | 80      | 85      | 82      | 99      | 62      | 90      |
| A16658     | LFS025.D | 81      | 90      | 88      | 91      | 84      | 90      |
| A16659     | LFS026.D | 72      | 78      | 84      | 90      | 68      | 92      |
| A16660     | LFS027.D | 81      | 85      | 88      | 94      | 83      | 96      |
| A16660 MS  | LFS028.D | 78      | 87      | 83      | 99      | 91      | 101     |
| A16660 MSD | LFS029.D | 89      | 104     | 100     | 106     | 106     | 115     |
| A16661     | LFS030.D | 75      | 85      | 84      | 96      | 81      | 95      |
| A16662     | LFS031.D | 82      | 98      | 93      | 103     | 101     | 104     |
| A16663     | LFS032.D | 70      | 89      | 89      | 96      | 99      | 108     |
| A16664     | LFS033.D | 80      | 91      | 91      | 99      | 95      | 102     |
| A16665     | LFS034.D | 72      | 86      | 90      | 98      | 63      | 102     |
| A16666     | LFS035.D | 67      | 85      | 103     | 109     | 73      | 109     |
| A16667     | LFS036.D | 86      | 95      | 97      | 111     | 90      | 110     |
| A16668     | LFS037.D | 83      | 97      | 95      | 107     | 93      | 110     |
| SBLK042099 | LFS041.D | 82      | 92      | 98      | 97      | 84      | 104     |
| A16669     | LFS042.D | 72      | 82      | 88      | 91      | 89      | 101     |
| A16670     | LFS043.D | 69      | 83      | 85      | 87      | 82      | 100     |
| A16671     | LFS044.D | 81      | 87      | 97      | 90      | 80      | 94      |
| A16672     | LFS045.D | 76      | 82      | 87      | 97      | 88      | 103     |
| A16672 MS  | LFS046.D | 64      | 78      | 79      | 83      | 87      | 108     |
| A16672 MSD | LFS047.D | 73      | 83      | 89      | 98      | 92      | 113     |
| A16673     | LFS048.D | 75      | 82      | 87      | 89      | 77      | 95      |
| A16674     | LFS049.D | 77      | 90      | 99      | 104     | 87      | 107     |
| A16675     | LFS050.D | 81      | 90      | 95      | 96      | 90      | 99      |
| A16676     | LFS051.D | 76      | 87      | 91      | 92      | 79      | 102     |
| A16677     | LFS052.D | 82      | 87      | 88      | 92      | 81      | 94      |
| A16678     | LFS053.D | 75      | 87      | 88      | 88      | 87      | 99      |
| A16679     | LFS054.D | 82      | 93      | 94      | 99      | 97      | 111     |

Surrogate Limits

|                            | Soil     |
|----------------------------|----------|
| Surr 1 = 2-Fluorobenzene   | (25-121) |
| Surr 2 = Phenol-d5         | (24-113) |
| Surr 3 = Nitrobenzene-d5   | (23-120) |
| Surr 4 = 2-Fluorobiphenyl  | (30-115) |
| Surr 5 = 2,4-Dibromophenol | (19-122) |
| Surr 6 = Terphenyl-d14     | (18-137) |

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Table 2.9 Results of the MS/MSD Analysis for BNA in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 Results Based on Dry Weight

Sample ID: A16552

| Compound Name              | Sample Conc<br>µg/kg | MS                   |                   |              | MSD                  |                    |               | RPD | QC Limits |     |     |    |
|----------------------------|----------------------|----------------------|-------------------|--------------|----------------------|--------------------|---------------|-----|-----------|-----|-----|----|
|                            |                      | Spike Added<br>µg/kg | MS Conc.<br>µg/kg | MS %<br>Rec. | Spike Added<br>µg/kg | MSD Conc.<br>µg/kg | MSD %<br>Rec. |     | % Rec.    | RPD |     |    |
| Phenol                     | 22700                | 3683                 | 34000             | NC           | 3683                 | 29460              | NC            | NC  | 26        | -   | 90  | 35 |
| 2-Chlorophenol             | U                    | 3683                 | 4007              | 109          | 3683                 | 3175               | 86            | 23  | 25        | -   | 102 | 50 |
| 1,4-Dichlorobenzene        | U                    | 1842                 | 1764              | 96           | 1842                 | 1444               | 78            | 20  | 28        | -   | 104 | 27 |
| N-Nitroso-Di-N-Propylamine | U                    | 1842                 | 2024              | 110          | 1842                 | 1667               | 91            | 19  | 41        | -   | 126 | 38 |
| 1,2,4-Trichlorobenzene     | U                    | 1842                 | 1937              | 105          | 1842                 | 1610               | 87            | 18  | 38        | -   | 107 | 23 |
| 4-Chloro-3-Methylphenol    | U                    | 3683                 | 4536              | 123          | 3683                 | 4031               | 109           | 12  | 26        | -   | 103 | 33 |
| Acenaphthene               | U                    | 1842                 | 2149              | 117          | 1842                 | 1843               | 100           | 15  | 31        | -   | 137 | 19 |
| 4-Nitrophenol              | U                    | 3683                 | 4042              | 110          | 3683                 | 3823               | 104           | 6   | 11        | -   | 114 | 50 |
| 2,4-Dinitrotoluene         | U                    | 1842                 | 1941              | 105          | 1842                 | 1762               | 96            | 10  | 28        | -   | 89  | 47 |
| Pentachlorophenol          | U                    | 3683                 | 3313              | 90           | 3683                 | 3168               | 86            | 5   | 17        | -   | 109 | 47 |
| Pyrene                     | U                    | 1842                 | 2464              | 134          | 1842                 | 2131               | 116           | 14  | 35        | -   | 142 | 36 |

Sample ID: A16660

| Compound Name              | Sample Conc<br>µg/kg | MS                   |                   |              | MSD                  |                    |               | RPD | QC Limits |     |     |    |
|----------------------------|----------------------|----------------------|-------------------|--------------|----------------------|--------------------|---------------|-----|-----------|-----|-----|----|
|                            |                      | Spike Added<br>µg/kg | MS Conc.<br>µg/kg | MS %<br>Rec. | Spike Added<br>µg/kg | MSD Conc.<br>µg/kg | MSD %<br>Rec. |     | % Rec.    | RPD |     |    |
| Phenol                     | 1840                 | 3370                 | 4963              | 93           | 3370                 | 5738               | 116           | 22  | 26        | -   | 90  | 35 |
| 2-Chlorophenol             | U                    | 3370                 | 2668              | 79           | 3370                 | 2979               | 88            | 11  | 25        | -   | 102 | 50 |
| 1,4-Dichlorobenzene        | U                    | 1685                 | 1242              | 74           | 1685                 | 1478               | 88            | 17  | 28        | -   | 104 | 27 |
| N-Nitroso-Di-N-Propylamine | U                    | 1685                 | 1535              | 91           | 1685                 | 1798               | 107           | 16  | 41        | -   | 126 | 38 |
| 1,2,4-Trichlorobenzene     | U                    | 1685                 | 1333              | 79           | 1685                 | 1567               | 93            | 16  | 38        | -   | 107 | 23 |
| 4-Chloro-3-Methylphenol    | U                    | 3370                 | 3231              | 96           | 3370                 | 3696               | 110           | 13  | 26        | -   | 103 | 33 |
| Acenaphthene               | U                    | 1685                 | 1657              | 98           | 1685                 | 1823               | 108           | 10  | 31        | -   | 137 | 19 |
| 4-Nitrophenol              | U                    | 3370                 | 2787              | 83           | 3370                 | 3134               | 93            | 12  | 11        | -   | 114 | 50 |
| 2,4-Dinitrotoluene         | U                    | 1685                 | 1626              | 97           | 1685                 | 1665               | 99            | 2   | 28        | -   | 89  | 47 |
| Pentachlorophenol          | U                    | 3370                 | 2540              | 75           | 3370                 | 2755               | 82            | 8   | 17        | -   | 109 | 47 |
| Pyrene                     | U                    | 1685                 | 1749              | 104          | 1685                 | 1899               | 113           | 8   | 35        | -   | 142 | 36 |

Sample ID: A16672

| Compound Name              | Sample Conc<br>µg/kg | MS                   |                   |              | MSD                  |                    |               | RPD | QC Limits |     |     |    |
|----------------------------|----------------------|----------------------|-------------------|--------------|----------------------|--------------------|---------------|-----|-----------|-----|-----|----|
|                            |                      | Spike Added<br>µg/kg | MS Conc.<br>µg/kg | MS %<br>Rec. | Spike Added<br>µg/kg | MSD Conc.<br>µg/kg | MSD %<br>Rec. |     | % Rec.    | RPD |     |    |
| Phenol                     | U                    | 4095                 | 3280              | 80           | 4095                 | 3575               | 87            | 9   | 26        | -   | 90  | 35 |
| 2-Chlorophenol             | U                    | 4095                 | 2858              | 70           | 4095                 | 3251               | 79            | 13  | 25        | -   | 102 | 50 |
| 1,4-Dichlorobenzene        | U                    | 2048                 | 1298              | 63           | 2048                 | 1462               | 71            | 12  | 28        | -   | 104 | 27 |
| N-Nitroso-Di-N-Propylamine | U                    | 2048                 | 1562              | 76           | 2048                 | 1736               | 85            | 11  | 41        | -   | 126 | 38 |
| 1,2,4-Trichlorobenzene     | U                    | 2048                 | 1607              | 79           | 2048                 | 1755               | 86            | 9   | 38        | -   | 107 | 23 |
| 4-Chloro-3-Methylphenol    | U                    | 4095                 | 3786              | 93           | 4095                 | 3954               | 97            | 4   | 26        | -   | 103 | 33 |
| Acenaphthene               | U                    | 2048                 | 1744              | 85           | 2048                 | 1982               | 97            | 13  | 31        | -   | 137 | 19 |
| 4-Nitrophenol              | U                    | 4095                 | 3661              | 89           | 4095                 | 3628               | 89            | 1   | 11        | -   | 114 | 50 |
| 2,4-Dinitrotoluene         | U                    | 2048                 | 1564              | 76           | 2048                 | 1747               | 85            | 11  | 28        | -   | 89  | 47 |
| Pentachlorophenol          | U                    | 4095                 | 3149              | 77           | 4095                 | 3301               | 81            | 5   | 17        | -   | 109 | 47 |
| Pyrene                     | 4120                 | 2048                 | 6095              | 95           | 2048                 | 6658               | 124           | 26  | 35        | -   | 142 | 36 |

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Roy F. Weston, Inc.  
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Accutest Labs  
Fresh Ponds Corp Village, Bldg B  
2235 Route 130  
Dayton, NJ 08810

Attn: Matt Cordova

9 April 1999

Project # 3347-143-001-3439 Lackawanna Foundry

As per Weston REAC Purchase Order number 102694, please analyze samples according to the following parameters:

| Analysis/Method                                 | Matrix | # of samples |
|---|--------|--------------|
| Cyanide\ SW-846-9012                            | Soil   | 30           |
| Data package: Package with Diskette Deliverable |        |              |

Samples are expected to arrive at your laboratory on April 15, 1999. **All applicable QA/QC analysis as per method, will be performed on our sample matrix. Preliminary sample and QC result tables plus a signed copy of our Chain of Custody must be faxed to REAC 10 business days after receipt of the last samples.** The complete data package is due 21 business days after receipt of last batch of samples. The complete data package must include all items on the deliverables checklist. **Expect all samples to be difficult matrix and all raw data must be included in final analytical report.**

All sample and QC results (ie: LCS, Duplicates, and Blanks) must be summarized in a ExCel diskette deliverable.

Please submit all reports and technical questions concerning this project to **John Johnson** at (732) 321-4248 or fax to (732) 494-4020. Any contractual question, please call Cynthia Lentini at (732) 321-4296.

Sincerely,

Misty Barkley  
Data Validation and Report Writing Group Leader  
Roy F. Weston, Inc. / REAC Project

MB:jj Attachments

cc R. Singhvi  
P. Campagna  
3439\non\mem\9904\sub\3439Con

V. Kansal  
Subcontracting File  
D. Angwenyi

C. Lentini  
S. Fama  
M. Barkley

00058





REAC, bn, NJ  
 (908) 321-4200  
 EPA Contract 68-C4-0022

CHA OF CUSTODY RECORD  
 Project Name Chloroform Feedlot Site  
 Project Number 03347-198-001-3489-01  
 RFW Contact J. ... Phone: 782-521-4200

No: 016909  
 SHEET NO 2 OF 3

Sample Identification

| REAC # | Sample No. | Sampling Location | Matrix | Date Collected | # of Bottles | Container/Preservative | RNA RES, RB | VOC | Analyses Requested |
|--------|------------|-------------------|--------|----------------|--------------|------------------------|-------------|-----|--------------------|
| -774   | C16656     | 6. N. ...         | S      | 4/14/91        | 1            | 402/40C                | RNA RES, RB | VOC | TALMPLABS          |
| -775   | A16657     | 7                 |        |                |              | 802/40C                | ✓           | ✓   |                    |
| -776   | B16657     | 7                 |        |                |              | 402/40C                |             | ✓   |                    |
| -777   | C16657     | 7                 |        |                |              | 402/40C                | ✓           | ✓   |                    |
| -778   | A16658     | 8                 |        |                |              | 802/40C                | ✓           | ✓   |                    |
| -779   | B16658     | 8                 |        |                |              | 402/40C                | ✓           | ✓   |                    |
| -780   | C16658     | 8                 |        |                |              | 402/40C                | ✓           | ✓   |                    |
| -781   | A16659     | 9                 |        |                |              | 802/40C                | ✓           | ✓   |                    |
| -782   | B16659     | 9                 |        |                |              | 402/40C                | ✓           | ✓   |                    |
| -783   | C16659     | 9                 |        |                |              | 802/40C                | ✓           | ✓   |                    |
| -784   | A16660     | 10                |        |                |              | 402/40C                | ✓           | ✓   |                    |
| -785   | B16660     | 10                |        |                |              | 402/40C                | ✓           | ✓   |                    |
| -786   | C16660     | 10                |        |                |              | 802/40C                | ✓           | ✓   |                    |
| -787   | A16661     | 11                |        |                |              | 402/40C                | ✓           | ✓   |                    |
| -788   | B16661     | 11                |        |                |              | 402/40C                | ✓           | ✓   |                    |
| -789   | C16661     | 11                |        |                |              | 802/40C                | ✓           | ✓   |                    |
| -790   | A16662     | 12                |        |                |              | 402/40C                | ✓           | ✓   |                    |
| -791   | B16662     | 12                |        |                |              | 802/40C                | ✓           | ✓   |                    |
| -792   | C16662     | 12                |        |                |              | 402/40C                | ✓           | ✓   |                    |
| -793   | A16663     | 13                |        |                |              | 802/40C                | ✓           | ✓   |                    |

Special Instructions

Lot to select MS/MSO  
 BNA - Beer near Am. Spec. on 1st available  
 PPS+ pesticide  
 DCF - Pol. + H. ...  
 VOC - Volatile organic compounds

FOR SUBCONTRACTING USE ONLY  
 FROM CHAIN OF CUSTODY #

| Items/Reason | Relinquished By | Date    | Received By | Date    | Relinquished By | Date    | Received By | Date    | Time  |
|--------------|-----------------|---------|-------------|---------|-----------------|---------|-------------|---------|-------|
| Cellulose    | David Agnew     | 4/15/91 | David Agnew | 4/15/91 | David Agnew     | 4/15/91 | David Agnew | 4/15/91 | 11:27 |
|              | David Agnew     | 4/15/91 | David Agnew | 4/15/91 | David Agnew     | 4/15/91 | David Agnew | 4/15/91 | 11:40 |
|              | David Agnew     | 4/15/91 | David Agnew | 4/15/91 | David Agnew     | 4/15/91 | David Agnew | 4/15/91 | 15:52 |

02-00009  
 ml - n. 11.129  
 cc - dehydrated cells

REAC, Edison, NJ  
 (908) 321-4200  
 EPA Contract 68-C4-0022

**CHAIN OF CUSTODY RECORD**

Project Name Littleton Foundry Site  
 Project Number 033747-4300-3437-01  
 RFW Contact James G. Gage Phone: 732-381-4200

No: 06910

SHEET NO. 3 OF 3

**Sample Identification**

**Analyses Requested**

| REAC # | Sample No. | Sampling Location | Matrix | Date Collected | # of Bottles | Container/Preservative | BVA/FEH/BB | VOC | TA/Me/als |
|--------|------------|-------------------|--------|----------------|--------------|------------------------|------------|-----|-----------|
| -794   | B16663     | 13                | S      | 4/14/99        | 1            | 40ml/40c               |            | ✓   |           |
| -795   | C16663     | 13                |        |                |              | 40ml/40c               |            |     |           |
| -796   | A16664     | 14                |        |                |              | 80ml/40c               | ✓          |     |           |
| -797*  | B16664     | 14                |        |                |              | 40ml/40c               |            |     |           |
| -798   | C16664     | 14                |        |                |              | 40ml/40c               |            | ✓   |           |
| -799   | A16665     | 15                |        |                |              | 80ml/40c               | ✓          |     |           |
| -800   | B16665     | 15                |        |                |              | 40ml/40c               |            |     |           |
| -801   | C16665     | 15                |        |                |              | 40ml/40c               |            |     |           |

**Matrix:**

SD - Sediment  
 DS - Drum Solids  
 DL - Drum Liquids  
 X - Other  
 PW - Potable Water  
 GW - Groundwater  
 SW - Surface Water  
 SL - Sludge  
 S - Soil  
 W - Water  
 O - Oil  
 A - Air  
 VOC: volatile organic compounds  
 ml - milliliters  
 oz - ounce

Special Instructions: do not select MS/MSD  
Water BNA - Base neutral/acid extractable  
PEST - pesticides  
PCR - poly chlorinated biphenyls  
ml - milliliters  
oz - ounce

**FOR SUBCONTRACTING USE ONLY**  
**FROM CHAIN OF CUSTODY #**

| Items/Reason  | Relinquished By | Date    | Received By | Date    | Relinquished By | Date    | Received By | Date    | Time     |
|---------------|-----------------|---------|-------------|---------|-----------------|---------|-------------|---------|----------|
| cell analysis | David Gagne     | 4/14/99 | David Gagne | 4/15/99 | David Gagne     | 4/15/99 | Y. Exume    | 4/15/99 | 11:25    |
|               |                 |         |             |         | David Gagne     | 4/15/99 | M. Yopp     | 4/15/99 | 11:40 am |
|               |                 |         |             |         | David Gagne     | 4/15/99 | J. P. Gage  | 4/15/99 | 1:35     |

REAC, on, NJ  
 (908) 321-4200  
 EPA Contract 68-C4-0022

**CH/ OF CUSTODY RECORD**

Project Name LAWANNA FOUNDRY  
 Project Number 03347-143-001-347-01  
 RFW Contact Salvatore Ferraro Phone: 131 311 4300

No. 069111

SHEET NO. 1 OF 3

**Sample Identification**

041541-

| REAC # | Sample No. | Sampling Location | Matrix | Date Collected | # of Bottles | Container/Preservative | INUP/REST/IRIS | Analyses Requested | Metals |
|--------|------------|-------------------|--------|----------------|--------------|------------------------|----------------|--------------------|--------|
| -802   | A166666    | LOCATION 16       | S      | 14-APR-99      | 1            | 8oz glass / ICE        | ✓              | ✓                  |        |
| -803   | B166666    | ↓                 |        |                | 1            | 40ml glass / ICE       | ✓              | ✓                  |        |
| -804   | C166666    | ↓                 |        |                | 1            | 4oz glass / ICE        | ✓              | ✓                  |        |
| -805   | A166667    | LOCATION 17       |        |                | 1            | 8oz / ICE              | ✓              | ✓                  |        |
| -806   | B166667    | ↓                 |        |                | 1            | 40ml / ICE             | ✓              | ✓                  |        |
| -807   | C166667    | ↓                 |        |                | 1            | 40oz / ICE             | ✓              | ✓                  |        |
| -808   | A166668    | LOCATION 18       |        |                | 1            | 8oz / ICE              | ✓              | ✓                  |        |
| -809   | B166668    | ↓                 |        |                | 1            | 40ml / ICE             | ✓              | ✓                  |        |
| -810   | C166668    | ↓                 |        |                | 1            | 4oz / ICE              | ✓              | ✓                  |        |
| -811   | A166669    | LOCATION 19       |        |                | 1            | 8oz / ICE              | ✓              | ✓                  |        |
| -812   | B166669    | ↓                 |        |                | 1            | 40ml / ICE             | ✓              | ✓                  |        |
| -813   | C166669    | ↓                 |        |                | 1            | 4oz / ICE              | ✓              | ✓                  |        |
| -814   | A166670    | LOCATION 20       |        |                | 1            | 8oz / ICE              | ✓              | ✓                  |        |
| -815   | B166670    | ↓                 |        |                | 1            | 40ml / ICE             | ✓              | ✓                  |        |
| -816   | C166670    | ↓                 |        |                | 1            | 4oz / ICE              | ✓              | ✓                  |        |
| -817   | A166671    | LOCATION 21       |        |                | 1            | 8oz / ICE              | ✓              | ✓                  |        |
| -818   | B166671    | ↓                 |        |                | 1            | 40ml / ICE             | ✓              | ✓                  |        |
| -819   | C166671    | ↓                 |        |                | 1            | 4oz / ICE              | ✓              | ✓                  |        |
| -820   | A166672    | LOCATION 22       |        |                | 1            | 8oz / ICE              | ✓              | ✓                  |        |
| -821   | B166672    | ↓                 |        |                | 1            | 40ml / ICE             | ✓              | ✓                  |        |

Matrix:

- SD - Sediment
- DS - Drum Solids
- DL - Drum Liquids
- X - Other
- PW - Potable Water
- GW - Groundwater
- SW - Surface Water
- SL - Sludge
- S - Soil
- W - Water
- O - Oil
- A - Air

Special Instructions: Lab to choose M/S/M/S/D  
 02-ounce  
 ml-milliliters  
 BNA - Benzene  
 Pest - pesticides  
 PCB - polychlorinated biphenyls  
 VOC - volatile organic compounds

**FOR SUBCONTRACTING USE ONLY**  
**FROM CHAIN OF CUSTODY #**

| Items/Reason | Relinquished By | Date    | Received By  | Date    | Time    | Items/Reason   | Relinquished By | Date    | Received By  | Date    | Time    |
|--------------|-----------------|---------|--------------|---------|---------|----------------|-----------------|---------|--------------|---------|---------|
| all analysis | Sal Ferraro     | 4/14/99 | David Argeny | 4/15/99 | 11:00am | Metal Analysis | David Argeny    | 4/15/99 | Y. Et...     | 4/15/99 | 11:05   |
|              |                 |         |              |         |         | Acc Analysis   | David Argeny    | 4/15/99 | David Argeny | 4/15/99 | 11:40am |
|              |                 |         |              |         |         | all analysis   | David Argeny    | 4/15/99 | David Argeny | 4/15/99 | 1:53    |

REAC, L son, NJ  
(908) 321-4200

**CHA OF CUSTODY RECORD**

Project Name: LAKHAWANA FOUNDRY  
 Project Number: 03341 143 001 3439 01  
 RFW Contact: SALVATORE FAMA Phone: 732 301 4200

No: 06912

SHEET NO. OF 3

**Sample Identification**

| REAC # | Sample No. | Sampling Location | Matrix | Date Collected | # of Bottles | Container/Preservative | DNA/RES/PCB | VOCs | METHALS |
|--------|------------|-------------------|--------|----------------|--------------|------------------------|-------------|------|---------|
| -822   | A16672     | LOCATION 22       | S      | 14-APR-99      | 1            | 4 OZ glass / ICE       |             |      |         |
| -823   | A16673     | LOCATION 23       | S      |                | 1            | 8 OZ glass /           | ✓           |      |         |
| -824   | B16673     | ↓                 |        |                | 1            | 40 ml glass /          |             |      |         |
| -825   | C16673     | ↓                 |        |                | 1            | 4 OZ glass /           | ✓           |      |         |
| -826   | A16674     | LOCATION 24       |        |                | 1            | 8 OZ /                 | ✓           |      |         |
| -827   | B16674     | ↓                 |        |                | 1            | 40 ml /                |             |      |         |
| -828   | C16674     | ↓                 |        |                | 1            | 4 OZ /                 |             |      |         |
| -829   | A16675     | LOCATION 25       |        |                | 1            | 8 OZ /                 | ✓           |      |         |
| -830   | B16675     | ↓                 |        |                | 1            | 40 ml /                |             |      |         |
| -831   | C16675     | ↓                 |        |                | 1            | 4 OZ /                 |             |      |         |
| -832   | A16676     | LOCATION 25-D     |        |                | 1            | 8 OZ /                 | ✓           |      |         |
| -833   | B16676     | ↓                 |        |                | 1            | 40 ml /                |             |      |         |
| -834   | C16676     | ↓                 |        |                | 1            | 4 OZ /                 |             |      |         |
| -835   | A16677     | LOCATION 26       |        |                | 1            | 8 OZ /                 | ✓           |      |         |
| -836   | B16677     | ↓                 |        |                | 1            | 40 ml /                |             |      |         |
| -837   | C16677     | ↓                 |        |                | 1            | 4 OZ /                 |             |      |         |
| -838   | A16678     | LOCATION 26-D     |        |                | 1            | 8 OZ /                 | ✓           |      |         |
| -839   | B16678     | ↓                 |        |                | 1            | 40 ml /                |             |      |         |
| -840   | C16678     | ↓                 |        |                | 1            | 4 OZ /                 |             |      |         |
| -841   | A16679     | LOCATION 27       |        |                | 1            | 8 OZ                   | ✓           |      |         |

**Analyses Requested**

Special Instructions: Light to select ms/msd

**FOR SUBCONTRACTING USE ONLY**  
**FROM CHAIN OF CUSTODY #**

Soil: 02-ounce  
 Water: ml - milliliters  
 Oil: BNA - Benzene, naphthalene, acetophenone  
 Air: volatile organic compounds  
PCB Polychlorinated biphenyls

PW - Potable Water  
 GW - Groundwater  
 SW - Surface Water  
 SL - Sludge  
 SD - Sediment  
 DS - Drum Solids  
 DL - Drum Liquids  
 X - Other

| Items/Reason        | Relinquished By | Date           | Received By          | Date           | Time           | Items/Reason        | Relinquished By      | Date           | Received By       | Date           | Time         |
|---------------------|-----------------|----------------|----------------------|----------------|----------------|---------------------|----------------------|----------------|-------------------|----------------|--------------|
| <u>all analyses</u> | <u>Sal Fama</u> | <u>4/14/99</u> | <u>David Angerey</u> | <u>4/15/99</u> | <u>11:00am</u> | <u>Acc/Analysis</u> | <u>David Angerey</u> | <u>4/15/99</u> | <u>Y. E. Fama</u> | <u>4/15/99</u> | <u>11:25</u> |
|                     |                 |                |                      |                |                | <u>Acc/Analysis</u> | <u>David Angerey</u> | <u>4/15/99</u> | <u>M. P. Fama</u> | <u>4/15/99</u> | <u>11:40</u> |
|                     |                 |                |                      |                |                | <u>Acc/Analysis</u> | <u>David Angerey</u> | <u>4/15/99</u> | <u>C. F. Fama</u> | <u>4/15/99</u> | <u>13:52</u> |



REAC: on, NJ  
(908) 321-4200

**CHA OF CUSTODY RECORD**

Project Name LICKENHAWNS FOUNDRY  
 Project Number 03047 H3 001 3439 01  
 RFW Contact Sal Fama Phone 732 321 4200

No: 06913

SHEET NO 3 OF 3

**Sample Identification**

**Analyses Requested**

| REAC #  | Sample No. | Sampling Location | Matrix | Date Collected | # of Bottles | Container/Preservative | PCRB | PCRB | VOCs | Metals |
|---|------------|-------------------|--------|----------------|--------------|------------------------|------|------|------|--------|
| 842   | B16671     | Location 27       | S      | 14-APR-11      | 1            | 40 ml glass / ICE      |      |      | ✓    |        |
| 843   | B16671     | ↓                 | ↓      | ↓              | 1            | 40 ml glass / ICE      |      |      | ✓    |        |
| 844   | A16680     | Fire Pit (VOC)    | ↓      | ↓              | 1            | 40 ml /                |      |      | ✓    |        |
| 845   | B16680     | Fire Pit (VOC)    | ↓      | ↓              | 1            | 40 ml /                |      |      | ✓    |        |
| <p><i>(The entire table below this point is crossed out with a large X)</i></p> |            |                   |        |                |              |                        |      |      |      |        |

Special Instructions: Lab to select ms/ms  
 PMA - base metal/acid extractable  
 PCB - PCBs in soil  
 OR - organic  
 0.1 - 0.11, 1.2

- Matrix:
- SD - Sediment
  - DS - Drum Solids
  - DL - Drum Liquids
  - X - Other
  - PW - Potable Water
  - GW - Groundwater
  - SW - Surface Water
  - SL - Sludge
  - S - Soil
  - W - Water
  - O - Oil
  - A - Air

**FOR SUBCONTRACTING USE ONLY**  
**FROM CHAIN OF CUSTODY #**

| Items/Reason      | Relinquished By | Date    | Received By    | Date    | Relinquished By | Date    | Received By    | Date    | Time |
|-------------------|-----------------|---------|----------------|---------|-----------------|---------|----------------|---------|------|
| all/analy 1, 1, 1 | Sal Fama        | 4/14/11 | David Angewany | 4/15/11 | David Angewany  | 4/15/11 | David Angewany | 4/15/11 | 1125 |
|                   |                 |         |                |         |                 |         |                |         |      |
|                   |                 |         |                |         |                 |         |                |         |      |
|                   |                 |         |                |         |                 |         |                |         |      |
|                   |                 |         |                |         |                 |         |                |         |      |

REAC, Ec. JR, NJ  
 (908) 321-4200

PA Contract 68-C4-0022

**CHAI, JF CUSTODY RECORD**

Project Name: LAKAWANNA FOUNDRY  
 Project Number: Q3217 143 001 3A51 01  
 RFW Contact: SALVATORE FAMA Phone: 732 321 4200

No: 06914

SHEET NO. 1 OF 2

**Sample Identification**

| REAC #  | Sample No. | Sampling Location | Matrix | Date Collected | # of Bottles | Container/Preservative | Analyses Requested |
|---------|------------|-------------------|--------|----------------|--------------|------------------------|--------------------|
| 84091-1 | D16650     | Location 1        | S      | 14-APR-97      | 1            | 8oz Glass / Ice        | CYANIDE            |
| -2      | D16651     | 2                 |        |                |              |                        |                    |
| -3      | D16652     | 2-D               |        |                |              |                        |                    |
| -4      | D16653     | 3                 |        |                |              |                        |                    |
| -5      | D16654     | 4                 |        |                |              |                        |                    |
| -6      | D16655     | 5                 |        |                |              |                        |                    |
| -7      | D16656     | 6                 |        |                |              |                        |                    |
| -8      | D16657     | 7                 |        |                |              |                        |                    |
| -9      | D16658     | 8                 |        |                |              |                        |                    |
| -10     | D16659     | 9                 |        |                |              |                        |                    |
| -11     | D16660     | 10                |        |                |              |                        |                    |
| -12     | D16661     | 11                |        |                |              |                        |                    |
| -13     | D16662     | 12                |        |                |              |                        |                    |
| -14     | D16663     | 13                |        |                |              |                        |                    |
| -15     | D16664     | 14                |        |                |              |                        |                    |
| -16     | D16665     | 15                |        |                |              |                        |                    |
| -17     | D16666     | 16                |        |                |              |                        |                    |
| -18     | D16667     | 17                |        |                |              |                        |                    |
| -19     | D16668     | 18                |        |                |              |                        |                    |
| -20     | D16669     | 19                |        |                |              |                        |                    |

FOR SUBCONTRACTING USE ONLY  
 FROM CHAIN OF CUSTODY #

RECEIVED AT ACQUISITION THE FOLLOWING TEMPERATURE 1221

Special Instructions:  
 S - Soil  
 W - Water  
 O - Oil  
 A - Air

PW - Potable Water  
 GW - Groundwater  
 SW - Surface Water  
 SL - Sludge

| Items/Reason | Relinquished By | Date    | Received By | Date    | Relinquished By | Date | Received By | Date | Time |
|--------------|-----------------|---------|-------------|---------|-----------------|------|-------------|------|------|
| Handwritten  | Handwritten     | 4/14/99 | Handwritten | 4/15/99 | Fed Ex          |      | Handwritten |      |      |

11821 0048

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## Case Narrative

The data in this report have been validated to two significant figures. Any other representation of the data is the responsibility of the user.

### Metals in Soil Package I 171

The percent recoveries for C 16655 MS/MSD exceeded the acceptable QC limits for antimony (0%, 0%), copper (20%, 63%), lead (12%, 38%), selenium (42%, 54%) and arsenic C 16655 MSD (126%). The antimony results for samples C 16650, C 16651, C 16652, C 16653, C 16654, C 16655, C 16657, C 16658, C 16659, C 16660 should be regarded as unusable. The result for antimony in sample C 16656 should be regarded as estimated. The results for copper, lead, selenium and arsenic for samples C 16650, C 16651, C 16652, C 16653, C 16654, C 16655, C 16656, C 16657, C 16658, C 16659 and C 16660 should be regarded as estimated.

The percent recoveries for C 16661 MS/MSD exceeded the acceptable QC limits for antimony (64%, 58%), manganese (10%, 0%), zinc (44%, 26%) and thallium C 16661 MS (69%). The antimony, manganese, zinc and thallium results for samples C 16661, C 16662, C 16663, C 16664, C 16665, C 16666, C 16667 and C 16668 should be regarded as estimated.

The percent recoveries for C 16669 MS/MSD exceeded the acceptable QC limits for antimony (46%, 39%), chromium (219%, 139%), nickel (173%, 136%), selenium (43%, 45%), thallium (20%, 21%) and C 16669 MS for arsenic (145%) and vanadium (126%). The thallium results for samples C 16669, C 16670, C 16671, C 16672, C 16673, C 16674, C 16675, C 16676, C 16677, C 16678 and C 16679 should be regarded as unusable. The results for antimony, chromium, nickel, selenium, arsenic and vanadium for samples C 16669, C 16670, C 16671, C 16672, C 16673, C 16674, C 16675, C 16676, C 16677, C 16678 and C 16679 should be regarded as estimated.

### Cyanide in Soil Package I 182

The data were examined and were found to be acceptable.

### BNA in Soil Package I 190

In the continuing calibration check standard of 4/28/99, the acceptable QC limits were exceeded for bis (2-chloroisopropyl) ether (29%). The data are not affected because this analyte was not detected in the associated samples.

The acceptable QC limits were exceeded for the percent recoveries of one base/neutral surrogate for A 16652 MS. The data are not affected.

## Analytical Procedure for Metals in Soil

### Sample Preparation

A representative 1-2 g (wet weight) sample, weighed to 0.01 g accuracy, was mixed with 10-mL 1:1 nitric acid, placed in a clean beaker and digested in nitric acid and hydrogen peroxide according to SW-846, Method 3050. The final reflux was either nitric acid or hydrochloric acid depending on the metals to be determined. After digestion, the samples were allowed to cool to room temperature and transferred to 100 mL volumetric flasks and diluted to volume with ASTM Type II water. The samples were analyzed for all metals, except mercury, by USEPA SW-846, Method 7000 (Atomic absorption) or Method 6010 (Inductively Coupled Argon Plasma-ICAP) procedures.

A representative 0.5-0.6 g (wet weight) sample, weighed to 0.01-g accuracy, was prepared and analyzed separately for mercury on a Varian SpectrAA-300 Atomic Absorption Spectrophotometer equipped with a Varian VGA-76 vapor gas analyzer according to SW-846, Method 7471.

A separate sample was used to determine total solids.

A reagent blank and a blank spike sample were carried through the sample preparation procedure for each batch of samples processed. One matrix spike (MS) and one matrix spike duplicate (MSD) were analyzed for each batch or for every ten samples.

### Analysis and Calculations

The instruments were calibrated and operated according to SW-846, Method 7000/7471/6010 and the manufacturers operating instructions. After calibration, initial calibration verification (ICV), initial calibration blank (ICB) and quality control check standards were run to verify proper calibration. The continuing calibration verification (CCV) and continuing calibration blank (CCB) were run after every ten samples to assure proper operation during sample analysis.

The metal concentrations in solution, in micrograms per liter ( $\mu\text{g/L}$ ) were taken from the read-out systems of the Atomic Absorption instruments. The results were converted to milligrams per kilogram ( $\text{mg/kg}$ ) by correcting the reading for the sample weight and percent solids. The ICAP results ( $\text{mg/kg}$ ) were corrected for sample weight prior to instrument read-out; the instrument read-out was then corrected for percent solids.

Final concentrations, based on wet weight are given by:

$$\text{mg metal/kg sample} = [(A \times V) / W] \times DF \times CF$$

where:

- A = Instrument read-out ( $\mu\text{g/L}$ , AA:  $\text{mg/kg}$ , ICAP )
- V = final volume of processed sample (mL, AA = 100 ICAP)
- W = weight of sample (g, AA: 1.00 ICAP)
- DF = Dilution Factor (1.00 for no dilution)
- CF = conversion factor (0.001, AA: 1.00, ICAP)

## Analytical Procedure for BNA in Soil

### Extraction Procedure

Prior to extraction each sample was spiked with a six component surrogate mixture consisting of nitrobenzene-d<sub>5</sub>, 2-fluorobiphenyl, terphenyl-d<sub>14</sub>, phenol-d<sub>5</sub>, 2-fluorophenol, and 2,4,6-tribromophenol. Thirty grams of sample was mixed with 30 g anhydrous sodium sulfate, and Soxhlet extracted for 16 hours with 300 mL of methylene chloride. The extract was concentrated to 1.0 mL, an internal standard mixture consisting of 1,4-dichlorobenzene-d<sub>4</sub>, naphthalene-d<sub>8</sub>, acenaphthene-d<sub>10</sub>, phenanthrene-d<sub>10</sub>, chrysene-d<sub>12</sub>, and perylene-d<sub>12</sub> was added, and analyzed.

### Analysis Procedure

An HP 6890/5972 Gas Chromatograph/Mass Spectrometer (GC/MS), equipped with a 6890 autosampler and controlled by a PC computer equipped with Enviroquant software was used to analyze the samples.

The instrument conditions were:

|  |  |
|--|--|
| Column   | Restek Rtx-5 (crossbonded SE-54)<br>30 meter x 0.25mm ID, 0.50 µm<br>film thickness  |
| Injection Temperature                          | 280° C   |
| Transfer Temperature                           | 280° C   |
| Source Temperature and<br>Analyzer Temperature | Controlled by thermal transfer of heat from transfer<br>line   |
| Temperature Program                            | 50°C for 0.5 min<br>20° C/min to 295° C<br>hold for 8.5 min<br>25° C/min to 310° C<br>hold for 15 min                      |
| Pulsed Split Injection                         | Split time = 2.00 min @ 8:1 split ratio<br>Pressure Pulse = 16 psi for 0.5 min, then normal                                |
| Injection Volume                               | 1 µL<br><br>Must use 4 mm ID single gooseneck liners packed<br>with 10 mm plug of silanized and conditioned glass<br>wool. |

The GC/MS system was calibrated using 5 BNA standard mixtures at 20, 50, 80, 120, and 160 µg/mL. Before analysis each day, the system was tuned with 50-ng decafluorotriphenylphosphine (DFTPP) and passed a continuing calibration check when analyzing a 50-µg/mL standard mixture in which the responses were evaluated by comparison to the average response of the calibration curve.

Table 1.1 Results of the Analysis for Metals in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 Results Based on Dry Weight

| Client ID |                 | Method Blank |           | C16650       |                  | C16651     |              | C16652       |           | C16653     |           | C16654     |
|-----------|-----------------|--------------|-----------|--------------|------------------|------------|--------------|--------------|-----------|------------|-----------|------------|
| Location  |                 | Lab          | NA        | 1 West Fence | 2 Surface Middle | 2D middle  | 3 West Fence | 4 west Fence |           |            |           |            |
| % Solids  |                 |              |           | 64.02        | 94.99            | 94.41      | 79.41        | 80.67        |           |            |           |            |
| Parameter | Analysis Method | Conc mg/kg   | MDL mg/kg | Conc mg/kg   | MDL mg/kg        | Conc mg/kg | MDL mg/kg    | Conc mg/kg   | MDL mg/kg | Conc mg/kg | MDL mg/kg | Conc mg/kg |
| Aluminum  | ICAP            | U            | 18        | 11000        | 24               | 1300       | 16           | 1200         | 16        | 7800       | 21        | 13000      |
| Antimony  | ICAP            | U            | 6.0       | U            | 8.0              | U          | 5.2          | U            | 5.4       | U          | 7.1       | U          |
| Arsenic   | AA-Fur          | U            | 0.50      | 14           | 1.5              | 1.6        | 0.50         | 2.5          | 0.48      | 11         | 0.62      | 13         |
| Barium    | ICAP            | U            | 1.0       | 160          | 1.3              | 18         | 0.87         | 17           | 0.90      | 140        | 1.2       | 150        |
| Beryllium | ICAP            | U            | 0.50      | 0.93         | 0.67             | U          | 0.44         | U            | 0.45      | 0.81       | 0.59      | 2.2        |
| Cadmium   | ICAP            | U            | 0.50      | 3.2          | 0.67             | 0.48       | 0.44         | 0.66         | 0.45      | 3.7        | 0.59      | 2.4        |
| Calcium   | ICAP            | U            | 50        | 21000        | 67               | 1900       | 44           | 3300         | 45        | 94000      | 59        | 120000     |
| Chromium  | ICAP            | U            | 0.50      | 45           | 0.67             | 25         | 0.44         | 15           | 0.45      | 33         | 0.59      | 28         |
| Cobalt    | ICAP            | U            | 1.0       | 12           | 1.3              | 2.1        | 0.87         | 1.2          | 0.90      | 7.7        | 1.2       | 6.2        |
| Copper    | ICAP            | U            | 1.0       | 74           | 1.3              | 57         | 0.87         | 35           | 0.90      | 75         | 1.2       | 67         |
| Iron      | ICAP            | U            | 10        | 43000        | 13               | 24000      | 8.7          | 13000        | 9.0       | 46000      | 12        | 30000      |
| Lead      | ICAP            | U            | 4.0       | 300          | 5.3              | 44         | 3.5          | 38           | 3.6       | 200        | 4.7       | 260        |
| Magnesium | ICAP            | U            | 50        | 4800         | 67               | 600        | 44           | 560          | 45        | 7200       | 59        | 11000      |
| Manganese | ICAP            | U            | 1.0       | 1200         | 1.3              | 250        | 0.87         | 140          | 0.90      | 800        | 1.2       | 1300       |
| Mercury   | Cold Vapor      | U            | 0.04      | 0.21         | 0.06             | 0.06       | 0.03         | 0.1          | 0.04      | 0.11       | 0.05      | 0.16       |
| Nickel    | ICAP            | U            | 1.0       | 42           | 1.3              | 15         | 0.87         | 9.2          | 0.90      | 28         | 1.2       | 32         |
| Potassium | ICAP            | U            | 200       | 1500         | 270              | U          | 170          | U            | 180       | 920        | 240       | 1500       |
| Selenium  | AA-Fur          | U            | 0.50      | 1.0          | 0.75             | U          | 0.50         | U            | 0.48      | 0.79       | 0.62      | 0.98       |
| Silver    | ICAP            | U            | 0.50      | U            | 0.67             | U          | 0.44         | U            | 0.45      | U          | 0.59      | U          |
| Sodium    | ICAP            | U            | 50        | 110          | 67               | U          | 44           | 64           | 45        | 140        | 59        | 520        |
| Thallium  | AA-Fur          | U            | 0.50      | U            | 0.75             | U          | 0.50         | U            | 0.48      | U          | 0.62      | U          |
| Vanadium  | ICAP            | U            | 2.0       | 27           | 2.7              | 5.2        | 1.7          | 3.7          | 1.8       | 17         | 2.4       | 18         |
| Zinc      | ICAP            | U            | 2.0       | 760          | 2.7              | 140        | 1.7          | 79           | 1.8       | 640        | 2.4       | 540        |

**Table 1.1 (Cont) Results of the Analysis for Metals in Soil  
WA # 3-439 Lackawanna Foundry Site  
Results Based on Dry Weight**

| Client ID |                 | C16655       |           | C16656          |           | C16657     |           | C16658     |           | C16659     |           | C16660     |           |
|-----------|-----------------|--------------|-----------|-----------------|-----------|------------|-----------|------------|-----------|------------|-----------|------------|-----------|
| Location  |                 | 5 west Fence |           | 6 Near Entrance |           | 7          |           | 8          |           | 9          |           | 10         |           |
| % Solids  |                 | 81.63        |           | 82.59           |           | 75.99      |           | 67.39      |           | 88.49      |           | 99.00      |           |
| Parameter | Analysis Method | Conc mg/kg   | MDL mg/kg | Conc mg/kg      | MDL mg/kg | Conc mg/kg | MDL mg/kg | Conc mg/kg | MDL mg/kg | Conc mg/kg | MDL mg/kg | Conc mg/kg | MDL mg/kg |
| Aluminum  | ICAP            | 12000        | 20        | 4600            | 15        | 5300       | 22        | 3900       | 23        | 5800       | 18        | 2100       | 16        |
| Antimony  | ICAP            | U            | 6.6       | 6.4             | 4.9       | U          | 7.4       | U          | 7.5       | U          | 6.0       | U          | 5.4       |
| Arsenic   | AA-Fur          | 15           | 0.55      | 20              | 0.50      | 6.3        | 0.63      | 6.5        | 0.69      | 16         | 0.51      | 1.3        | 0.49      |
| Barium    | ICAP            | 130          | 1.1       | 270             | 0.81      | 120        | 1.2       | 83         | 1.3       | 170        | 1.0       | 22         | 0.90      |
| Beryllium | ICAP            | 0.85         | 0.55      | U               | 0.41      | U          | 0.62      | U          | 0.63      | U          | 0.50      | U          | 0.45      |
| Cadmium   | ICAP            | 2.4          | 0.55      | 20              | 0.41      | 3.6        | 0.62      | 2.4        | 0.63      | 4.4        | 0.50      | 0.59       | 0.45      |
| Calcium   | ICAP            | 14000        | 55        | 21000           | 41        | 18000      | 62        | 8300       | 63        | 14000      | 50        | 2200       | 45        |
| Chromium  | ICAP            | 29           | 0.55      | 190             | 0.41      | 55         | 0.62      | 36         | 0.63      | 63         | 0.50      | 14         | 0.45      |
| Cobalt    | ICAP            | 14           | 1.1       | 20              | 0.81      | 6.7        | 1.2       | 8.5        | 1.3       | 8.4        | 1.0       | 2.0        | 0.90      |
| Copper    | ICAP            | 200          | 1.1       | 440             | 0.81      | 110        | 1.2       | 81         | 1.3       | 200        | 1.0       | 41         | 0.90      |
| Iron      | ICAP            | 35000        | 11        | 200000          | 41        | 47000      | 12        | 73000      | 13        | 76000      | 10        | 19000      | 9.0       |
| Lead      | ICAP            | 180          | 4.4       | 610             | 3.3       | 2100       | 5.0       | 230        | 5.0       | 340        | 4.0       | 34         | 3.6       |
| Magnesium | ICAP            | 4800         | 55        | 3100            | 41        | 3300       | 62        | 1800       | 63        | 4300       | 50        | 610        | 45        |
| Manganese | ICAP            | 700          | 1.1       | 1800            | 0.81      | 1100       | 1.2       | 790        | 1.3       | 1100       | 1.0       | 180        | 0.90      |
| Mercury   | Cold Vapor      | 0.11         | 0.04      | 1.1             | 0.04      | 0.22       | 0.05      | 0.13       | 0.05      | 0.95       | 0.04      | 0.27       | 0.04      |
| Nickel    | ICAP            | 41           | 1.1       | 160             | 0.81      | 110        | 1.2       | 41         | 1.3       | 160        | 1.0       | 15         | 0.90      |
| Potassium | ICAP            | 1200         | 220       | 630             | 160       | 570        | 250       | 540        | 250       | 390        | 200       | 500        | 180       |
| Selenium  | AA-Fur          | 0.58         | 0.55      | 0.77            | 0.50      | U          | 0.63      | U          | 0.69      | U          | 0.51      | U          | 0.49      |
| Silver    | ICAP            | U            | 0.55      | U               | 2.1       | U          | 0.62      | U          | 0.63      | U          | 0.50      | U          | 0.45      |
| Sodium    | ICAP            | 79           | 55        | 410             | 41        | 230        | 62        | 230        | 63        | 270        | 50        | 1500       | 45        |
| Thallium  | AA-Fur          | U            | 0.55      | U               | 0.50      | U          | 0.63      | U          | 0.69      | U          | 0.51      | U          | 0.49      |
| Vanadium  | ICAP            | 23           | 2.2       | 20              | 1.6       | 20         | 2.5       | 14         | 2.5       | 16         | 2.0       | 4.2        | 1.8       |
| Zinc      | ICAP            | 340          | 2.2       | 1100            | 1.6       | 1000       | 2.5       | 450        | 2.5       | 550        | 2.0       | 42         | 1.8       |



**Table 1.1 (Cont) Results of the Analysis for Metals in Soil  
WA # 3-439 Lackawanna Foundry Site  
Results Based on Dry Weight**

| Client ID<br>Location<br>% Solids |                    | C18673<br>Location 23<br>81.54 |              | C16674<br>Location 24<br>77.01 |              | C16675<br>Location 25<br>80.54 |              | C16676<br>Location 25-D<br>78.92 |              | C16677<br>Location 26<br>74.19 |              | C166<br>Location<br>63.2 |
|-----------------------------------|--------------------|--------------------------------|--------------|--------------------------------|--------------|--------------------------------|--------------|----------------------------------|--------------|--------------------------------|--------------|--------------------------|
| Parameter                         | Analysis<br>Method | Conc<br>mg/kg                  | MDL<br>mg/kg | Conc<br>mg/kg                  | MDL<br>mg/kg | Conc<br>mg/kg                  | MDL<br>mg/kg | Conc<br>mg/kg                    | MDL<br>mg/kg | Conc<br>mg/kg                  | MDL<br>mg/kg | Conc<br>mg/kg            |
| Aluminum                          | ICAP               | 9300                           | 19           | 7300                           | 19           | 12000                          | 18           | 17000                            | 21           | 14000                          | 23           | 14000                    |
| Antimony                          | ICAP               | U                              | 6.5          | U                              | 6.5          | U                              | 6.1          | U                                | 7.0          | U                              | 7.6          | U                        |
| Arsenic                           | AA-Fur             | 8.4                            | 0.46         | 5                              | 0.98         | 21                             | 0.49         | 25                               | 0.50         | 66                             | 0.53         | 48                       |
| Barium                            | ICAP               | 94                             | 1.1          | 64                             | 1.1          | 130                            | 1.0          | 140                              | 1.2          | 250                            | 1.3          | 280                      |
| Beryllium                         | ICAP               | 0.66                           | 0.54         | U                              | 0.54         | 0.73                           | 0.50         | 0.97                             | 0.59         | 1.3                            | 0.64         | 0.67                     |
| Cadmium                           | ICAP               | 1.5                            | 0.54         | 1.2                            | 0.54         | 2.3                            | 0.50         | U                                | 0.59         | 5.2                            | 0.64         | 1.3                      |
| Calcium                           | ICAP               | 12000                          | 54           | 10000                          | 54           | 39000                          | 50           | 11000                            | 59           | 53000                          | 64           | 120000                   |
| Chromium                          | ICAP               | 24                             | 0.54         | 18                             | 0.54         | 31                             | 0.50         | 25                               | 0.59         | 51                             | 0.64         | 23                       |
| Cobalt                            | ICAP               | 8.0                            | 1.1          | 4.8                            | 1.1          | 9.7                            | 1.0          | 15                               | 1.2          | 9.7                            | 1.3          | 7.3                      |
| Copper                            | ICAP               | 52                             | 1.1          | 49                             | 1.1          | 65                             | 1.0          | 27                               | 1.2          | 110                            | 1.3          | 34                       |
| Iron                              | ICAP               | 30000                          | 11           | 25000                          | 11           | 29000                          | 10           | 32000                            | 12           | 45000                          | 13           | 20000                    |
| Lead                              | ICAP               | 160                            | 4.3          | 130                            | 4.3          | 230                            | 4.0          | 40                               | 4.7          | 740                            | 5.1          | 100                      |
| Magnesium                         | ICAP               | 3600                           | 54           | 3500                           | 54           | 5400                           | 50           | 5400                             | 59           | 6300                           | 64           | 7300                     |
| Manganese                         | ICAP               | 560                            | 1.1          | 390                            | 1.1          | 710                            | 1.0          | 610                              | 1.2          | 1500                           | 1.3          | 570                      |
| Mercury                           | Cold Vapor         | 0.13                           | 0.04         | 0.13                           | 0.04         | 0.17                           | 0.04         | 0.08                             | 0.04         | 1.6                            | 0.04         | 0.17                     |
| Nickel                            | ICAP               | 28                             | 1.1          | 18                             | 1.1          | 38                             | 1.0          | 41                               | 1.2          | 39                             | 1.3          | 22                       |
| Potassium                         | ICAP               | 1100                           | 220          | 1100                           | 220          | 1400                           | 200          | 1500                             | 230          | 1600                           | 250          | 1400                     |
| Selenium                          | AA-Fur             | 0.57                           | 0.46         | U                              | 0.49         | U                              | 0.49         | U                                | 0.50         | 1.6                            | 0.53         | U                        |
| Silver                            | ICAP               | U                              | 0.54         | U                              | 0.54         | U                              | 0.50         | U                                | 0.59         | U                              | 0.64         | U                        |
| Sodium                            | ICAP               | 130                            | 54           | 250                            | 54           | 170                            | 50           | U                                | 59           | 190                            | 64           | 190                      |
| Thallium                          | AA-Fur             | U                              | 0.46         | U                              | 0.49         | U                              | 0.49         | U                                | 0.50         | U                              | 0.53         | U                        |
| Vanadium                          | ICAP               | 19                             | 2.2          | 14                             | 2.2          | 23                             | 2.0          | 27                               | 2.3          | 29                             | 2.5          | 20                       |
| Zinc                              | ICAP               | 660                            | 2.2          | 350                            | 2.2          | 760                            | 2.0          | 150                              | 2.3          | 1600                           | 2.5          | 330                      |

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**Table 1.1 (Cont) Results of the Analysis for Metals in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 Results Based on Dry Weight**

| Client ID | C16679          |            |           |
|-----------|-----------------|------------|-----------|
| Location  | Location 27     |            |           |
| % Solids  | 83.78           |            |           |
| Parameter | Analysis Method | Conc mg/kg | MDL mg/kg |
| Aluminum  | ICAP            | 2900       | 17        |
| Antimony  | ICAP            | U          | 5.6       |
| Arsenic   | AA-Fur          | 19         | 0.34      |
| Barium    | ICAP            | 59         | 0.94      |
| Beryllium | ICAP            | U          | 0.47      |
| Cadmium   | ICAP            | 0.84       | 0.47      |
| Calcium   | ICAP            | 15000      | 47        |
| Chromium  | ICAP            | 95         | 0.47      |
| Cobalt    | ICAP            | 8.7        | 0.94      |
| Copper    | ICAP            | 280        | 0.94      |
| Iron      | ICAP            | 78000      | 9.4       |
| Lead      | ICAP            | 140        | 3.8       |
| Magnesium | ICAP            | 2500       | 47        |
| Manganese | ICAP            | 770        | 0.94      |
| Mercury   | Cold Vapor      | 0.09       | 0.04      |
| Nickel    | ICAP            | 89         | 0.94      |
| Potassium | ICAP            | 320        | 190       |
| Selenium  | AA-Fur          | U          | 0.34      |
| Silver    | ICAP            | U          | 0.47      |
| Sodium    | ICAP            | 120        | 47        |
| Thallium  | AA-Fur          | U          | 0.34      |
| Vanadium  | ICAP            | 15         | 1.9       |
| Zinc      | ICAP            | 250        | 1.9       |

Table 1.3 (cont.) Results of the Analysis for BNA in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 (Results are Based on Dry Weight)

| Sample No                   | A16654         |              | A16655         |              | A16656          |              |
|-----------------------------|----------------|--------------|----------------|--------------|-----------------|--------------|
| Sample Location             | 4 West Fence   |              | 5 West Fence   |              | 6 Near Entrance |              |
| GC/MS File Name             | LFS009         |              | LFS010         |              | LFS012          |              |
| Dilution Factor             | 10             |              | 1              |              | 10              |              |
| % Solid                     | 82             |              | 81             |              | 85              |              |
| Compound Name               | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg  | MDL<br>µg/kg |
| Phenol                      | U              | 4100         | U              | 410          | U               | 3900         |
| bis(-2-Chloroethyl)Ether    | U              | 4100         | U              | 410          | U               | 3900         |
| 2-Chlorophenol              | U              | 4100         | U              | 410          | U               | 3900         |
| 1,3-Dichlorobenzene         | U              | 4100         | U              | 410          | U               | 3900         |
| 1,4-Dichlorobenzene         | U              | 4100         | U              | 410          | U               | 3900         |
| Benzyl alcohol              | U              | 4100         | U              | 410          | U               | 3900         |
| 1,2-Dichlorobenzene         | U              | 4100         | U              | 410          | U               | 3900         |
| 2-Methylphenol              | U              | 4100         | U              | 410          | U               | 3900         |
| bis(2-Chloroisopropyl)ether | U              | 4100         | U              | 410          | U               | 3900         |
| 4-Methylphenol              | U              | 4100         | U              | 410          | U               | 3900         |
| N-Nitroso-Di-n-propylamine  | U              | 4100         | U              | 410          | U               | 3900         |
| Hexachloroethane            | U              | 4100         | U              | 410          | U               | 3900         |
| Nitrobenzene                | U              | 4100         | U              | 410          | U               | 3900         |
| Isophorone                  | U              | 4100         | U              | 410          | U               | 3900         |
| 2-Nitrophenol               | U              | 4100         | U              | 410          | U               | 3900         |
| 2,4-Dimethylphenol          | U              | 4100         | U              | 410          | U               | 3900         |
| bis(2-Chloroethoxy)methane  | U              | 4100         | U              | 410          | U               | 3900         |
| 2,4-Dichlorophenol          | U              | 4100         | U              | 410          | U               | 3900         |
| 1,2,4-Trichlorobenzene      | U              | 4100         | U              | 410          | U               | 3900         |
| Naphthalene                 | U              | 4100         | 310            | J 410        | U               | 3900         |
| 4-Chloroaniline             | U              | 4100         | U              | 410          | U               | 3900         |
| Hexachlorobutadiene         | U              | 4100         | U              | 410          | U               | 3900         |
| 4-Chloro-3-methylphenol     | U              | 4100         | U              | 410          | U               | 3900         |
| 2-Methylnaphthalene         | U              | 4100         | 270            | J 410        | U               | 3900         |
| Hexachlorocyclopentadiene   | U              | 4100         | U              | 410          | U               | 3900         |
| 2,4,6-Trichlorophenol       | U              | 4100         | U              | 410          | U               | 3900         |
| 2,4,5-Trichlorophenol       | U              | 4100         | U              | 410          | U               | 3900         |
| 2-Chloronaphthalene         | U              | 4100         | U              | 410          | U               | 3900         |
| 2-Nitroaniline              | U              | 4100         | U              | 410          | U               | 3900         |
| Dimethylphthalate           | U              | 4100         | U              | 410          | U               | 3900         |
| Acenaphthylene              | U              | 4100         | 200            | J 410        | U               | 3900         |
| 2,6-Dinitrotoluene          | U              | 4100         | U              | 410          | U               | 3900         |
| 3-Nitroaniline              | U              | 4100         | U              | 410          | U               | 3900         |
| Acenaphthene                | U              | 4100         | 740            | 410          | U               | 3900         |
| 2,4-Dinitrophenol           | U              | 4100         | U              | 410          | U               | 3900         |
| 4-Nitrophenol               | U              | 4100         | U              | 410          | U               | 3900         |
| Dibenzofuran                | U              | 4100         | 490            | 410          | U               | 3900         |
| 2,4-Dinitrotoluene          | U              | 4100         | U              | 410          | U               | 3900         |
| Diethylphthalate            | U              | 4100         | U              | 410          | U               | 3900         |
| 4-Chlorophenyl-phenylether  | U              | 4100         | U              | 410          | U               | 3900         |
| Fluorene                    | U              | 4100         | 880            | 410          | U               | 3900         |
| 4-Nitroaniline              | U              | 4100         | U              | 410          | U               | 3900         |
| 4,6-Dinitro-2-methylphenol  | U              | 4100         | U              | 410          | U               | 3900         |
| N-Nitrosodiphenylamine      | U              | 4100         | U              | 410          | U               | 3900         |
| 4-Bromophenyl-phenylether   | U              | 4100         | U              | 410          | U               | 3900         |
| Hexachlorobenzene           | U              | 4100         | U              | 410          | U               | 3900         |
| Pentachlorophenol           | U              | 4100         | U              | 410          | U               | 3900         |
| Phenanthrene                | U              | 4100         | 7900           | 410          | U               | 3900         |
| Anthracene                  | U              | 4100         | 1800           | 410          | U               | 3900         |
| Carbazole                   | U              | 4100         | 1200           | 410          | U               | 3900         |
| Di-n-butylphthalate         | U              | 4100         | U              | 410          | U               | 3900         |
| Fluoranthene                | 2300           | J 4100       | 9900           | 410          | 2000            | J 3900       |
| Pyrene                      | 1900           | J 4100       | 7800           | 410          | 1700            | J 3900       |
| Butylbenzylphthalate        | U              | 4100         | U              | 410          | U               | 3900         |
| Benzo(a)anthracene          | 1100           | J 4100       | 5000           | 410          | 1000            | J 3900       |
| 3,3'-Dichlorobenzidine      | U              | 4100         | U              | 410          | U               | 3900         |
| Chrysene                    | 1400           | J 4100       | 5300           | 410          | 1100            | J 3900       |
| Bis(2-Ethylhexyl)phthalate  | 1000           | J 4100       | 260            | J 410        | 810             | J 3900       |
| Di-n-octylphthalate         | U              | 4100         | U              | 410          | U               | 3900         |
| Benzo(b)fluoranthene        | 1700           | J 4100       | 5800           | 410          | 1400            | J 3900       |
| Benzo(k)fluoranthene        | 1400           | J 4100       | 4500           | 410          | 1200            | J 3900       |
| Benzo(a)pyrene              | 1700           | J 4100       | 5800           | 410          | 1300            | J 3900       |
| Indeno(1,2,3-cd)pyrene      | 1300           | J 4100       | 3400           | 410          | 990             | J 3900       |
| Dibenzo(a,h)anthracene      | U              | 4100         | 1300           | 410          | U               | 3900         |
| Benzo(g,h,i)perylene        | 1600           | J 4100       | 3700           | 410          | 1100            | J 3900       |

Table 1.3 (cont.) Results of the Analysis for BNA in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 (Results are Based on Dry Weight)

| Sample No                   | A16661         | A16662       | A16663         | A16664       | A16665         |              |                |              |                |              |
|-----------------------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|
| Sample Location             | 11             | 12           | 13             | 14           | 15             |              |                |              |                |              |
| GC/MS File Name             | LS030          | LS031        | LS032          | LS033        | LS034          |              |                |              |                |              |
| Dilution Factor             | 10             | 5            | 5              | 5            | 5              |              |                |              |                |              |
| % Solid                     | 85             | 80           | 91             | 74           | 62             |              |                |              |                |              |
| Compound Name               | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg |
| Phenol                      | 9700           | 3900         | 4700           | 2100         | 2200           | 1800         | 7000           | 2300         | 1600           | J 2700       |
| bis(-2-Chloroethyl)Ether    | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 2-Chlorophenol              | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 1,3-Dichlorobenzene         | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 1,4-Dichlorobenzene         | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Benzyl alcohol              | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 1,2-Dichlorobenzene         | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 2-Methylphenol              | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| bis(2-Chloroisopropyl)ether | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 4-Methylphenol              | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| N-Nitroso-Di-n-propylamine  | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Hexachloroethane            | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Nitrobenzene                | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Isophorone                  | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 2-Nitrophenol               | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 2,4-Dimethylphenol          | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| bis(2-Chloroethoxy)methane  | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 2,4-Dichlorophenol          | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 1,2,4-Trichlorobenzene      | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Naphthalene                 | 2100 J         | 3900         | 960 J          | 2100         | 670 J          | 1800         | 1400 J         | 2300         | U              | 2700         |
| 4-Chloroaniline             | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Hexachlorobutadiene         | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 4-Chloro-3-methylphenol     | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 2-Methylnaphthalene         | U              | 3900         | 500 J          | 2100         | 400 J          | 1800         | U              | 2300         | U              | 2700         |
| Hexachlorocyclopentadiene   | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 2,4,6-Trichlorophenol       | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 2,4,5-Trichlorophenol       | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 2-Chloronaphthalene         | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 2-Nitroaniline              | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Dimethylphthalate           | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Acenaphthylene              | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 2,6-Dinitrotoluene          | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 3-Nitroaniline              | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Acenaphthene                | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 2,4-Dinitrophenol           | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 4-Nitrophenol               | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Dibenzofuran                | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 2,4-Dinitrotoluene          | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Diethylphthalate            | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 4-Chlorophenyl-phenylether  | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Fluorene                    | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 4-Nitroaniline              | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 4,6-Dinitro-2-methylphenol  | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| N-Nitrosodiphenylamine      | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 4-Bromophenyl-phenylether   | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Hexachlorobenzene           | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Pentachlorophenol           | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Phenanthrene                | U              | 3900         | 530 J          | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Anthracene                  | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Carbazole                   | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Di-n-butylphthalate         | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Fluoranthene                | U              | 3900         | 600 J          | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Pyrene                      | U              | 3900         | 900 J          | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Butylbenzylphthalate        | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Benzo(a)anthracene          | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| 3,3'-Dichlorobenzidine      | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Chrysene                    | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Bis(2-Ethylhexyl)phthalate  | 16000          | 3900         | 1300 J         | 2100         | 540 J          | 1800         | U              | 2300         | U              | 2700         |
| Di-n-octylphthalate         | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Benzo(b)fluoranthene        | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Benzo(k)fluoranthene        | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Benzo(a)pyrene              | U              | 3900         | U              | 2100         | U              | 1800         | 460 J          | 2300         | U              | 2700         |
| Indeno(1,2,3-cd)pyrene      | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Dibenzo(a,h)anthracene      | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |
| Benzo(g,h,i)perylene        | U              | 3900         | U              | 2100         | U              | 1800         | U              | 2300         | U              | 2700         |

### Introduction

REAC, in response to WA 3-439, provided analytical support for environmental samples collected from Lackawanna Foundry Site located in Lackawanna, NY as described in the following table. The support also included QA/QC, data review, and preparation of an analytical report containing a summary of the analytical methods, the results, and the QA/QC results.

The samples were treated with procedures consistent with those described in SOP # 1008 and are summarized in the following table:

| COC # | Number of Samples | Sampling Date | Date Received | Matrix | Analysis | Laboratory |
|-------|-------------------|---------------|---------------|--------|----------|------------|
| 06908 | 7                 | 4/14/99       | 4/15/99       | Soil   | BNA      | REAC       |
| 06908 | 6                 | 4/14/99       | 4/15/99       | Soil   | Metals   | REAC       |
| 06909 | 7                 | 4/14/99       | 4/15/99       | Soil   | BNA      | REAC       |
| 06909 | 7                 | 4/14/99       | 4/15/99       | Soil   | Metals   | REAC       |
| 06910 | 2                 | 4/14/99       | 4/15/99       | Soil   | BNA      | REAC       |
| 06910 | 3                 | 4/14/99       | 4/15/99       | Soil   | Metals   | REAC       |
| 06911 | 7                 | 4/14/99       | 4/15/99       | Soil   | BNA      | REAC       |
| 06911 | 6                 | 4/14/99       | 4/15/99       | Soil   | Metals   | REAC       |
| 06912 | 7                 | 4/14/99       | 4/15/99       | Soil   | BNA      | REAC       |
| 06912 | 7                 | 4/14/99       | 4/15/99       | Soil   | Metals   | REAC       |
| 06913 | 1                 | 4/14/99       | 4/15/99       | Soil   | Metals   | REAC       |
| 06914 | 20                | 4/14/99       | 4/15/99       | Soil   | Cyanide  | Accutest   |
| 06915 | 10                | 4/14/99       | 4/15/99       | Soil   | Cyanide  | Accutest   |

Analyses requested on these chains of custody, but not presented in this report, will be given in a separate report.

ANALYTICAL REPORT

Prepared by  
Roy F. Weston, Inc.

Lackawanna Foundry Site  
Lackawanna, NY

May 1999

EPA Work Assignment No. 3-439  
WESTON Work Order No. 03347-143-001-3439-01  
EPA Contract No. 68-C4-0022

Submitted to  
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EPA-ERTC

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APPENDIX D  
FINAL ANALYTICAL REPORT  
LACKAWANNA FOUNDRY SITE  
LACKAWANNA, NEW YORK  
MAY 1999





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### Introduction

REAC, in response to WA 3-439, provided analytical support for environmental samples collected from Lackawanna Foundry Site located in Lackawanna, NY as described in the following table. The support also included QA/QC, data review, and preparation of an analytical report containing a summary of the analytical methods, the results, and the QA/QC results.

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|-------|-------------------|---------------|---------------|--------|----------|------------|
| 06908 | 7                 | 4/14/99       | 4/15/99       | Soil   | BNA -    | REAC       |
| 06908 | 6                 | 4/14/99       | 4/15/99       | Soil   | Metals   | REAC       |
| 06909 | 7                 | 4/14/99       | 4/15/99       | Soil   | BNA      | REAC       |
| 06909 | 7                 | 4/14/99       | 4/15/99       | Soil   | Metals   | REAC       |
| 06910 | 2                 | 4/14/99       | 4/15/99       | Soil   | BNA      | REAC       |
| 06910 | 3                 | 4/14/99       | 4/15/99       | Soil   | Metals   | REAC       |
| 06911 | 7                 | 4/14/99       | 4/15/99       | Soil   | BNA      | REAC       |
| 06911 | 6                 | 4/14/99       | 4/15/99       | Soil   | Metals   | REAC       |
| 06912 | 7                 | 4/14/99       | 4/15/99       | Soil   | BNA      | REAC       |
| 06912 | 7                 | 4/14/99       | 4/15/99       | Soil   | Metals   | REAC       |
| 06913 | 1                 | 4/14/99       | 4/15/99       | Soil   | Metals   | REAC       |
| 06914 | 20                | 4/14/99       | 4/15/99       | Soil   | Cyanide  | Accutest   |
| 06915 | 10                | 4/14/99       | 4/15/99       | Soil   | Cyanide  | Accutest   |

Analyses requested on these chains of custody, but not presented in this report, will be given in a separate report.

## Case Narrative

The data in this report have been validated to two significant figures. Any other representation of the data is the responsibility of the user.

### Metals in Soil Package I 171

The percent recoveries for C 16655 MS/MSD exceeded the acceptable QC limits for antimony (0%, 0%), copper (20%, 63%), lead (12%, 38%), selenium (42%, 54%) and arsenic C 16655 MSD (126%). The antimony results for samples C 16650, C 16651, C 16652, C 16653, C 16654, C 16655, C 16657, C 16658, C 16659, C 16660 should be regarded as unusable. The result for antimony in sample C 16656 should be regarded as estimated. The results for copper, lead, selenium and arsenic for samples C 16650, C 16651, C 16652, C 16653, C 16654, C 16655, C 16656, C 16657, C 16658, C 16659 and C 16660 should be regarded as estimated.

The percent recoveries for C 16661 MS/MSD exceeded the acceptable QC limits for antimony (64%, 58%), manganese (10%, 0%), zinc (44%, 26%) and thallium C 16661 MS (69%). The antimony, manganese, zinc and thallium results for samples C 16661, C 16662, C 16663, C 16664, C 16665, C 16666, C 16667 and C 16668 should be regarded as estimated.

The percent recoveries for C 16669 MS/MSD exceeded the acceptable QC limits for antimony (46%, 39%), chromium (219%, 139%), nickel (173%, 136%), selenium (43%, 45%), thallium (20%, 21%) and C 16669 MS for arsenic (145%) and vanadium (126%). The thallium results for samples C 16669, C 16670, C 16671, C 16672, C 16673, C 16674, C 16675, C 16676, C 16677, C 16678 and C 16679 should be regarded as unusable. The results for antimony, chromium, nickel, selenium, arsenic and vanadium for samples C 16669, C 16670, C 16671, C 16672, C 16673, C 16674, C 16675, C 16676, C 16677, C 16678 and C 16679 should be regarded as estimated.

### Cyanide in Soil Package I 182

The data were examined and were found to be acceptable.

### BNA in Soil Package I 190

In the continuing calibration check standard of 4/28 99, the acceptable QC limits were exceeded for bis (2-chloroisopropyl) ether (29%). The data are not affected because this analyte was not detected in the associated samples.

The acceptable QC limits were exceeded for the percent recoveries of one base/neutral surrogate for A 16652 MS. The data are not affected.



### Summary of Abbreviations

|                |   |    |           |    |           |
|----------------|---|----|-----------|----|-----------|
| AA             | Atomic Absorption   |    |           |    |           |
| B              | The analyte was found in the blank  |    |           |    |           |
| BFB            | Bromofluorobenzene  |    |           |    |           |
| C              | Centigrade  |    |           |    |           |
| D              | (Surrogate Table) this value is from a diluted sample and was not calculated<br>(Result Table) this result was obtained from a diluted sample |    |           |    |           |
| Dioxin         | denotes Polychlorinated Dibenzo-p-dioxins and Polychlorinated Dibenzofurans and/or  |    |           |    |           |
| PCDD           | and PCDF  |    |           |    |           |
| CLP            | Contract Laboratory Protocol  |    |           |    |           |
| COC            | Chain of Custody  |    |           |    |           |
| CONC           | Concentration   |    |           |    |           |
| CRDL           | Contract Required Detection Limit   |    |           |    |           |
| CRQL           | Contract Required Quantitation Limit  |    |           |    |           |
| DFTPP          | Decafluorotriphenylphosphine  |    |           |    |           |
| DL             | Detection Limit   |    |           |    |           |
| E              | The value is greater than the highest linear standard and is estimated  |    |           |    |           |
| EMPC           | Estimated maximum possible concentration  |    |           |    |           |
| ICAP           | Inductively Coupled Argon Plasma  |    |           |    |           |
| ISTD           | Internal Standard   |    |           |    |           |
| J              | The value is below the method detection limit and is estimated  |    |           |    |           |
| LCS            | Laboratory Control Sample   |    |           |    |           |
| LCSD           | Laboratory Control Sample Duplicate   |    |           |    |           |
| MDL            | Method Detection Limit  |    |           |    |           |
| MI             | Matrix Interference   |    |           |    |           |
| MS             | Matrix Spike  |    |           |    |           |
| MSD            | Matrix Spike Duplicate  |    |           |    |           |
| MW             | Molecular Weight  |    |           |    |           |
| NA             | either Not Applicable or Not Available  |    |           |    |           |
| NC             | Not Calculated  |    |           |    |           |
| NR             | Not Requested   |    |           |    |           |
| NS             | Not Spiked  |    |           |    |           |
| % D            | Percent Difference  |    |           |    |           |
| % REC          | Percent Recovery  |    |           |    |           |
| PPB            | Parts per billion   |    |           |    |           |
| PPBV           | Parts per billion by volume   |    |           |    |           |
| PQL            | Practical Quantitation Limit  |    |           |    |           |
| QL             | Quantitation Limit  |    |           |    |           |
| RPD            | Relative Percent Difference   |    |           |    |           |
| RSD            | Relative Standard Deviation   |    |           |    |           |
| SIM            | Selected Ion Monitoring   |    |           |    |           |
| TCLP           | Toxic Characteristics Leaching Procedure  |    |           |    |           |
| U              | Denotes not detected  |    |           |    |           |
| W              | Weathered sample; the results should be regarded as estimated   |    |           |    |           |
| m <sup>3</sup> | cubic meter   | kg | kilogram  | μg | microgram |
| L              | liter   | g  | gram      | pg | picogram  |
| mL             | milliliter  | mg | milligram | ng | nanogram  |
| μL             | microliter  |    |           |    |           |
| *              | denotes a value that exceeds the acceptable QC limit  |    |           |    |           |
|                | Abbreviations that are specific to a particular table are explained in footnotes on that table  |    |           |    |           |

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## Analytical Procedure for Metals in Soil

### Sample Preparation

A representative 1-2 g (wet weight) sample, weighed to 0.01 g accuracy, was mixed with 10-mL 1:1 nitric acid, placed in a clean beaker and digested in nitric acid and hydrogen peroxide according to SW-846, Method 3050. The final reflux was either nitric acid or hydrochloric acid depending on the metals to be determined. After digestion, the samples were allowed to cool to room temperature and transferred to 100 mL volumetric flasks and diluted to volume with ASTM Type II water. The samples were analyzed for all metals, except mercury, by USEPA SW-846, Method 7000 (Atomic absorption) or Method 6010 (Inductively Coupled Argon Plasma-ICAP) procedures.

A representative 0.5-0.6 g (wet weight) sample, weighed to 0.01-g accuracy, was prepared and analyzed separately for mercury on a Varian SpectrAA-300 Atomic Absorption Spectrophotometer equipped with a Varian VGA-76 vapor gas analyzer according to SW-846, Method 7471.

A separate sample was used to determine total solids.

A reagent blank and a blank spike sample were carried through the sample preparation procedure for each batch of samples processed. One matrix spike (MS) and one matrix spike duplicate (MSD) were analyzed for each batch or for every ten samples.

### Analysis and Calculations

The instruments were calibrated and operated according to SW-846, Method 7000/7471/6010 and the manufacturers operating instructions. After calibration, initial calibration verification (ICV), initial calibration blank (ICB) and quality control check standards were run to verify proper calibration. The continuing calibration verification (CCV) and continuing calibration blank (CCB) were run after every ten samples to assure proper operation during sample analysis.

The metal concentrations in solution, in micrograms per liter ( $\mu\text{g/L}$ ) were taken from the read-out systems of the Atomic Absorption instruments. The results were converted to milligrams per kilogram ( $\text{mg/kg}$ ) by correcting the reading for the sample weight and percent solids. The ICAP results ( $\text{mg/kg}$ ) were corrected for sample weight prior to instrument read-out; the instrument read-out was then corrected for percent solids.

Final concentrations, based on wet weight are given by:

$$\text{mg metal/kg sample} = [(A \times V) / W] \times DF \times CF$$

where:

- A = Instrument read-out ( $\mu\text{g/L}$ , AA:  $\text{mg/kg}$ , ICAP )
- V = final volume of processed sample (mL, AA: .00 ICAP)
- W = weight of sample (g, AA: 1.00 ICAP)
- DF = Dilution Factor (1.00 for no dilution)
- CF = conversion factor (0.001, AA: 1.00, ICAP)

### Summary of Abbreviations

|                |   |    |           |    |           |
|----------------|---|----|-----------|----|-----------|
| AA             | Atomic Absorption   |    |           |    |           |
| B              | The analyte was found in the blank  |    |           |    |           |
| BFB            | Bromofluorobenzene  |    |           |    |           |
| C              | Centigrade  |    |           |    |           |
| D              | (Surrogate Table) this value is from a diluted sample and was not calculated<br>(Result Table) this result was obtained from a diluted sample |    |           |    |           |
| Dioxin         | denotes Polychlorinated Dibenzo-p-dioxins and Polychlorinated Dibenzofurans and/or  |    |           |    |           |
| PCDD           | and PCDF  |    |           |    |           |
| CLP            | Contract Laboratory Protocol  |    |           |    |           |
| COC            | Chain of Custody  |    |           |    |           |
| CONC           | Concentration   |    |           |    |           |
| CRDL           | Contract Required Detection Limit   |    |           |    |           |
| CRQL           | Contract Required Quantitation Limit  |    |           |    |           |
| DFTPP          | Decafluorotriphenylphosphine  |    |           |    |           |
| DL             | Detection Limit   |    |           |    |           |
| E              | The value is greater than the highest linear standard and is estimated  |    |           |    |           |
| EMPC           | Estimated maximum possible concentration  |    |           |    |           |
| ICAP           | Inductively Coupled Argon Plasma  |    |           |    |           |
| ISTD           | Internal Standard   |    |           |    |           |
| J              | The value is below the method detection limit and is estimated  |    |           |    |           |
| LCS            | Laboratory Control Sample   |    |           |    |           |
| LCSD           | Laboratory Control Sample Duplicate   |    |           |    |           |
| MDL            | Method Detection Limit  |    |           |    |           |
| MI             | Matrix Interference   |    |           |    |           |
| MS             | Matrix Spike  |    |           |    |           |
| MSD            | Matrix Spike Duplicate  |    |           |    |           |
| MW             | Molecular Weight  |    |           |    |           |
| NA             | either Not Applicable or Not Available  |    |           |    |           |
| NC             | Not Calculated  |    |           |    |           |
| NR             | Not Requested   |    |           |    |           |
| NS             | Not Spiked  |    |           |    |           |
| % D            | Percent Difference  |    |           |    |           |
| % REC          | Percent Recovery  |    |           |    |           |
| PPB            | Parts per billion   |    |           |    |           |
| PPBV           | Parts per billion by volume   |    |           |    |           |
| PQL            | Practical Quantitation Limit  |    |           |    |           |
| QL             | Quantitation Limit  |    |           |    |           |
| RPD            | Relative Percent Difference   |    |           |    |           |
| RSD            | Relative Standard Deviation   |    |           |    |           |
| SIM            | Selected Ion Monitoring   |    |           |    |           |
| TCLP           | Toxic Characteristics Leaching Procedure  |    |           |    |           |
| U              | Denotes not detected  |    |           |    |           |
| W              | Weathered sample; the results should be regarded as estimated   |    |           |    |           |
| m <sup>3</sup> | cubic meter   | kg | kilogram  | μg | microgram |
| L              | liter   | g  | gram      | pg | picogram  |
| mL             | milliliter  | mg | milligram | ng | nanogram  |
| μL             | microliter  |    |           |    |           |
| *              | denotes a value that exceeds the acceptable QC limit  |    |           |    |           |
|                | Abbreviations that are specific to a particular table are explained in footnotes on that table  |    |           |    |           |

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## Analytical Procedure for Metals in Soil

### Sample Preparation

A representative 1-2 g (wet weight) sample, weighed to 0.01 g accuracy, was mixed with 10-mL 1:1 nitric acid, placed in a clean beaker and digested in nitric acid and hydrogen peroxide according to SW-846, Method 3050. The final reflux was either nitric acid or hydrochloric acid depending on the metals to be determined. After digestion, the samples were allowed to cool to room temperature and transferred to 100 mL volumetric flasks and diluted to volume with ASTM Type II water. The samples were analyzed for all metals, except mercury, by USEPA SW-846, Method 7000 (Atomic absorption) or Method 6010 (Inductively Coupled Argon Plasma-ICAP) procedures.

A representative 0.5-0.6 g (wet weight) sample, weighed to 0.01-g accuracy, was prepared and analyzed separately for mercury on a Varian SpectrAA-300 Atomic Absorption Spectrophotometer equipped with a Varian VGA-76 vapor gas analyzer according to SW-846, Method 7471.

A separate sample was used to determine total solids.

A reagent blank and a blank spike sample were carried through the sample preparation procedure for each batch of samples processed. One matrix spike (MS) and one matrix spike duplicate (MSD) were analyzed for each batch or for every ten samples.

### Analysis and Calculations

The instruments were calibrated and operated according to SW-846, Method 7000/7471/6010 and the manufacturers operating instructions. After calibration, initial calibration verification (ICV), initial calibration blank (ICB) and quality control check standards were run to verify proper calibration. The continuing calibration verification (CCV) and continuing calibration blank (CCB) were run after every ten samples to assure proper operation during sample analysis.

The metal concentrations in solution, in micrograms per liter ( $\mu\text{g/L}$ ) were taken from the read-out systems of the Atomic Absorption instruments. The results were converted to milligrams per kilogram ( $\text{mg/kg}$ ) by correcting the reading for the sample weight and percent solids. The ICAP results ( $\text{mg/kg}$ ) were corrected for sample weight prior to instrument read-out; the instrument read-out was then corrected for percent solids.

Final concentrations, based on wet weight are given by:

$$\text{mg metal/kg sample} = [(A \times V) / W] \times DF \times CF$$

where:

- A = Instrument read-out ( $\mu\text{g/L}$ , AA;  $\text{mg/kg}$ , ICAP)
- V = final volume of processed sample (mL, AA = 100 ICAP)
- W = weight of sample (g, AA; 1.00 ICAP)
- DF = Dilution Factor (1.00 for no dilution)
- CF = conversion factor (0.001, AA; 1.00, ICAP)



For samples that required dilution to be within the instrument calibration range, DF is given by:

$$DF = (C+B)/C$$

where:

B = acid blank matrix used for dilution (mL)

C = sample blank aliquot (mL)

Final concentrations, based on dry weight, are given by:

$$\text{mg/kg(dry)} = [\text{mg/kg (wet)} \times 100] / S$$

where

S = percent solids

The results of the analysis are listed in Table 1.1.

### Analytical Procedure for Cyanide in Soil

The subcontract laboratory determined the concentration of cyanide in the samples by analyzing them for cyanide by USEPA Method 9012 M found in SW-846. The results of both analyses are listed in Table 1.2.

The BNA results, based on dry weight, are listed in Table 1.3; the tentatively identified compounds are listed in Table 1.4. The concentration of the detected compounds was calculated using the following equation:

$$C_u = \frac{DF \times A_u \times I_{is} \times V_i}{A_{is} \times RF \text{ ( or } RF_{ave} \text{ )} \times V_i \times W \times D}$$

where

- $C_u$  = Concentration of target analyte ( $\mu\text{g}/\text{Kg}$ )
- DF = Dilution Factor
- $A_u$  = Area of target analyte
- $I_{is}$  = Mass of specific internal standard (ng)
- $V_i$  = Volume of extract ( $\mu\text{L}$ )
- $A_{is}$  = Area of specific internal standard
- RF = Response Factor (unitless)
- $RF_{ave}$  = average Response Factor
- $V_i$  = Volume of extract injected ( $\mu\text{L}$ )
- W = Weight of sample (g)
- D = Decimal per cent solids

The  $RF_{ave}$  is used when a sample is associated with an initial calibration curve. The RF is used when a sample is associated with a continuing calibration.

Response Factor calculation:

The RF for each specific analyte is quantitated based on the area response from the continuing calibration check as follows:

$$RF = \frac{A_c \times I_{is}}{A_{is} \times I_c}$$

where

- RF = Response factor for a specific analyte
- $A_c$  = Area of the analyte in the standard
- $I_{is}$  = Mass of the specific internal standard
- $A_{is}$  = Area of the specific internal standard
- $I_c$  = Mass of the analyte in the standard

$$RF_{ave} = \frac{RF_1 + \dots + RF_n}{n}$$

and

n = number of Samples

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Table 1.1 (Cont) Results of the Analysis for Metals in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 Results Based on Dry Weight

| Client ID |                 | C16661     |           | C16662     |           | C16663     |           | C16664     |           | C16665     |           | C16666      |           |
|-----------|-----------------|------------|-----------|------------|-----------|------------|-----------|------------|-----------|------------|-----------|-------------|-----------|
| Location  |                 | 11         |           | 12         |           | 13         |           | 14         |           | 15         |           | Location 16 |           |
| % Solids  |                 | 85.56      |           | 81.01      |           | 89.86      |           | 78.52      |           | 94.98      |           | 91.68       |           |
| Parameter | Analysis Method | Conc mg/kg | MDL mg/kg | Conc mg/kg | MDL mg/kg | Conc mg/kg | MDL mg/kg | Conc mg/kg | MDL mg/kg | Conc mg/kg | MDL mg/kg | Conc mg/kg  | MDL mg/kg |
| Aluminum  | ICAP            | 740        | 18        | 4200       | 12        | 4600       | 19        | 2300       | 15        | 2200       | 17        | 1300        | 17        |
| Antimony  | ICAP            | U          | 5.9       | U          | 4.2       | 13         | 6.3       | U          | 4.8       | U          | 5.6       | U           | 5.7       |
| Arsenic   | AA-Fur          | 1.8        | 0.40      | 7.0        | 0.35      | 16         | 0.39      | 4.4        | 0.41      | 2.7        | 0.48      | 1.5         | 0.45      |
| Barium    | ICAP            | 14         | 0.98      | 84         | 0.69      | 100        | 1.0       | 39         | 0.81      | 50         | 0.93      | 15          | 0.96      |
| Beryllium | ICAP            | U          | 0.49      | U          | 0.35      | U          | 0.52      | U          | 0.40      | U          | 0.47      | U           | 0.48      |
| Cadmium   | ICAP            | 4.2        | 0.49      | 2.0        | 0.35      | 3.4        | 0.52      | 0.92       | 0.40      | 0.48       | 0.47      | U           | 0.48      |
| Calcium   | ICAP            | 1100       | 49        | 21000      | 35        | 1200       | 52        | 24000      | 40        | 1600       | 47        | 1200        | 48        |
| Chromium  | ICAP            | 13         | 0.49      | 100        | 0.35      | 79         | 0.52      | 86         | 0.40      | 14         | 0.47      | 6.1         | 0.48      |
| Cobalt    | ICAP            | 1.7        | 0.98      | 7.0        | 0.69      | 11         | 1.0       | 3.6        | 0.81      | 0.98       | 0.93      | 1.4         | 0.96      |
| Copper    | ICAP            | 33         | 0.98      | 270        | 0.69      | 370        | 1.0       | 93         | 0.81      | 100        | 0.93      | 44          | 0.96      |
| Iron      | ICAP            | 22000      | 9.8       | 56000      | 6.9       | 140000     | 53        | 39000      | 8.1       | 11000      | 9.3       | 16000       | 9.6       |
| Lead      | ICAP            | 48         | 3.9       | 150        | 2.8       | 540        | 4.2       | 110        | 3.2       | 58         | 3.7       | 58          | 3.8       |
| Magnesium | ICAP            | 240        | 49        | 2400       | 35        | 480        | 52        | 3400       | 40        | 460        | 47        | 410         | 48        |
| Manganese | ICAP            | 160        | 0.98      | 810        | 0.69      | 770        | 1.0       | 2500       | 0.81      | 120        | 0.93      | 210         | 0.96      |
| Mercury   | Cold Vapor      | 0.04       | 0.04      | 0.17       | 0.03      | 0.07       | 0.03      | 0.12       | 0.03      | U          | 0.04      | U           | 0.04      |
| Nickel    | ICAP            | 13         | 0.98      | 75         | 0.69      | 72         | 1.0       | 25         | 0.81      | 31         | 0.93      | 9.5         | 0.96      |
| Potassium | ICAP            | U          | 200       | 430        | 140       | 500        | 210       | 260        | 160       | U          | 190       | U           | 190       |
| Selenium  | AA-Fur          | U          | 0.40      | U          | 0.35      | 0.39       | 0.39      | U          | 0.41      | U          | 0.48      | U           | 0.45      |
| Silver    | ICAP            | U          | 0.49      | U          | 0.35      | U          | 0.52      | U          | 0.40      | U          | 0.47      | U           | 0.48      |
| Sodium    | ICAP            | U          | 49        | 210        | 35        | 200        | 52        | 92         | 40        | 72         | 47        | U           | 48        |
| Thallium  | AA-Fur          | U          | 0.40      | U          | 0.35      | U          | 0.39      | U          | 0.41      | U          | 0.48      | U           | 0.45      |
| Vanadium  | ICAP            | 2.7        | 2.0       | 14         | 1.4       | 16         | 2.1       | 43         | 1.6       | 5.4        | 1.9       | 3.1         | 1.9       |
| Zinc      | ICAP            | 81         | 2.0       | 280        | 1.4       | 660        | 2.1       | 190        | 1.6       | 110        | 1.9       | 98          | 1.9       |

00011

**Table 1.1 (Cont) Results of the Analysis for Metals in Soil  
WA # 3-439 Lackawanna Foundry Site  
Results Based on Dry Weight**

| Client ID |                 | C16667      |           | C16668      |           | C16669      |           | C16670      |           | C16671      |           | C16672      |           |
|-----------|-----------------|-------------|-----------|-------------|-----------|-------------|-----------|-------------|-----------|-------------|-----------|-------------|-----------|
| Location  |                 | Location 17 |           | Location 18 |           | Location 19 |           | Location 20 |           | Location 21 |           | Location 22 |           |
| % Solids  |                 | 90.83       |           | 92.93       |           | 95.49       |           | 88.29       |           | 51.93       |           | 78.77       |           |
| Parameter | Analysis Method | Conc mg/kg  | MDL mg/kg | Conc mg/kg  | MDL mg/kg | Conc mg/kg  | MDL mg/kg | Conc mg/kg  | MDL mg/kg | Conc mg/kg  | MDL mg/kg | Conc mg/kg  | MDL mg/kg |
| Aluminum  | ICAP            | 5900        | 17        | 790         | 17        | 1400        | 15        | 2400        | 17        | 12000       | 23        | 15000       | 20        |
| Antimony  | ICAP            | 6.5         | 5.8       | U           | 5.6       | U           | 5.0       | U           | 5.7       | U           | 7.8       | U           | 6.7       |
| Arsenic   | AA-Fur          | 7.1         | 0.95      | 4.4         | 0.45      | 6.7         | 0.51      | 6.5         | 0.46      | 20          | 0.54      | 77          | 0.58      |
| Barium    | ICAP            | 180         | 0.97      | 9.0         | 0.94      | 29          | 0.83      | 130         | 0.94      | 200         | 1.3       | 170         | 1.1       |
| Beryllium | ICAP            | U           | 0.48      | U           | 0.47      | U           | 0.42      | U           | 0.47      | 0.96        | 0.65      | 1.2         | 0.56      |
| Cadmium   | ICAP            | U           | 0.48      | U           | 0.47      | 0.59        | 0.42      | 2.3         | 0.47      | 13          | 0.65      | 2.3         | 0.56      |
| Calcium   | ICAP            | 130000      | 48        | 1200        | 47        | 850         | 42        | 3800        | 47        | 23000       | 65        | 34000       | 56        |
| Chromium  | ICAP            | 33          | 0.48      | 8.9         | 0.47      | 97          | 0.42      | 130         | 0.47      | 95          | 0.65      | 32          | 0.56      |
| Cobalt    | ICAP            | 47          | 0.97      | U           | 0.94      | 4.4         | 0.83      | 12          | 0.94      | 15          | 1.3       | 11          | 1.1       |
| Copper    | ICAP            | 240         | 0.97      | 28          | 0.94      | 300         | 0.83      | 720         | 0.94      | 270         | 1.3       | 47          | 1.1       |
| Iron      | ICAP            | 25000       | 9.7       | 5800        | 9.4       | 58000       | 8.3       | 93000       | 9.4       | 110000      | 13        | 32000       | 11        |
| Lead      | ICAP            | 300         | 3.9       | 21          | 3.7       | 140         | 3.3       | 280         | 3.8       | 660         | 5.2       | 180         | 4.5       |
| Magnesium | ICAP            | 19000       | 48        | 340         | 47        | 410         | 42        | 1000        | 47        | 4700        | 65        | 5900        | 56        |
| Manganese | ICAP            | 840         | 0.97      | 75          | 0.94      | 760         | 0.83      | 690         | 0.94      | 1600        | 1.3       | 1200        | 1.1       |
| Mercury   | Cold Vapor      | 0.06        | 0.04      | 0.04        | 0.04      | U           | 0.03      | 1.3         | 0.04      | 0.77        | 0.05      | 0.73        | 0.04      |
| Nickel    | ICAP            | 38          | 0.97      | 2.7         | 0.94      | 57          | 0.83      | 150         | 0.94      | 100         | 1.3       | 36          | 1.1       |
| Potassium | ICAP            | 470         | 190       | U           | 190       | U           | 170       | 280         | 190       | 900         | 260       | 1500        | 220       |
| Selenium  | AA-Fur          | U           | 0.47      | U           | 0.45      | U           | 0.51      | U           | 0.46      | 1.0         | 0.54      | U           | 0.58      |
| Silver    | ICAP            | U           | 0.48      | U           | 0.47      | U           | 0.42      | U           | 0.47      | U           | 0.65      | U           | 0.56      |
| Sodium    | ICAP            | 330         | 48        | U           | 47        | 68          | 42        | 68          | 47        | 360         | 65        | 120         | 56        |
| Thallium  | AA-Fur          | U           | 0.47      | U           | 0.45      | U           | 0.51      | U           | 0.46      | U           | 0.54      | U           | 0.58      |
| Vanadium  | ICAP            | 11          | 1.9       | 2.6         | 1.9       | 14          | 1.7       | 21          | 1.9       | 36          | 2.6       | 27          | 2.2       |
| Zinc      | ICAP            | 1300        | 1.9       | 55          | 1.9       | 310         | 1.7       | 290         | 1.9       | 1000        | 2.6       | 1100        | 2.2       |

Table 1.1 (Cont) Results of the Analysis for Metals in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 Results Based on Dry Weight

| Client ID |                 | C16667      |           | C16668      |           | C16669      |           | C16670      |           | C16671      |           | C16672      |           |
|-----------|-----------------|-------------|-----------|-------------|-----------|-------------|-----------|-------------|-----------|-------------|-----------|-------------|-----------|
| Location  |                 | Location 17 |           | Location 18 |           | Location 19 |           | Location 20 |           | Location 21 |           | Location 22 |           |
| % Solids  |                 | 90.83       |           | 92.93       |           | 95.49       |           | 88.29       |           | 51.93       |           | 78.77       |           |
| Parameter | Analysis Method | Conc mg/kg  | MDL mg/kg | Conc mg/kg  | MDL mg/kg | Conc mg/kg  | MDL mg/kg | Conc mg/kg  | MDL mg/kg | Conc mg/kg  | MDL mg/kg | Conc mg/kg  | MDL mg/kg |
| Aluminum  | ICAP            | 5900        | 17        | 790         | 17        | 1400        | 15        | 2400        | 17        | 12000       | 23        | 15000       | 20        |
| Antimony  | ICAP            | 6.5         | 5.8       | U           | 5.6       | U           | 5.0       | U           | 5.7       | U           | 7.8       | U           | 6.7       |
| Arsenic   | AA-Fur          | 7.1         | 0.95      | 4.4         | 0.45      | 6.7         | 0.51      | 6.5         | 0.46      | 20          | 0.54      | 77          | 0.58      |
| Barium    | ICAP            | 180         | 0.97      | 9.0         | 0.94      | 29          | 0.83      | 130         | 0.94      | 200         | 1.3       | 170         | 1.1       |
| Beryllium | ICAP            | U           | 0.48      | U           | 0.47      | U           | 0.42      | U           | 0.47      | 0.96        | 0.65      | 1.2         | 0.56      |
| Cadmium   | ICAP            | U           | 0.48      | U           | 0.47      | 0.59        | 0.42      | 2.3         | 0.47      | 13          | 0.65      | 2.3         | 0.56      |
| Calcium   | ICAP            | 130000      | 48        | 1200        | 47        | 850         | 42        | 3800        | 47        | 23000       | 65        | 34000       | 56        |
| Chromium  | ICAP            | 33          | 0.48      | 8.9         | 0.47      | 97          | 0.42      | 130         | 0.47      | 95          | 0.65      | 32          | 0.56      |
| Cobalt    | ICAP            | 47          | 0.97      | U           | 0.94      | 4.4         | 0.83      | 12          | 0.94      | 15          | 1.3       | 11          | 1.1       |
| Copper    | ICAP            | 240         | 0.97      | 28          | 0.94      | 300         | 0.83      | 720         | 0.94      | 270         | 1.3       | 47          | 1.1       |
| Iron      | ICAP            | 25000       | 9.7       | 5800        | 9.4       | 58000       | 8.3       | 93000       | 9.4       | 110000      | 13        | 32000       | 11        |
| Lead      | ICAP            | 300         | 3.9       | 21          | 3.7       | 140         | 3.3       | 280         | 3.8       | 660         | 5.2       | 180         | 4.5       |
| Magnesium | ICAP            | 19000       | 48        | 340         | 47        | 410         | 42        | 1000        | 47        | 4700        | 65        | 5900        | 56        |
| Manganese | ICAP            | 840         | 0.97      | 75          | 0.94      | 760         | 0.83      | 690         | 0.94      | 1600        | 1.3       | 1200        | 1.1       |
| Mercury   | Cold Vapor      | 0.06        | 0.04      | 0.04        | 0.04      | U           | 0.03      | 1.3         | 0.04      | 0.77        | 0.05      | 0.73        | 0.04      |
| Nickel    | ICAP            | 38          | 0.97      | 2.7         | 0.94      | 57          | 0.83      | 150         | 0.94      | 100         | 1.3       | 36          | 1.1       |
| Potassium | ICAP            | 470         | 190       | U           | 190       | U           | 170       | 280         | 190       | 900         | 260       | 1500        | 220       |
| Selenium  | AA-Fur          | U           | 0.47      | U           | 0.45      | U           | 0.51      | U           | 0.46      | 1.0         | 0.54      | U           | 0.58      |
| Silver    | ICAP            | U           | 0.48      | U           | 0.47      | U           | 0.42      | U           | 0.47      | U           | 0.65      | U           | 0.56      |
| Sodium    | ICAP            | 330         | 48        | U           | 47        | 68          | 42        | 68          | 47        | 360         | 65        | 120         | 56        |
| Thallium  | AA-Fur          | U           | 0.47      | U           | 0.45      | U           | 0.51      | U           | 0.46      | U           | 0.54      | U           | 0.58      |
| Vanadium  | ICAP            | 11          | 1.9       | 2.6         | 1.9       | 14          | 1.7       | 21          | 1.9       | 36          | 2.6       | 27          | 2.2       |
| Zinc      | ICAP            | 1300        | 1.9       | 55          | 1.9       | 310         | 1.7       | 290         | 1.9       | 1000        | 2.6       | 1100        | 2.2       |

00012

**Table 1.1 (Cont) Results of the Analysis for Metals in Soil  
WA # 3-439 Lackawanna Foundry Site  
Results Based on Dry Weight**

| Client ID |                 | C16673      |           | C16674      |           | C16675      |           | C16676        |           | C16677      |           | C          |
|-----------|-----------------|-------------|-----------|-------------|-----------|-------------|-----------|---------------|-----------|-------------|-----------|------------|
| Location  |                 | Location 23 |           | Location 24 |           | Location 25 |           | Location 25-D |           | Location 26 |           | Loca       |
| % Solids  |                 | 81.54       |           | 77.01       |           | 80.54       |           | 78.92         |           | 74.19       |           | (          |
| Parameter | Analysis Method | Conc mg/kg  | MDL mg/kg | Conc mg/kg  | MDL mg/kg | Conc mg/kg  | MDL mg/kg | Conc mg/kg    | MDL mg/kg | Conc mg/kg  | MDL mg/kg | Conc mg/kg |
| Aluminum  | ICAP            | 9300        | 19        | 7300        | 19        | 12000       | 18        | 17000         | 21        | 14000       | 23        | 1400       |
| Antimony  | ICAP            | U           | 6.5       | U           | 6.5       | U           | 6.1       | U             | 7.0       | U           | 7.6       |            |
| Arsenic   | AA-Fur          | 8.4         | 0.46      | 5           | 0.98      | 21          | 0.49      | 25            | 0.50      | 66          | 0.53      | 4          |
| Barium    | ICAP            | 94          | 1.1       | 64          | 1.1       | 130         | 1.0       | 140           | 1.2       | 250         | 1.3       | 28         |
| Beryllium | ICAP            | 0.66        | 0.54      | U           | 0.54      | 0.73        | 0.50      | 0.97          | 0.59      | 1.3         | 0.64      | 0.6        |
| Cadmium   | ICAP            | 1.5         | 0.54      | 1.2         | 0.54      | 2.3         | 0.50      | U             | 0.59      | 5.2         | 0.64      | 1.         |
| Calcium   | ICAP            | 12000       | 54        | 10000       | 54        | 39000       | 50        | 11000         | 59        | 53000       | 64        | 12000      |
| Chromium  | ICAP            | 24          | 0.54      | 18          | 0.54      | 31          | 0.50      | 25            | 0.59      | 51          | 0.64      | 2          |
| Cobalt    | ICAP            | 8.0         | 1.1       | 4.8         | 1.1       | 9.7         | 1.0       | 15            | 1.2       | 9.7         | 1.3       | 7          |
| Copper    | ICAP            | 52          | 1.1       | 49          | 1.1       | 65          | 1.0       | 27            | 1.2       | 110         | 1.3       | 3          |
| Iron      | ICAP            | 30000       | 11        | 25000       | 11        | 29000       | 10        | 32000         | 12        | 45000       | 13        | 2000       |
| Lead      | ICAP            | 160         | 4.3       | 130         | 4.3       | 230         | 4.0       | 40            | 4.7       | 740         | 5.1       | 10         |
| Magnesium | ICAP            | 3600        | 54        | 3500        | 54        | 5400        | 50        | 5400          | 59        | 6300        | 64        | 730        |
| Manganese | ICAP            | 560         | 1.1       | 390         | 1.1       | 710         | 1.0       | 610           | 1.2       | 1500        | 1.3       | 57         |
| Mercury   | Cold Vapor      | 0.13        | 0.04      | 0.13        | 0.04      | 0.17        | 0.04      | 0.08          | 0.04      | 1.6         | 0.04      | 0.1        |
| Nickel    | ICAP            | 28          | 1.1       | 18          | 1.1       | 38          | 1.0       | 41            | 1.2       | 39          | 1.3       | 2          |
| Potassium | ICAP            | 1100        | 220       | 1100        | 220       | 1400        | 200       | 1500          | 230       | 1600        | 250       | 140        |
| Selenium  | AA-Fur          | 0.57        | 0.46      | U           | 0.49      | U           | 0.49      | U             | 0.50      | 1.6         | 0.53      |            |
| Silver    | ICAP            | U           | 0.54      | U           | 0.54      | U           | 0.50      | U             | 0.59      | U           | 0.64      |            |
| Sodium    | ICAP            | 130         | 54        | 250         | 54        | 170         | 50        | U             | 59        | 190         | 64        | 19         |
| Thallium  | AA-Fur          | U           | 0.46      | U           | 0.49      | U           | 0.49      | U             | 0.50      | U           | 0.53      |            |
| Vanadium  | ICAP            | 19          | 2.2       | 14          | 2.2       | 23          | 2.0       | 27            | 2.3       | 29          | 2.5       | 2          |
| Zinc      | ICAP            | 660         | 2.2       | 350         | 2.2       | 760         | 2.0       | 150           | 2.3       | 1600        | 2.5       | 30         |

00013

Table 1.1 (Cont) Results of the Analysis for Metals in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 Results Based on Dry Weight

| Client ID | C16679          |            |           |
|-----------|-----------------|------------|-----------|
| Location  | Location 27     |            |           |
| % Solids  | 83.78           |            |           |
| Parameter | Analysis Method | Conc mg/kg | MDL mg/kg |
| Aluminum  | ICAP            | 2900       | 17        |
| Antimony  | ICAP            | U          | 5.6       |
| Arsenic   | AA-Fur          | 19         | 0.34      |
| Barium    | ICAP            | 59         | 0.94      |
| Beryllium | ICAP            | U          | 0.47      |
| Cadmium   | ICAP            | 0.84       | 0.47      |
| Calcium   | ICAP            | 15000      | 47        |
| Chromium  | ICAP            | 95         | 0.47      |
| Cobalt    | ICAP            | 8.7        | 0.94      |
| Copper    | ICAP            | 280        | 0.94      |
| Iron      | ICAP            | 78000      | 9.4       |
| Lead      | ICAP            | 140        | 3.8       |
| Magnesium | ICAP            | 2500       | 47        |
| Manganese | ICAP            | 770        | 0.94      |
| Mercury   | Cold Vapor      | 0.09       | 0.04      |
| Nickel    | ICAP            | 89         | 0.94      |
| Potassium | ICAP            | 320        | 190       |
| Selenium  | AA-Fur          | U          | 0.34      |
| Silver    | ICAP            | U          | 0.47      |
| Sodium    | ICAP            | 120        | 47        |
| Thallium  | AA-Fur          | U          | 0.34      |
| Vanadium  | ICAP            | 15         | 1.9       |
| Zinc      | ICAP            | 250        | 1.9       |

03014



Table 1.2 Results of the Analysis for Cyanide in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 Results Based on Dry Weight

| Sample ID    | Location    | % Solids | Cyanide       |              |
|--------------|-------------|----------|---------------|--------------|
|              |             |          | Conc<br>mg/kg | MDL<br>mg/kg |
| Method Blank | -           | 100      | U             | 1.0          |
| Method Blank | -           | 100      | U             | 1.0          |
| D16650       | 1           | 56.3     | U             | 1.8          |
| D16651       | 2           | 94.1     | U             | 1.1          |
| D16652       | 2-D         | 95.2     | U             | 1.0          |
| D16653       | 3           | 81.4     | U             | 1.2          |
| D16654       | 4           | 86.6     | U             | 1.2          |
| D16655       | 5           | 84.0     | U             | 1.2          |
| D16656       | 6           | 86.6     | U             | 1.2          |
| D16657       | 7           | 79.9     | U             | 1.2          |
| D16658       | 8           | 79.1     | U             | 1.3          |
| D16659       | 9           | 89.7     | U             | 1.1          |
| D16660       | 10          | 99.2     | U             | 1.0          |
| D16661       | 11          | 85.8     | U             | 1.2          |
| D16662       | 12          | 77.3     | U             | 1.3          |
| D16663       | 13          | 92.0     | U             | 1.1          |
| D16664       | 14          | 75.9     | U             | 1.3          |
| D16665       | 15          | 94.7     | U             | 1.0          |
| D16666       | 16          | 90.3     | U             | 1.1          |
| D16667       | 17          | 91.8     | U             | 1.1          |
| D16668       | 18          | 93.1     | U             | 1.1          |
| D16669       | 19          | 93.2     | U             | 1.1          |
| D16670       | Location 20 | 90.8     | U             | 1.1          |
| D16671       | 21          | 43.8     | U             | 2.3          |
| D16672       | 22          | 81.8     | U             | 1.2          |
| D16673       | 23          | 84.0     | U             | 1.2          |
| D16674       | 24          | 81.4     | U             | 1.2          |
| D16675       | 25          | 87.5     | U             | 1.1          |
| D16676       | 25-D        | 78.2     | U             | 1.3          |
| D16677       | 26          | 76.7     | U             | 1.3          |
| D16678       | 26-D        | 75.2     | U             | 1.3          |
| D16679       | 27          | 84.7     | U             | 1.2          |

00015

Table 1.3 Results of the Analysis for BNA in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 (Results are Based on Dry Weight)

| Sample No                   | SBLK041599     | A16650       | A16651           | A16652       | A16653        |              |               |              |               |              |
|-----------------------------|----------------|--------------|------------------|--------------|---------------|--------------|---------------|--------------|---------------|--------------|
| Sample Location             | Lab Blank      | 1 West Fence | 2 Surface Middle | 2D Middle    | 3 West Fence  |              |               |              |               |              |
| GC/MS File Name             | LFS002         | LFS003       | LFS004           | LFS005       | LFS008        |              |               |              |               |              |
| Dilution Factor             | 1              | 10           | 1                | 5            | 10            |              |               |              |               |              |
| % Solid                     | 100            | 57           | 95               | 91           | 73            |              |               |              |               |              |
| Compound Name               | Conc.<br>µg/kg | MDL<br>µg/kg | Conc<br>µg/kg    | MDL<br>µg/kg | Conc<br>µg/kg | MDL<br>µg/kg | Conc<br>µg/kg | MDL<br>µg/kg | Conc<br>µg/kg | MDL<br>µg/kg |
| Phenol                      | U              | 330          | U                | 5800         | 16000         | 350          | 23000         | 1800         | U             | 4600         |
| bis(-2-Chloroethyl)Ether    | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| 2-Chlorophenol              | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| 1,3-Dichlorobenzene         | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| 1,4-Dichlorobenzene         | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| Benzyl alcohol              | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| 1,2-Dichlorobenzene         | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| 2-Methylphenol              | U              | 330          | U                | 5800         | 370           | 350          | 410           | J 1800       | U             | 4600         |
| bis(2-Chloroisopropyl)ether | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| 4-Methylphenol              | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| N-Nitroso-Di-n-propylamine  | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| Hexachloroethane            | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| Nitrobenzene                | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| Isophorone                  | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| 2-Nitrophenol               | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| 2,4-Dimethylphenol          | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| bis(2-Chloroethoxy)methane  | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| 2,4-Dichlorophenol          | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| 1,2,4-Trichlorobenzene      | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| Naphthalene                 | U              | 330          | U                | 5800         | 970           | 350          | 960           | J 1800       | U             | 4600         |
| 4-Chloroaniline             | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| Hexachlorobutadiene         | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| 4-Chloro-3-methylphenol     | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| 2-Methylnaphthalene         | U              | 330          | U                | 5800         | 360           | 350          | 640           | J 1800       | U             | 4600         |
| Hexachlorocyclopentadiene   | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| 2,4,6-Trichlorophenol       | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| 2,4,5-Trichlorophenol       | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| 2-Chloronaphthalene         | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| 2-Nitroaniline              | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| Dimethylphthalate           | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| Acenaphthylene              | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| 2,6-Dinitrotoluene          | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| 3-Nitroaniline              | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| Acenaphthene                | U              | 330          | U                | 5800         | 95            | J 350        | U             | 1800         | U             | 4600         |
| 2,4-Dinitrophenol           | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| 4-Nitrophenol               | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| Dibenzofuran                | U              | 330          | U                | 5800         | 100           | J 350        | U             | 1800         | U             | 4600         |
| 2,4-Dinitrotoluene          | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| Diethylphthalate            | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| 4-Chlorophenyl-phenylether  | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| Fluorene                    | U              | 330          | U                | 5800         | 87            | J 350        | U             | 1800         | U             | 4600         |
| 4-Nitroaniline              | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| 4,6-Dinitro-2-methylphenol  | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| N-Nitrosodiphenylamine      | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| 4-Bromophenyl-phenylether   | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| Hexachlorobenzene           | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| Pentachlorophenol           | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| Phenanthrene                | U              | 330          | U                | 5800         | 1400          | 350          | U             | 1800         | U             | 4600         |
| Anthracene                  | U              | 330          | U                | 5800         | 220           | J 350        | U             | 1800         | U             | 4600         |
| Carbazole                   | U              | 330          | U                | 5800         | 190           | J 350        | U             | 1800         | U             | 4600         |
| Di-n-butylphthalate         | U              | 330          | U                | 5800         | 150           | J 350        | U             | 1800         | U             | 4600         |
| Fluoranthene                | U              | 330          | U                | 5800         | 2000          | 350          | 380           | J 1800       | 1000          | J 4600       |
| Pyrene                      | U              | 330          | U                | 5800         | 1500          | 350          | U             | 1800         | U             | 4600         |
| Butylbenzylphthalate        | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| Benzo(a)anthracene          | U              | 330          | U                | 5800         | 800           | 350          | U             | 1800         | U             | 4600         |
| 3,3'-Dichlorobenzidine      | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| Chrysene                    | U              | 330          | U                | 5800         | 920           | 350          | U             | 1800         | U             | 4600         |
| Bis(2-Ethylhexyl)phthalate  | U              | 330          | U                | 5800         | 640           | 350          | 1300          | J 1800       | 1000          | J 4600       |
| Di-n-octylphthalate         | U              | 330          | U                | 5800         | U             | 350          | U             | 1800         | U             | 4600         |
| Benzo(b)fluoranthene        | U              | 330          | U                | 5800         | 1100          | 350          | U             | 1800         | U             | 4600         |
| Benzo(k)fluoranthene        | U              | 330          | U                | 5800         | 880           | 350          | U             | 1800         | U             | 4600         |
| Benzo(a)pyrene              | U              | 330          | U                | 5800         | 1100          | 350          | U             | 1800         | U             | 4600         |
| Indeno(1,2,3-cd)pyrene      | U              | 330          | U                | 5800         | 750           | 350          | U             | 1800         | U             | 4600         |
| Dibenzo(a,h)anthracene      | U              | 330          | U                | 5800         | 300           | J 350        | U             | 1800         | U             | 4600         |
| Benzo(g,h,i)perylene        | U              | 330          | U                | 5800         | 890           | 350          | U             | 1800         | U             | 4600         |

Table 1.3 (cont.) Results of the Analysis for BNA in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 (Results are Based on Dry Weight)

| Sample No                   | SBLK041999     | A16657       | A16658         | A16659       | A16660         |              |                |              |                |              |
|-----------------------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|
| Sample Location             | Lab Blank      | 7            | 8              | 9            | 10             |              |                |              |                |              |
| GC/MS File Name             | LFS023         | LFS024       | LFS025         | LFS026       | LFS027         |              |                |              |                |              |
| Dilution Factor             | 1              | 10           | 10             | 10           | 5              |              |                |              |                |              |
| % Solid                     | 100            | 76           | 87             | 91           | 99             |              |                |              |                |              |
| Compound Name               | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg |
| Phenol                      | U              | 330          | 1900           | J 4400       | 1500           | J 3800       | 4900           | 3700         | 1800           | 1700         |
| bis(-2-Chloroethyl)Ether    | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| 2-Chlorophenol              | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| 1,3-Dichlorobenzene         | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| 1,4-Dichlorobenzene         | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| Benzyl alcohol              | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| 1,2-Dichlorobenzene         | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| 2-Methylphenol              | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | 570            | J 1700       |
| bis(2-Chloroisopropyl)ether | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| 4-Methylphenol              | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| N-Nitroso-Di-n-propylamine  | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| Hexachloroethane            | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| Nitrobenzene                | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| Isophorone                  | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| 2-Nitrophenol               | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| 2,4-Dimethylphenol          | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| bis(2-Chloroethoxy)methane  | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| 2,4-Dichlorophenol          | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| 1,2,4-Trichlorobenzene      | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| Naphthalene                 | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | 3100           | 1700         |
| 4-Chloroaniline             | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| Hexachlorobutadiene         | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| 4-Chloro-3-methylphenol     | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| 2-Methylnaphthalene         | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | 1500           | J 1700       |
| Hexachlorocyclopentadiene   | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| 2,4,6-Trichlorophenol       | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| 2,4,5-Trichlorophenol       | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| 2-Chloronaphthalene         | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| 2-Nitroaniline              | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| Dimethylphthalate           | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| Acenaphthylene              | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| 2,6-Dinitrotoluene          | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| 3-Nitroaniline              | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| Acenaphthene                | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| 2,4-Dinitrophenol           | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| 4-Nitrophenol               | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| Dibenzofuran                | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| 2,4-Dinitrotoluene          | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| Diethylphthalate            | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| 4-Chlorophenyl-phenylether  | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| Fluorene                    | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| 4-Nitroaniline              | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| 4,6-Dinitro-2-methylphenol  | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| N-Nitrosodiphenylamine      | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| 4-Bromophenyl-phenylether   | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| Hexachlorobenzene           | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| Pentachlorophenol           | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| Phenanthrene                | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | 680            | J 1700       |
| Anthracene                  | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| Carbazole                   | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| Di-n-butylphthalate         | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| Fluoranthene                | U              | 330          | 2400           | J 4400       | 1300           | J 3800       | 1000           | J 3700       | U              | 1700         |
| Pyrene                      | U              | 330          | 2200           | J 4400       | 1200           | J 3800       | 880            | J 3700       | U              | 1700         |
| Butylbenzylphthalate        | U              | 330          | U              | 4400         | U              | 3800         | 6000           | 3700         | U              | 1700         |
| Benzo(a)anthracene          | U              | 330          | 1500           | J 4400       | U              | 3800         | U              | 3700         | U              | 1700         |
| 3,3'-Dichlorobenzidine      | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| Chrysene                    | U              | 330          | 1900           | J 4400       | 1000           | J 3800       | U              | 3700         | U              | 1700         |
| Bis(2-Ethylhexyl)phthalate  | U              | 330          | U              | 4400         | 8700           | 3800         | 1200           | J 3700       | 1300           | J 1700       |
| Di-n-octylphthalate         | U              | 330          | U              | 4400         | U              | 3800         | U              | 3700         | U              | 1700         |
| Benzo(b)fluoranthene        | U              | 330          | 3100           | J 4400       | 1600           | J 3800       | 940            | J 3700       | U              | 1700         |
| Benzo(k)fluoranthene        | U              | 330          | 2400           | J 4400       | 1400           | J 3800       | 830            | J 3700       | U              | 1700         |
| Benzo(a)pyrene              | U              | 330          | 3300           | J 4400       | 1400           | J 3800       | 920            | J 3700       | U              | 1700         |
| Indeno(1,2,3-cd)pyrene      | U              | 330          | 3200           | J 4400       | 1700           | J 3800       | 770            | J 3700       | U              | 1700         |
| Dibenzo(a,h)anthracene      | U              | 330          | 1200           | J 4400       | U              | 3800         | U              | 3700         | U              | 1700         |
| Benzo(g,h,i)perylene        | U              | 330          | 3600           | J 4400       | 1900           | J 3800       | 920            | J 3700       | U              | 1700         |

ANALYTICAL REPORT

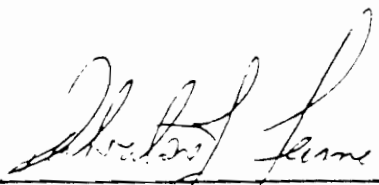
Prepared by  
Roy F. Weston, Inc.


Lackawanna Foundry Site  
Lackawanna, NY

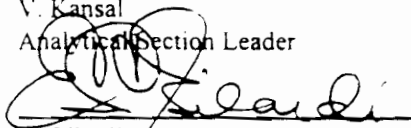
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### Introduction

REAC, in response to WA 3-439, provided analytical support for environmental samples collected from Lackawanna Foundry Site located in Lackawanna, NY as described in the following table. The support also included QA/QC, data review, and preparation of an analytical report containing a summary of the analytical methods, the results, and the QA/QC results.

The samples were treated with procedures consistent with those described in SOP # 1008 and are summarized in the following table:

| COC # | Number of Samples | Sampling Date | Date Received | Matrix | Analysis | Laboratory |
|-------|-------------------|---------------|---------------|--------|----------|------------|
| 06505 | 16                | 4/14/99       | 4/19/99       | Soil   | VOC      | Paragon    |
| 06507 | 16                | 4/14/99       | 4/19/99       | Soil   | VOC      | Paragon    |
| 06908 | -                 | 4/14/99       | 4/15/99       | Soil   | Pest/PCB | REAC       |
| 06909 | -                 | 4/14/99       | 4/15/99       | Soil   | Pest/PCB | REAC       |
| 06910 | 2                 | 4/14/99       | 4/15/99       | Soil   | Pest/PCB | REAC       |
| 06911 | 7                 | 4/14/99       | 4/15/99       | Soil   | Pest/PCB | REAC       |
| 06912 | -                 | 4/14/99       | 4/15/99       | Soil   | Pest/PCB | REAC       |

Analyses requested on these chains of custody, but not presented in this report, were given in previous reports.

## Case Narrative

The data in this report have been validated to two significant figures. Any other representation of the data is the responsibility of the user.

### VOC in Soil Package 1 200

The method blank of 4/23/99 contained 1.5 µg/kg methylene chloride. The methylene chloride in samples 16650, 16651, 16652, 16653, 16655, 16656, 16657 and 16658 should be regarded as not detected because the concentration of methylene chloride in these samples is less than ten times as much as that of the method blank.

The method blank of 4/26/99 contained 2.1 µg/kg methylene chloride. The methylene chloride in samples 16654, 16661, 16662, 16664, 16660, 16659 and 16663 should be regarded as not detected because the concentration of methylene chloride in these samples is less than ten times as much as that of the method blank.

The method blank of 4/27/99 contained 3.6 µg/kg methylene chloride. The methylene chloride in samples 16666, 16667, 16665, 16668, 16669, 16670, 16671, 16675, 16676, 16677, 16679, 16680, 16678, 16672, 16673 and 16674 should be regarded as not detected because the concentration of methylene chloride in these samples is less than ten times as much as that of the method blank.

In the initial calibration of 4/21/99 the acceptable QC limits were exceeded by iodomethane (43%) and 2-chloroethyl vinyl ether (32%). The data are not affected because these compounds were not detected in the associated samples.

In the continuing calibration check standard of 4/22/99 the acceptable QC limits were exceeded by acetone (38%), iodomethane (57%), 2-butanone (34%), dibromomethane (30%), cis-1,3-dichloropropene (40%), 4-methyl-2-pentanone (27%), trans-1,3-dichloropropene (28%), 2-hexanone (41%), 1,2-dibromoethane (27%) and 1,2-dibromo-3-chloropropane (25.6%). The data for acetone, 2-butanone, dibromomethane, cis-1,3-dichloropropene, 4-methyl-2-pentanone, trans-1,3-dichloropropene, 2-hexanone, 1,2-dibromoethane and 1,2-dibromo-3-chloropropane in samples B 16650, B 16651, B 16652, B 16653, B 16655, B 16656, B 16657 and B 16658 are not affected. The data for iodomethane in samples B 16650, B 16651, B 16652, B 16653, B 16655, B 16656, B 16657 and B 16658 should be regarded as estimated.

In the continuing calibration check standard of 4/26/99 the acceptable QC limits were exceeded by chloromethane (28%), iodomethane (53%), carbon tetrachloride (38%), cis-1,3-dichloropropene (53%), trans-1,3-dichloropropene (38%), 1,3-dichloropropane (28%), dibromochloromethane (33%) and 1,2-dibromoethane (37%). The data for chloromethane, carbon tetrachloride, trans-1,3-dichloropropene, 1,3-dichloropropane, dibromochloromethane and 1,2-dibromoethane in samples B 16654, B 16661, B 16662, B 16664, B 16659, B 16660 and B 16663 are not affected. The data for iodomethane and cis-1,3-dichloropropene in samples B 16654, B 16661, B 16662, B 16664, B 16659, B 16660 and B 16663 should be regarded as estimated.

In the continuing calibration check standard of 4/27/99 the acceptable QC limits were exceeded by iodomethane (68%), carbon tetrachloride (36%), cis-1,3-dichloropropene (49%), trans-1,3-dichloropropene (48%), 2-hexanone (27%), 1,3-dichloropropane (33%), dibromoethane (36%) and 1,2-dibromoethane (57%). The data for carbon tetrachloride, cis-1,3-dichloropropene, trans-1,3-dichloropropene, 2-hexanone, 1,3-dichloropropane, and dibromoethane (36%) in samples B 16666, B 16667, B 16665, B 16668, B 16669, B 16670, B 16671, B 16675, B 16676, B 16677, B 16679, B 16680, B 16678, B 16672, B 16673 and B 16674 are not affected. The data for iodomethane and 1,2-dibromoethane in samples B 16666, B 16667, B 16665, B 16668, B 16669, B 16670, B 16671, B 16675,

B 16676, B 16677, B 16679, B 16680, B 16678, B 16672, B 16673 and B 16674 should be regarded as estimated.

The percent recoveries of one or more surrogates exceeded the acceptable QC limits for samples B 16669, B 16670, B 16675, B 16673, B 16674 and B 16665 MS. All results should be regarded as estimated.

The areas of two or fewer internal standards exceeded the acceptable QC limits for samples B 16653, B 16656, B 16658, B 16661, B 16660, B 16671, B 16672 and B 16673. The data that were quantified under these internal standards should be regarded as estimated.

The areas of all four internal standards exceeded the acceptable QC limits for samples B 16675 and B 16677. The data that were quantified under these internal standards should be regarded as estimated.

Only one MS/MSD pair was analyzed. The subcontract laboratory did not analyze an additional MS/MSD pair as required.

#### Pesticides/PCBs in Soil Package I 194

In the continuing calibration check standard of 4/29/99, the acceptable QC limits were exceeded for one of the five peaks used for Aroclor 1260 (26%). The data are not affected because the average of the five peaks was 10.6%.

In the end of sequence check standard of 4/20/99, the acceptable QC limits were exceeded for heptachlor (44%), endosulfan II (29%), p,p'-DDT (94%), endrin aldehyde (44%), endosulfan sulfate (29%), methoxychlor (100%), endrin ketone (56%) and decachlorobiphenyl (69%). The data are not affected because this calibration was not used to quantify any analytes.

In the end of sequence check standard of 4/23/99, the acceptable QC limits were exceeded for  $\delta$ -BHC (29%), p,p'-DDD (36%), p,p'-DDT (81%), methoxychlor (77%), endrin ketone (33%) and decachlorobiphenyl (53%). The data are not affected because this calibration was not used to quantify any analytes.

In the end of sequence check standard of 4/26/99, the acceptable QC limits were exceeded for  $\beta$ -BHC (30%), heptachlor (52%),  $\gamma$ -chlordane (28%),  $\alpha$ -chlordane (29%), endosulfan I (28%), p,p'-DDE (34%), dieldrin (30%), endrin (46%), endosulfan II (39%), p,p'-DDT (92%), endrin aldehyde (45%), endosulfan sulfate (33%), methoxychlor (87%), endrin ketone (52%) and decachlorobiphenyl (59%). The data are not affected because this calibration was not used to quantify any analytes.

In the end of sequence check standard of 4/28/99, the acceptable QC limits were exceeded for the average of the five peaks used to quantify Aroclor 1254 (74%). The data are not affected because this calibration was not used to quantify Aroclor 1254 in any of the samples.

In the end of sequence check standard of 4/29/99, the acceptable QC limits were exceeded for the average of the five peaks used to quantify Aroclor 1260 (27.2%), specifically, peak 4 was 32% and peak 5 was 52%. The data are not affected because this calibration was not used to quantify Aroclor 1254 in any of the samples.

In the end of sequence check standard of 4/28/99, the acceptable QC limits were exceeded for peak 5 (26%) of Aroclor 1260. The data are not affected because this calibration was not used to quantify Aroclor 1260 in any of the samples and the average of the five peaks was 9.6%.

The acceptable percent recoveries were exceeded for both surrogates for samples A 16650 and A 16664. The data for these samples should be regarded as estimated.



The acceptable percent recoveries were exceeded for the surrogate decachlorobiphenyl for samples A 16652, A 16653, A 16653 MS, A 16653 MSD, A 16654, A 16655, A 16656, A 16658, A 16660, A 16661, A 16662, A 16663, A 16665, A 16665 MS, A 16665 MSD, A 16666, A 16667, A 16668, A 16672, A 16674, A 16675, A 16676, A 16677, A 16678, A 16679, A 16677 MS and A 16677 MSD. The data for these samples are not affected.

Additional Aroclors are likely to be present in various samples such as A 16672, A 16673, A 16674, A 16677, A 16679. The laboratory did not report these in the final results because their concentrations were below the detection limit and very low in comparison to the Aroclors that were reported.

### Summary of Abbreviations

|                |   |    |           |    |           |
|----------------|---|----|-----------|----|-----------|
| AA             | Atomic Absorption   |    |           |    |           |
| B              | The analyte was found in the blank  |    |           |    |           |
| BFB            | Bromofluorobenzene  |    |           |    |           |
| C              | Centigrade  |    |           |    |           |
| D              | (Surrogate Table) this value is from a diluted sample and was not calculated<br>(Result Table) this result was obtained from a diluted sample |    |           |    |           |
| Dioxin         | denotes Polychlorinated Dibenzop-dioxins and Polychlorinated Dibenzofurans and/or   |    |           |    |           |
| PCDD           | and PCDF  |    |           |    |           |
| CLP            | Contract Laboratory Protocol  |    |           |    |           |
| COC            | Chain of Custody  |    |           |    |           |
| CONC           | Concentration   |    |           |    |           |
| CRDL           | Contract Required Detection Limit   |    |           |    |           |
| CRQL           | Contract Required Quantitation Limit  |    |           |    |           |
| DFTPP          | Decafluorotriphenylphosphine  |    |           |    |           |
| DL             | Detection Limit   |    |           |    |           |
| E              | The value is greater than the highest linear standard and is estimated  |    |           |    |           |
| EMPC           | Estimated maximum possible concentration  |    |           |    |           |
| ICAP           | Inductively Coupled Argon Plasma  |    |           |    |           |
| ISTD           | Internal Standard   |    |           |    |           |
| J              | The value is below the method detection limit and is estimated  |    |           |    |           |
| LCS            | Laboratory Control Sample   |    |           |    |           |
| LCSD           | Laboratory Control Sample Duplicate   |    |           |    |           |
| MDL            | Method Detection Limit  |    |           |    |           |
| MI             | Matrix Interference   |    |           |    |           |
| MS             | Matrix Spike  |    |           |    |           |
| MSD            | Matrix Spike Duplicate  |    |           |    |           |
| MW             | Molecular Weight  |    |           |    |           |
| NA             | either Not Applicable or Not Available  |    |           |    |           |
| NC             | Not Calculated  |    |           |    |           |
| NR             | Not Requested   |    |           |    |           |
| NS             | Not Spiked  |    |           |    |           |
| % D            | Percent Difference  |    |           |    |           |
| % REC          | Percent Recovery  |    |           |    |           |
| PPB            | Parts per billion   |    |           |    |           |
| PPBV           | Parts per billion by volume   |    |           |    |           |
| PQL            | Practical Quantitation Limit  |    |           |    |           |
| QL             | Quantitation Limit  |    |           |    |           |
| RPD            | Relative Percent Difference   |    |           |    |           |
| RSD            | Relative Standard Deviation   |    |           |    |           |
| SIM            | Selected Ion Monitoring   |    |           |    |           |
| TCLP           | Toxic Characteristics Leaching Procedure  |    |           |    |           |
| U              | Denotes not detected  |    |           |    |           |
| W              | Weathered sample; the results should be regarded as estimated   |    |           |    |           |
| m <sup>3</sup> | cubic meter   | kg | kilogram  | μg | microgram |
| L              | liter   | g  | gram      | pg | picogram  |
| mL             | milliliter  | mg | milligram | ng | nanogram  |
| μL             | microliter  |    |           |    |           |
| *              | denotes a value that exceeds the acceptable QC limit  |    |           |    |           |
|                | Abbreviations that are specific to a particular table are explained in footnotes on that table  |    |           |    |           |

Revision 7/23/98

### Analytical Procedure for VOC in Soil

The subcontract laboratory determined the concentration of volatile organic compounds in the samples by analyzing them by modified USEPA Method 8260 B found in SW-846. The results of the analyses are listed in Table 1.1.

## Analytical Procedure for Pesticides/PCBs in Soil

### Extraction Procedure

The soil samples were extracted by the Soxhlet method. A thirty gram aliquot was spiked with a surrogate solution consisting of tetrachloro-m-xylene and decachlorobiphenyl, mixed with 30 g anhydrous sodium sulfate and Soxhlet extracted for 16 hours with 300-mL hexane. The extract was concentrated to 5 mL.

### Gas Chromatographic Analysis - Pesticide Quantitation and PCB Screening

The extract was analyzed for pesticides using simultaneous dual column injections. The analysis was done on an HP-6890, equipped with an HP-6890 automatic sampler, and controlled with an HP Chem-Station. The following conditions were employed:

|                                    |   |
|------------------------------------|---|
| First Column                       | DB-608, 30 meter, 0.32 mm fused silica capillary, 0.50 $\mu$ m film thickness           |
| Injector Temperature               | 200° C  |
| Detector Temperature               | 325° C  |
| Second Column                      | Rtx-CLPesticides, 30 meter, 0.53 mm fused silica capillary, 0.50 $\mu$ m film thickness |
| Injector Temperature               | 200° C  |
| Detector Temperature               | 325° C  |
| Temperature Program-(both columns) | 120 ° C for 1 minute<br>9 °C/min to 285°C, 10 min at 285°C                              |
| Injection Volume                   | 1 $\mu$ L   |

The gas chromatographs were calibrated using 5 pesticide standards at 20, 50, 100, 200, and 500  $\mu$ g/L. The results from each mixture were used to calculate the response factor (RF) of each analyte and the average Response Factor was used to calculate the concentration of pesticide in the sample. Quantification was based on the DB-608 column (signal 1) and the identity of the analyte was confirmed using the Rtx-CLPesticides column (signal 2). A fingerprint chromatogram was run using seven Aroclors and toxaphene; calibration curves were run only if a particular Aroclor or toxaphene was found in the sample.

The pesticide results, listed in Table 1.2, are calculated by using the following formula:

$$C_u = \frac{DF \times A_u \times V_i}{RF_{ave} \times V_i \times W \times D}$$

where

- $C_u$  = Concentration of analyte ( $\mu\text{g}/\text{kg}$ )
- DF = Dilution Factor
- $A_u$  = Area or peak height
- $V_i$  = Volume of sample (mL)
- $RF_{ave}$  = Average response factor
- $V_i$  = Volume of extract injected ( $\mu\text{L}$ )
- W = Weight of sample (g)
- D = Decimal percent solids

Response Factor calculation:

The RF for each specific analyte is quantitated based on the area response from the continuing calibration check as follows:

$$RF = \frac{A_u}{\text{total pg injected}}$$

where

$A_u$  = Area or peak height

and

$$RF_{ave} = \frac{RF_1 + \dots + RF_n}{n}$$

where

n = number of samples

Revision 7/22/97

Table 1.1 Results of the Analysis for VOC in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 (Results are Based on Dry Weight)

| Sample No.                 | Method Blank   |              | B16650         |              | B16651            |              | B16652         |              | B16653         |              |    |     |     |    |     |
|----------------------------|----------------|--------------|----------------|--------------|-------------------|--------------|----------------|--------------|----------------|--------------|----|-----|-----|----|-----|
| Sample Location            | 4/23/99        |              | 1 West Fence   |              | 2D Surface Middle |              | 2D Middle      |              | 3 West Fence   |              |    |     |     |    |     |
| % Solid                    | NA             |              | 65             |              | 93                |              | 92             |              | 76             |              |    |     |     |    |     |
| Compound Name              | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg    | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg |    |     |     |    |     |
| Dichlorodifluoromethane    | U              | 10           | U              | 15           | U                 | 11           | U              | 11           | U              | 13           |    |     |     |    |     |
| Chloromethane              | U              | 10           | U              | 15           | U                 | 11           | U              | 11           | U              | 13           |    |     |     |    |     |
| Vinyl chloride             | U              | 10           | U              | 15           | U                 | 11           | U              | 11           | U              | 13           |    |     |     |    |     |
| Bromomethane               | U              | 10           | U              | 15           | U                 | 11           | U              | 11           | U              | 13           |    |     |     |    |     |
| Chloroethane               | U              | 10           | U              | 15           | U                 | 11           | U              | 11           | U              | 13           |    |     |     |    |     |
| Trichlorofluoromethane     | U              | 5            | U              | 7.7          | U                 | 5.4          | U              | 5.4          | U              | 6.6          |    |     |     |    |     |
| Acrolein                   | U              | 50           | U              | 77           | U                 | 54           | U              | 54           | U              | 66           |    |     |     |    |     |
| 1,1-Dichloroethylene       | U              | 5            | U              | 7.7          | U                 | 5.4          | U              | 5.4          | U              | 6.6          |    |     |     |    |     |
| Trichlorotrifluoroethane   | U              | 5            | U              | 7.7          | U                 | 5.4          | U              | 5.4          | U              | 6.6          |    |     |     |    |     |
| Acetone                    | U              | 20           | U              | 31           | U                 | 21           | U              | 21           | U              | 26           |    |     |     |    |     |
| Iodomethane                | U              | 5            | U              | 7.7          | U                 | 5.4          | U              | 5.4          | U              | 6.6          |    |     |     |    |     |
| Carbon disulfide           | U              | 5            | U              | 7.7          | U                 | 5.4          | U              | 5.4          | U              | 6.6          |    |     |     |    |     |
| Methylene chloride         | 1.5            | J            | 5              | 1.9          | JB                | 7.7          | 1.4            | JB           | 5.4            | 1.2          | JB | 5.4 | 1.8 | JB | 6.6 |
| trans-1,2-Dichloroethylene | U              | 5            | U              | 7.7          | U                 | 5.4          | U              | 5.4          | U              | 6.6          |    |     |     |    |     |
| Methyl t-butyl ether       | U              | 5            | U              | 7.7          | U                 | 5.4          | U              | 5.4          | U              | 6.6          |    |     |     |    |     |
| Acrylonitrile              | U              | 50           | U              | 77           | U                 | 54           | U              | 54           | U              | 66           |    |     |     |    |     |
| 1,1-Dichloroethane         | U              | 5            | U              | 7.7          | U                 | 5.4          | U              | 5.4          | U              | 6.6          |    |     |     |    |     |
| Vinyl acetate              | U              | 20           | U              | 31           | U                 | 21           | U              | 21           | U              | 26           |    |     |     |    |     |
| cis-1,2-Dichloroethylene   | U              | 5            | U              | 7.7          | U                 | 5.4          | U              | 5.4          | U              | 6.6          |    |     |     |    |     |
| 2-Butanone                 | U              | 20           | U              | 31           | U                 | 21           | U              | 21           | U              | 26           |    |     |     |    |     |
| Bromochloromethane         | U              | 5            | U              | 7.7          | U                 | 5.4          | U              | 5.4          | U              | 6.6          |    |     |     |    |     |
| Chloroform                 | U              | 5            | U              | 7.7          | U                 | 5.4          | U              | 5.4          | U              | 6.6          |    |     |     |    |     |
| 1,1,1-Trichloroethane      | U              | 5            | U              | 7.7          | U                 | 5.4          | U              | 5.4          | U              | 6.6          |    |     |     |    |     |
| 2,2-Dichloropropane        | U              | 5            | U              | 7.7          | U                 | 5.4          | U              | 5.4          | U              | 6.6          |    |     |     |    |     |
| Carbon tetrachloride       | U              | 5            | U              | 7.7          | U                 | 5.4          | U              | 5.4          | U              | 6.6          |    |     |     |    |     |
| 1,1-Dichloropropene        | U              | 5            | U              | 7.7          | U                 | 5.4          | U              | 5.4          | U              | 6.6          |    |     |     |    |     |
| 1,2-Dichloroethane         | U              | 5            | U              | 7.7          | U                 | 5.4          | U              | 5.4          | U              | 6.6          |    |     |     |    |     |
| Benzene                    | U              | 5            | U              | 7.7          | U                 | 5.4          | U              | 5.4          | U              | 6.6          |    |     |     |    |     |
| Trichloroethylene          | U              | 5            | U              | 7.7          | U                 | 5.4          | U              | 5.4          | U              | 6.6          |    |     |     |    |     |
| 1,2-Dichloropropane        | U              | 5            | U              | 7.7          | U                 | 5.4          | U              | 5.4          | U              | 6.6          |    |     |     |    |     |
| Dibromomethane             | U              | 5            | U              | 7.7          | U                 | 5.4          | U              | 5.4          | U              | 6.6          |    |     |     |    |     |
| Bromochloromethane         | U              | 5            | U              | 7.7          | U                 | 5.4          | U              | 5.4          | U              | 6.6          |    |     |     |    |     |
| 2-Chloroethyl vinyl ether  | U              | 10           | U              | 15           | U                 | 11           | U              | 11           | U              | 13           |    |     |     |    |     |
| cis-1,3-Dichloropropylene  | U              | 5            | U              | 7.7          | U                 | 5.4          | U              | 5.4          | U              | 6.6          |    |     |     |    |     |
| 4-Methyl-2-pentanone       | U              | 20           | U              | 31           | U                 | 21           | U              | 21           | U              | 26           |    |     |     |    |     |
| Toluene                    | U              | 5            | U              | 7.7          | U                 | 5.4          | U              | 5.4          | U              | 6.6          |    |     |     |    |     |

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Table 1.1 Results of the Analysis for VOC in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 (Results are Based on Dry Weight)

| Sample No.<br>Sample Location<br>% Solid | Method Blank<br>4/23/99<br>NA |              | B16650<br>1 West Fence<br>65 |              | B16651<br>2D Surface Middle<br>93 |              | B16652<br>2D Middle<br>92 |              | B16653<br>3 West Fence<br>76 |              |
|--|-------------------------------|--------------|------------------------------|--------------|-----------------------------------|--------------|---------------------------|--------------|------------------------------|--------------|
| Compound Name                            | Conc.<br>µg/kg                | MDL<br>µg/kg | Conc.<br>µg/kg               | MDL<br>µg/kg | Conc.<br>µg/kg                    | MDL<br>µg/kg | Conc.<br>µg/kg            | MDL<br>µg/kg | Conc.<br>µg/kg               | MDL<br>µg/kg |
| trans-1,3-Dichloropropylene              | U                             | 5            | U                            | 7.7          | U                                 | 5.4          | U                         | 5.4          | U                            | 6.6          |
| 1,1,2-Trichloroethane                    | U                             | 5            | U                            | 7.7          | U                                 | 5.4          | U                         | 5.4          | U                            | 6.6          |
| 2-Hexanone                               | U                             | 20           | U                            | 31           | U                                 | 21           | U                         | 21           | U                            | 26           |
| Tetrachloroethylene                      | U                             | 5            | U                            | 7.7          | U                                 | 5.4          | U                         | 5.4          | U                            | 6.6          |
| 1,3-Dichloropropane                      | U                             | 5            | U                            | 7.7          | U                                 | 5.4          | U                         | 5.4          | U                            | 6.6          |
| Dibromochloroethane                      | U                             | 5            | U                            | 7.7          | U                                 | 5.4          | U                         | 5.4          | U                            | 6.6          |
| 1,2-Dibromoethane                        | U                             | 5            | U                            | 7.7          | U                                 | 5.4          | U                         | 5.4          | U                            | 6.6          |
| 1-Chlorohexane                           | U                             | 5            | U                            | 7.7          | U                                 | 5.4          | U                         | 5.4          | U                            | 6.6          |
| Chlorobenzene                            | U                             | 5            | U                            | 7.7          | U                                 | 5.4          | U                         | 5.4          | U                            | 6.6          |
| 1,1,1,2-Tetrachloroethane                | U                             | 5            | U                            | 7.7          | U                                 | 5.4          | U                         | 5.4          | U                            | 6.6          |
| Ethyl benzene                            | U                             | 5            | U                            | 7.7          | U                                 | 5.4          | U                         | 5.4          | U                            | 6.6          |
| mp- Xylene                               | U                             | 5            | U                            | 7.7          | U                                 | 5.4          | U                         | 5.4          | U                            | 6.6          |
| o- Xylene                                | U                             | 5            | U                            | 7.7          | U                                 | 5.4          | U                         | 5.4          | U                            | 6.6          |
| Styrene                                  | U                             | 5            | U                            | 7.7          | U                                 | 5.4          | U                         | 5.4          | U                            | 6.6          |
| Bromoform                                | U                             | 5            | U                            | 7.7          | U                                 | 5.4          | U                         | 5.4          | U                            | 6.6          |
| Isopropylbenzene                         | U                             | 5            | U                            | 7.7          | U                                 | 5.4          | U                         | 5.4          | U                            | 6.6          |
| 1,2,3-Trichloropropane                   | U                             | 5            | U                            | 7.7          | U                                 | 5.4          | U                         | 5.4          | U                            | 6.6          |
| 1,1,2,2-Tetrachloroethane                | U                             | 5            | U                            | 7.7          | U                                 | 5.4          | U                         | 5.4          | U                            | 6.6          |
| Bromobenzene                             | U                             | 5            | U                            | 7.7          | U                                 | 5.4          | U                         | 5.4          | U                            | 6.6          |
| n-Propylbenzene                          | U                             | 5            | U                            | 7.7          | U                                 | 5.4          | U                         | 5.4          | U                            | 6.6          |
| 2-Chlorotoluene                          | U                             | 5            | U                            | 7.7          | U                                 | 5.4          | U                         | 5.4          | U                            | 6.6          |
| 1,3,5-Trimethylbenzene                   | U                             | 5            | U                            | 7.7          | U                                 | 5.4          | U                         | 5.4          | U                            | 6.6          |
| 4-Chlorotoluene                          | U                             | 5            | U                            | 7.7          | U                                 | 5.4          | U                         | 5.4          | U                            | 6.6          |
| t-Butylbenzene                           | U                             | 5            | U                            | 7.7          | U                                 | 5.4          | U                         | 5.4          | U                            | 6.6          |
| 1,2,4-Trimethylbenzene                   | U                             | 5            | U                            | 7.7          | U                                 | 5.4          | U                         | 5.4          | U                            | 6.6          |
| sec-Butylbenzene                         | U                             | 5            | U                            | 7.7          | U                                 | 5.4          | U                         | 5.4          | U                            | 6.6          |
| 1,3-Dichlorobenzene                      | U                             | 5            | U                            | 7.7          | U                                 | 5.4          | U                         | 5.4          | U                            | 6.6          |
| p-Isopropyltoluene                       | U                             | 5            | U                            | 7.7          | U                                 | 5.4          | U                         | 5.4          | U                            | 6.6          |
| 1,4-Dichlorobenzene                      | U                             | 5            | U                            | 7.7          | U                                 | 5.4          | U                         | 5.4          | U                            | 6.6          |
| n-Butylbenzene                           | U                             | 5            | U                            | 7.7          | U                                 | 5.4          | U                         | 5.4          | U                            | 6.6          |
| 1,2-Dichlorobenzene                      | U                             | 5            | U                            | 7.7          | U                                 | 5.4          | U                         | 5.4          | U                            | 6.6          |
| 1,2-Dibromo-3-chloropropane              | U                             | 10           | U                            | 15           | U                                 | 11           | U                         | 11           | U                            | 13           |
| 1,2,4-Trichlorobenzene                   | U                             | 5            | U                            | 7.7          | U                                 | 5.4          | U                         | 5.4          | U                            | 6.6          |
| Hexachlorobutadiene                      | U                             | 5            | U                            | 7.7          | U                                 | 5.4          | U                         | 5.4          | U                            | 6.6          |
| Naphthalene                              | U                             | 5            | U                            | 7.7          | U                                 | 5.4          | U                         | 5.4          | 1.1                          | J 6.6        |
| 1,2,3-Trichlorobenzene                   | U                             | 5            | U                            | 7.7          | U                                 | 5.4          | U                         | 5.4          | U                            | 6.6          |

00010

Table 1.1 (Cont) Results of the Analysis for VOC in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 (Results are Based on Dry Weight)

| Sample No.<br>Sample Location<br>% Solid | B16655<br>5 West Fence<br>82 |              | B16656<br>5 Rear Entrance<br>86 |              | B16657<br>7<br>82 |              | B16658<br>8<br>62 |              |
|--|------------------------------|--------------|---------------------------------|--------------|-------------------|--------------|-------------------|--------------|
|  | Conc.<br>µg/kg               | MDL<br>µg/kg | Conc.<br>µg/kg                  | MDL<br>µg/kg | Conc.<br>µg/kg    | MDL<br>µg/kg | Conc.<br>µg/kg    | MDL<br>µg/kg |
| Dichlorodifluoromethane                  | U                            | 12           | U                               | 12           | U                 | 12           | U                 | 16           |
| Chloromethane                            | U                            | 12           | U                               | 12           | U                 | 12           | U                 | 16           |
| Vinyl chloride                           | U                            | 12           | U                               | 12           | U                 | 12           | U                 | 16           |
| Bromomethane                             | U                            | 12           | U                               | 12           | U                 | 12           | U                 | 16           |
| Chloroethane                             | U                            | 12           | U                               | 12           | U                 | 12           | U                 | 16           |
| Trichlorofluoromethane                   | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| Acrolein                                 | U                            | 61           | U                               | 58           | U                 | 61           | U                 | 81           |
| 1,1-Dichloroethylene                     | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| Trichlorotrifluoroethane                 | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| Acetone                                  | U                            | 24           | U                               | 23           | U                 | 24           | U                 | 32           |
| Iodomethane                              | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| Carbon disulfide                         | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| Methylene chloride                       | 1.6 JB                       | 6.1          | 1.6 JB                          | 5.8          | 1.4 JB            | 6.1          | 2.4 JB            | 8.1          |
| trans-1,2-Dichloroethylene               | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| Methyl t-butyl ether                     | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| Acrylonitrile                            | U                            | 61           | U                               | 58           | U                 | 61           | U                 | 81           |
| 1,1-Dichloroethane                       | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| Vinyl acetate                            | U                            | 24           | U                               | 23           | U                 | 24           | U                 | 32           |
| cis-1,2-Dichloroethylene                 | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| 2-Butanone                               | U                            | 24           | U                               | 23           | U                 | 24           | U                 | 32           |
| Bromochloromethane                       | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| Chloroform                               | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| 1,1,1,-Trichloroethane                   | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| 2,2-Dichloropropane                      | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| Carbon tetrachloride                     | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| 1,1-Dichloropropene                      | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| 1,2-Dichloroethane                       | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| Benzene                                  | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| Trichloroethylene                        | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| 1,2-Dichloropropane                      | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| Dibromomethane                           | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| Bromochloromethane                       | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| 2-Chloroethyl vinyl ether                | U                            | 12           | U                               | 12           | U                 | 12           | U                 | 16           |
| cis-1,3-Dichloropropylene                | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| 4-Methyl-2-pentanone                     | U                            | 24           | U                               | 23           | U                 | 24           | U                 | 32           |
| Toluene                                  | U                            | 6.1          | 4.0 J                           | 5.8          | U                 | 6.1          | U                 | 8.1          |

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Table 1.1 (Cont) Results of the Analysis for VOC in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 (Results are Based on Dry Weight)

| Sample No.<br>Sample Location<br>% Solid | B16655<br>5 West Fence<br>82 |              | B16656<br>5 Rear Entrance<br>86 |              | B16657<br>7<br>82 |              | B16658<br>8<br>62 |              |
|--|------------------------------|--------------|---------------------------------|--------------|-------------------|--------------|-------------------|--------------|
|  | Conc.<br>µg/kg               | MDL<br>µg/kg | Conc.<br>µg/kg                  | MDL<br>µg/kg | Conc.<br>µg/kg    | MDL<br>µg/kg | Conc.<br>µg/kg    | MDL<br>µg/kg |
| trans-1,3-Dichloropropylene              | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| 1,1,2-Trichloroethane                    | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| 2-Hexanone                               | U                            | 24           | U                               | 23           | U                 | 24           | U                 | 32           |
| Tetrachloroethylene                      | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| 1,3-Dichloropropane                      | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| Dibromochloromethane                     | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| 1,2-Dibromoethane                        | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| 1-Chlorohexane                           | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| Chlorobenzene                            | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| 1,1,1,2-Tetrachloroethane                | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| Ethyl benzene                            | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| mp- Xylene                               | U                            | 6.1          | 1.8 J                           | 5.8          | U                 | 6.1          | U                 | 8.1          |
| o- Xylene                                | U                            | 6.1          | 1.0 J                           | 5.8          | U                 | 6.1          | U                 | 8.1          |
| Styrene                                  | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| Bromoform                                | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| Isopropylbenzene                         | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| 1,2,3-Trichloropropane                   | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| 1,1,2,2-Tetrachloroethane                | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| Bromobenzene                             | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| n-Propylbenzene                          | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| 2-Chlorotoluene                          | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| 1,3,5-Tmethylbenzene                     | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| 4-Chlorotoluene                          | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| t-Butylbenzene                           | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| 1,2,4-Trimethylbenzene                   | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| sec-Butylbenzene                         | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| 1,3-Dichlorobenzene                      | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| p-Isopropyltoluene                       | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| 1,4-Dichlorobenzene                      | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| n-Butylbenzene                           | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| 1,2-Dichlorobenzene                      | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| 1,2-Dibromo-3-chloropropane              | U                            | 12           | U                               | 12           | U                 | 12           | U                 | 16           |
| 1,2,4-Trichlorobenzene                   | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| Hexachlorobutadiene                      | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| Naphthalene                              | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |
| 1,2,3-Trichlorobenzene                   | U                            | 6.1          | U                               | 5.8          | U                 | 6.1          | U                 | 8.1          |

00012

Table 1.1 (Cont) Results of the Analysis for VOC in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 (Results are Based on Dry Weight)

| Sample No.                 | Method Blank   |              | B16654         |              | B16661         |              | B16662         |              | B16664         |              |    |   |     |    |     |
|----------------------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----|---|-----|----|-----|
| Sample Location            | 4/26/99        |              | 4 West Fence   |              | 11             |              | 12             |              | 14             |              |    |   |     |    |     |
| % Solid                    | NA             |              | 87             |              | 84             |              | 83             |              | 82             |              |    |   |     |    |     |
| Compound Name              | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg |    |   |     |    |     |
| Dichlorodifluoromethane    | U              | 10           | U              | 12           | U              | 12           | U              | 12           | U              | 12           |    |   |     |    |     |
| Chloromethane              | U              | 10           | U              | 12           | U              | 12           | U              | 12           | U              | 12           |    |   |     |    |     |
| Vinyl chloride             | U              | 10           | U              | 12           | U              | 12           | U              | 12           | U              | 12           |    |   |     |    |     |
| Bromomethane               | U              | 10           | U              | 12           | U              | 12           | U              | 12           | U              | 12           |    |   |     |    |     |
| Chloroethane               | U              | 10           | U              | 12           | U              | 12           | U              | 12           | U              | 12           |    |   |     |    |     |
| Trichlorofluoromethane     | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |    |   |     |    |     |
| Acrolein                   | U              | 50           | U              | 58           | U              | 60           | U              | 60           | U              | 61           |    |   |     |    |     |
| 1,1-Dichloroethylene       | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |    |   |     |    |     |
| Trichlorotrifluoroethane   | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |    |   |     |    |     |
| Acetone                    | U              | 20           | U              | 23           | U              | 24           | U              | 24           | U              | 24           |    |   |     |    |     |
| Iodomethane                | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |    |   |     |    |     |
| Carbon disulfide           | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |    |   |     |    |     |
| Methylene chloride         | 2.1            | J            | 5              | 1.8          | JB             | 5.8          | 7.7            | B            | 6              | 3.9          | JB | 6 | 3.7 | JB | 6.1 |
| trans-1,2-Dichloroethylene | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |    |   |     |    |     |
| Methyl t-butyl ether       | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |    |   |     |    |     |
| Acrylonitrile              | U              | 50           | U              | 58           | U              | 60           | U              | 60           | U              | 61           |    |   |     |    |     |
| 1,1-Dichloroethane         | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |    |   |     |    |     |
| Vinyl acetate              | U              | 20           | U              | 23           | U              | 24           | U              | 24           | U              | 24           |    |   |     |    |     |
| cis-1,2-Dichloroethylene   | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |    |   |     |    |     |
| 2-Butanone                 | U              | 20           | U              | 23           | U              | 24           | U              | 24           | U              | 24           |    |   |     |    |     |
| Bromochloromethane         | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |    |   |     |    |     |
| Chloroform                 | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |    |   |     |    |     |
| 1,1,1,-Trichloroethane     | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |    |   |     |    |     |
| 2,2-Dichloropropane        | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |    |   |     |    |     |
| Carbon tetrachloride       | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |    |   |     |    |     |
| 1,1-Dichloropropene        | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |    |   |     |    |     |
| 1,2-Dichloroethane         | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |    |   |     |    |     |
| Benzene                    | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |    |   |     |    |     |
| Trichloroethylene          | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |    |   |     |    |     |
| 1,2-Dichloropropane        | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |    |   |     |    |     |
| Dibromomethane             | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |    |   |     |    |     |
| Bromochloromethane         | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |    |   |     |    |     |
| 2-Chloroethyl vinyl ether  | U              | 10           | U              | 12           | U              | 12           | U              | 12           | U              | 12           |    |   |     |    |     |
| cis-1,3-Dichloropropylene  | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |    |   |     |    |     |
| 4-Methyl-2-pentanone       | U              | 20           | U              | 23           | U              | 24           | U              | 24           | U              | 24           |    |   |     |    |     |
| Toluene                    | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |    |   |     |    |     |

00013

Table 1.1 (Cont) Results of the Analysis for VOC in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 (Results are Based on Dry Weight)

| Sample No.                  | Method Blank   | B16654       |                | B16661       |                | B16662       |                | B16664       |                |              |
|-----------------------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|
| Sample Location             | 4/26/99        | 4 West Fence |                | 11           |                | 12           |                | 14           |                |              |
| % Solid                     | NA             | 87           |                | 84           |                | 83           |                | 82           |                |              |
| Compound Name               | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg |
| trans-1,3-Dichloropropylene | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |
| 1,1,2-Trichloroethane       | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |
| 2-Hexanone                  | U              | 20           | U              | 23           | U              | 24           | U              | 24           | U              | 24           |
| Tetrachloroethylene         | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |
| 1,3-Dichloropropane         | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |
| Dibromochloromethane        | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |
| 1,2-Dibromoethane           | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |
| 1-Chlorohexane              | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |
| Chlorobenzene               | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |
| 1,1,1,2-Tetrachloroethane   | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |
| Ethyl benzene               | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |
| mp- Xylene                  | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |
| o- Xylene                   | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |
| Styrene                     | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |
| Bromoform                   | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |
| Isopropylbenzene            | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |
| 1,2,3-Trichloropropane      | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |
| 1,1,1,2,2-Tetrachloroethane | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |
| Bromobenzene                | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |
| n-Propylbenzene             | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |
| 2-Chlorotoluene             | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |
| 1,3,5-Trimethylbenzene      | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |
| 4-Chlorotoluene             | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |
| t-Butylbenzene              | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |
| 1,2,4-Trimethylbenzene      | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |
| sec-Butylbenzene            | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |
| 1,3-Dichlorobenzene         | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |
| p-Isopropyltoluene          | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |
| 1,4-Dichlorobenzene         | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |
| n-Butylbenzene              | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |
| 1,2-Dichlorobenzene         | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |
| 1,2-Dibromo-3-chloropropane | U              | 10           | U              | 12           | U              | 12           | U              | 12           | U              | 12           |
| 1,2,4-Trichlorobenzene      | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |
| Hexachlorobutadiene         | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |
| Naphthalene                 | U              | 5            | U              | 5.8          | 1.1            | 6            | U              | 6            | U              | 6.1          |
| 1,2,3-Trichlorobenzene      | U              | 5            | U              | 5.8          | U              | 6            | U              | 6            | U              | 6.1          |

00014

Table 1.1 (Cont) Results of the Analysis for VOC in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 (Results are Based on Dry Weight)

| Sample No.                 | B16659         |              | B16660         |              | B16663         |              |
|----------------------------|----------------|--------------|----------------|--------------|----------------|--------------|
| Sample Location            | 9              |              | 10             |              | 13             |              |
| % Solid                    | 90             |              | 99             |              | 90             |              |
| Compound Name              | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg |
| Dichlorodifluoromethane    | U              | 11           | U              | 10           | U              | 11           |
| Chloromethane              | U              | 11           | U              | 10           | U              | 11           |
| Vinyl chloride             | U              | 11           | U              | 10           | U              | 11           |
| Bromomethane               | U              | 11           | U              | 10           | U              | 11           |
| Chloroethane               | U              | 11           | U              | 10           | U              | 11           |
| Trichlorofluoromethane     | U              | 5.6          | U              | 5.1          | U              | 5.5          |
| Acrolein                   | U              | 56           | U              | 51           | U              | 55           |
| 1,1-Dichloroethylene       | U              | 5.6          | U              | 5.1          | U              | 5.5          |
| Trichlorotrifluoroethane   | U              | 5.6          | U              | 5.1          | U              | 5.5          |
| Acetone                    | U              | 22           | U              | 20           | U              | 22           |
| Iodomethane                | U              | 5.6          | U              | 5.1          | U              | 5.5          |
| Carbon disulfide           | U              | 5.6          | U              | 5.1          | U              | 5.5          |
| Methylene chloride         | 4.2            | JB           | 5.6            | 4.9          | JB             | 5.5          |
| trans-1,2-Dichloroethylene | U              | 5.6          | U              | 5.1          | U              | 5.5          |
| Methyl t-butyl ether       | U              | 5.6          | U              | 5.1          | U              | 5.5          |
| Acrylonitrile              | U              | 56           | U              | 51           | U              | 55           |
| 1,1-Dichloroethane         | U              | 5.6          | U              | 5.1          | U              | 5.5          |
| Vinyl acetate              | U              | 22           | U              | 20           | U              | 22           |
| cis-1,2-Dichloroethylene   | U              | 5.6          | U              | 5.1          | U              | 5.5          |
| 2-Butanone                 | U              | 22           | U              | 20           | U              | 22           |
| Bromochloromethane         | U              | 5.6          | U              | 5.1          | U              | 5.5          |
| Chloroform                 | U              | 5.6          | U              | 5.1          | U              | 5.5          |
| 1,1,1,-Trichloroethane     | U              | 5.6          | U              | 5.1          | U              | 5.5          |
| 2,2-Dichloropropane        | U              | 5.6          | U              | 5.1          | U              | 5.5          |
| Carbon tetrachloride       | U              | 5.6          | U              | 5.1          | U              | 5.5          |
| 1,1-Dichloropropene        | U              | 5.6          | U              | 5.1          | U              | 5.5          |
| 1,2-Dichloroethane         | U              | 5.6          | U              | 5.1          | U              | 5.5          |
| Benzene                    | U              | 5.6          | 0.6            | J            | 5.1            | 5.5          |
| Trichloroethylene          | U              | 5.6          | U              | 5.1          | U              | 5.5          |
| 1,2-Dichloropropane        | U              | 5.6          | U              | 5.1          | U              | 5.5          |
| Dibromomethane             | U              | 5.6          | U              | 5.1          | U              | 5.5          |
| Bromochloromethane         | U              | 5.6          | U              | 5.1          | U              | 5.5          |
| 2-Chloroethyl vinyl ether  | U              | 11           | U              | 10           | U              | 11           |
| cis-1,3-Dichloropropylene  | U              | 5.6          | U              | 5.1          | U              | 5.5          |
| 4-Methyl-2-pentanone       | U              | 22           | U              | 20           | U              | 22           |
| Toluene                    | U              | 5.6          | 27             | 5.1          | U              | 5.5          |

Table 1.1 (Cont) Results of the Analysis for VOC in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 (Results are Based on Dry Weight)

| Sample No.                  | B16659         |              | B16660         |              | B16663         |              |     |
|-----------------------------|----------------|--------------|----------------|--------------|----------------|--------------|-----|
| Sample Location             | 9              |              | 10             |              | 13             |              |     |
| % Solid                     | 90             |              | 99             |              | 90             |              |     |
| Compound Name               | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg |     |
| trans-1,3-Dichloropropylene | U              | 5.6          | U              | 5.1          | U              | 5.5          |     |
| 1,1,2-Trichloroethane       | U              | 5.6          | U              | 5.1          | U              | 5.5          |     |
| 2-Hexanone                  | U              | 22           | U              | 20           | U              | 22           |     |
| Tetrachloroethylene         | U              | 5.6          | U              | 5.1          | U              | 5.5          |     |
| 1,3-Dichloropropane         | U              | 5.6          | U              | 5.1          | U              | 5.5          |     |
| Dibromochloromethane        | U              | 5.6          | U              | 5.1          | U              | 5.5          |     |
| 1,2-Dibromoethane           | U              | 5.6          | U              | 5.1          | U              | 5.5          |     |
| 1-Chlorohexane              | U              | 5.6          | U              | 5.1          | U              | 5.5          |     |
| Chlorobenzene               | U              | 5.6          | U              | 5.1          | U              | 5.5          |     |
| 1,1,1,2-Tetrachloroethane   | U              | 5.6          | U              | 5.1          | U              | 5.5          |     |
| Ethyl benzene               | U              | 5.6          | 3.6            | J            | 5.1            | U            | 5.5 |
| mp- Xylene                  | U              | 5.6          | 13             | J            | 5.1            | U            | 5.5 |
| o- Xylene                   | U              | 5.6          | 4.7            | J            | 5.1            | U            | 5.5 |
| Styrene                     | U              | 5.6          | U              | 5.1          | U              | 5.5          |     |
| Bromofom                    | U              | 5.6          | U              | 5.1          | U              | 5.5          |     |
| Isopropylbenzene            | U              | 5.6          | U              | 5.1          | U              | 5.5          |     |
| 1,2,3-Trichloropropane      | U              | 5.6          | U              | 5.1          | U              | 5.5          |     |
| 1,1,1,2-Tetrachloroethane   | U              | 5.6          | U              | 5.1          | U              | 5.5          |     |
| Bromobenzene                | U              | 5.6          | U              | 5.1          | U              | 5.5          |     |
| n-Propylbenzene             | U              | 5.6          | 0.8            | J            | 5.1            | U            | 5.5 |
| 2-Chlorotoluene             | U              | 5.6          | U              | 5.1          | U              | 5.5          |     |
| 1,3,5-Tnmethylbenzene       | U              | 5.6          | 3              | J            | 5.1            | U            | 5.5 |
| 4-Chlorotoluene             | U              | 5.6          | U              | 5.1          | U              | 5.5          |     |
| t-Butylbenzene              | U              | 5.6          | U              | 5.1          | U              | 5.5          |     |
| 1,2,4-Tnmethylbenzene       | U              | 5.6          | 13             | 5.1          | U              | 5.5          |     |
| sec-Butylbenzene            | U              | 5.6          | U              | 5.1          | U              | 5.5          |     |
| 1,3-Dichlorobenzene         | U              | 5.6          | U              | 5.1          | U              | 5.5          |     |
| p-Isopropyltoluene          | U              | 5.6          | U              | 5.1          | U              | 5.5          |     |
| 1,4-Dichlorobenzene         | U              | 5.6          | U              | 5.1          | U              | 5.5          |     |
| n-Butylbenzene              | U              | 5.6          | 1.0            | J            | 5.1            | U            | 5.5 |
| 1,2-Dichlorobenzene         | U              | 5.6          | U              | 5.1          | U              | 5.5          |     |
| 1,2-Dibromo-3-chloropropane | U              | 11           | U              | 10           | U              | 11           |     |
| 1,2,4-Trichlorobenzene      | U              | 5.6          | U              | 5.1          | U              | 5.5          |     |
| Hexachlorobutadiene         | U              | 5.6          | U              | 5.1          | U              | 5.5          |     |
| Naphthalene                 | U              | 5.6          | 9.4            | 5.1          | U              | 5.5          |     |
| 1,2,3-Trichlorobenzene      | U              | 5.6          | U              | 5.1          | U              | 5.5          |     |

00016

Table 1.1 (Cont) Results of the Analysis for VOC in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 (Results are Based on Dry Weight)

| Sample No.<br>Sample Location<br>% Solid | Method Blank<br>4/27/99<br>NA |              | B16666<br>Location 16<br>86 |              | B16667<br>Location 17<br>92 |              | B16665<br>15<br>83 |              | B16668<br>Location 18<br>93 |              |   |    |     |    |     |
|--|-------------------------------|--------------|-----------------------------|--------------|-----------------------------|--------------|--------------------|--------------|-----------------------------|--------------|---|----|-----|----|-----|
| Compound Name                            | Conc.<br>µg/kg                | MDL<br>µg/kg | Conc.<br>µg/kg              | MDL<br>µg/kg | Conc.<br>µg/kg              | MDL<br>µg/kg | Conc.<br>µg/kg     | MDL<br>µg/kg | Conc.<br>µg/kg              | MDL<br>µg/kg |   |    |     |    |     |
| Dichlorodifluoromethane                  | U                             | 10           | U                           | 12           | U                           | 11           | U                  | 12           | U                           | 11           |   |    |     |    |     |
| Chloromethane                            | U                             | 10           | U                           | 12           | U                           | 11           | U                  | 12           | U                           | 11           |   |    |     |    |     |
| Vinyl chloride                           | U                             | 10           | U                           | 12           | U                           | 11           | U                  | 12           | U                           | 11           |   |    |     |    |     |
| Bromomethane                             | U                             | 10           | U                           | 12           | U                           | 11           | U                  | 12           | U                           | 11           |   |    |     |    |     |
| Chloroethane                             | U                             | 10           | U                           | 12           | U                           | 11           | U                  | 12           | U                           | 11           |   |    |     |    |     |
| Trichlorofluoromethane                   | U                             | 5            | U                           | 5.8          | U                           | 5.4          | U                  | 6            | U                           | 5.4          |   |    |     |    |     |
| Acrolein                                 | U                             | 50           | U                           | 58           | U                           | 54           | U                  | 60           | U                           | 54           |   |    |     |    |     |
| 1,1-Dichloroethylene                     | U                             | 5            | U                           | 5.8          | U                           | 5.4          | U                  | 6            | U                           | 5.4          |   |    |     |    |     |
| Trichlorotrifluoroethane                 | U                             | 5            | U                           | 5.8          | U                           | 5.4          | U                  | 6            | U                           | 5.4          |   |    |     |    |     |
| Acetone                                  | U                             | 20           | U                           | 23           | 16                          | J            | 22                 | U            | 24                          | 17           | J | 22 |     |    |     |
| Iodomethane                              | U                             | 5            | U                           | 5.8          | U                           | 5.4          | U                  | 6            | U                           | 5.4          |   |    |     |    |     |
| Carbon disulfide                         | U                             | 5            | U                           | 5.8          | U                           | 5.4          | U                  | 6            | U                           | 5.4          |   |    |     |    |     |
| Methylene chloride                       | 3.6                           | J            | 5                           | 7.4          | B                           | 5.8          | 4.7                | JB           | 5.4                         | 6.3          | B | 6  | 3.2 | JB | 5.4 |
| trans-1,2-Dichloroethylene               | U                             | 5            | U                           | 5.8          | U                           | 5.4          | U                  | 6            | U                           | 5.4          |   |    |     |    |     |
| Methyl t-butyl ether                     | U                             | 5            | U                           | 5.8          | U                           | 5.4          | U                  | 6            | U                           | 5.4          |   |    |     |    |     |
| Acrylonitrile                            | U                             | 50           | U                           | 58           | U                           | 54           | U                  | 60           | U                           | 54           |   |    |     |    |     |
| 1,1-Dichloroethane                       | U                             | 5            | U                           | 5.8          | U                           | 5.4          | U                  | 6            | U                           | 5.4          |   |    |     |    |     |
| Vinyl acetate                            | U                             | 20           | U                           | 23           | U                           | 22           | U                  | 24           | U                           | 22           |   |    |     |    |     |
| cis-1,2-Dichloroethylene                 | U                             | 5            | U                           | 5.8          | U                           | 5.4          | U                  | 6            | U                           | 5.4          |   |    |     |    |     |
| 2-Butanone                               | U                             | 20           | U                           | 23           | U                           | 22           | U                  | 24           | U                           | 22           |   |    |     |    |     |
| Bromochloromethane                       | U                             | 5            | U                           | 5.8          | U                           | 5.4          | U                  | 6            | U                           | 5.4          |   |    |     |    |     |
| Chloroform                               | U                             | 5            | U                           | 5.8          | U                           | 5.4          | U                  | 6            | U                           | 5.4          |   |    |     |    |     |
| 1,1,1-Trichloroethane                    | U                             | 5            | U                           | 5.8          | U                           | 5.4          | U                  | 6            | U                           | 5.4          |   |    |     |    |     |
| 2,2-Dichloropropane                      | U                             | 5            | U                           | 5.8          | U                           | 5.4          | U                  | 6            | U                           | 5.4          |   |    |     |    |     |
| Carbon tetrachloride                     | U                             | 5            | U                           | 5.8          | U                           | 5.4          | U                  | 6            | U                           | 5.4          |   |    |     |    |     |
| 1,1-Dichloropropene                      | U                             | 5            | U                           | 5.8          | U                           | 5.4          | U                  | 6            | U                           | 5.4          |   |    |     |    |     |
| 1,2-Dichloroethane                       | U                             | 5            | U                           | 5.8          | U                           | 5.4          | U                  | 6            | U                           | 5.4          |   |    |     |    |     |
| Benzene                                  | U                             | 5            | U                           | 5.8          | U                           | 5.4          | U                  | 6            | U                           | 5.4          |   |    |     |    |     |
| Trichloroethylene                        | U                             | 5            | U                           | 5.8          | U                           | 5.4          | U                  | 6            | U                           | 5.4          |   |    |     |    |     |
| 1,2-Dichloropropane                      | U                             | 5            | U                           | 5.8          | U                           | 5.4          | U                  | 6            | U                           | 5.4          |   |    |     |    |     |
| Dibromomethane                           | U                             | 5            | U                           | 5.8          | U                           | 5.4          | U                  | 6            | U                           | 5.4          |   |    |     |    |     |
| Bromochloromethane                       | U                             | 5            | U                           | 5.8          | U                           | 5.4          | U                  | 6            | U                           | 5.4          |   |    |     |    |     |
| 2-Chloroethyl vinyl ether                | U                             | 10           | U                           | 12           | U                           | 11           | U                  | 12           | U                           | 11           |   |    |     |    |     |
| cis-1,3-Dichloropropylene                | U                             | 5            | U                           | 5.8          | U                           | 5.4          | U                  | 6            | U                           | 5.4          |   |    |     |    |     |
| 4-Methyl-2-pentanone                     | U                             | 20           | U                           | 23           | U                           | 22           | U                  | 24           | U                           | 22           |   |    |     |    |     |
| Toluene                                  | U                             | 5            | U                           | 5.8          | U                           | 5.4          | U                  | 6            | U                           | 5.4          |   |    |     |    |     |

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Table 1.1 (Cont) Results of the Analysis for VOC in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 (Results are Based on Dry Weight)

| Sample No.<br>Sample Location<br>% Solid | Method Blank<br>4/27/99<br>NA | B16666<br>Location 16<br>86 | B16667<br>Location 17<br>92 | B16665<br>15<br>83 | B16668<br>Location 18<br>93 |              |                |              |                |              |
|--|-------------------------------|-----------------------------|-----------------------------|--------------------|-----------------------------|--------------|----------------|--------------|----------------|--------------|
| Compound Name                            | Conc.<br>µg/kg                | MDL<br>µg/kg                | Conc.<br>µg/kg              | MDL<br>µg/kg       | Conc.<br>µg/kg              | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg |
| trans-1,3-Dichloropropylene              | U                             | 5                           | U                           | 5.8                | U                           | 5.4          | U              | 6            | U              | 5.4          |
| 1,1,2-Trichloroethane                    | U                             | 5                           | U                           | 5.8                | U                           | 5.4          | U              | 6            | U              | 5.4          |
| 2-Hexanone                               | U                             | 20                          | U                           | 23                 | U                           | 22           | U              | 24           | U              | 22           |
| Tetrachloroethylene                      | U                             | 5                           | U                           | 5.8                | U                           | 5.4          | U              | 6            | U              | 5.4          |
| 1,3-Dichloropropane                      | U                             | 5                           | U                           | 5.8                | U                           | 5.4          | U              | 6            | U              | 5.4          |
| Dibromochloromethane                     | U                             | 5                           | U                           | 5.8                | U                           | 5.4          | U              | 6            | U              | 5.4          |
| 1,2-Dibromoethane                        | U                             | 5                           | U                           | 5.8                | U                           | 5.4          | U              | 6            | U              | 5.4          |
| 1-Chlorohexane                           | U                             | 5                           | U                           | 5.8                | U                           | 5.4          | U              | 6            | U              | 5.4          |
| Chlorobenzene                            | U                             | 5                           | U                           | 5.8                | U                           | 5.4          | U              | 6            | U              | 5.4          |
| 1,1,1,2-Tetrachloroethane                | U                             | 5                           | U                           | 5.8                | U                           | 5.4          | U              | 6            | U              | 5.4          |
| Ethyl benzene                            | U                             | 5                           | U                           | 5.8                | U                           | 5.4          | U              | 6            | U              | 5.4          |
| mp- Xylene                               | U                             | 5                           | U                           | 5.8                | U                           | 5.4          | U              | 6            | U              | 5.4          |
| o- Xylene                                | U                             | 5                           | U                           | 5.8                | U                           | 5.4          | U              | 6            | U              | 5.4          |
| Styrene                                  | U                             | 5                           | U                           | 5.8                | U                           | 5.4          | U              | 6            | U              | 5.4          |
| Bromoform                                | U                             | 5                           | U                           | 5.8                | U                           | 5.4          | U              | 6            | U              | 5.4          |
| Isopropylbenzene                         | U                             | 5                           | U                           | 5.8                | U                           | 5.4          | U              | 6            | U              | 5.4          |
| 1,2,3-Trichloropropane                   | U                             | 5                           | U                           | 5.8                | U                           | 5.4          | U              | 6            | U              | 5.4          |
| 1,1,1,2-Tetrachloroethane                | U                             | 5                           | U                           | 5.8                | U                           | 5.4          | U              | 6            | U              | 5.4          |
| Bromobenzene                             | U                             | 5                           | U                           | 5.8                | U                           | 5.4          | U              | 6            | U              | 5.4          |
| n-Propylbenzene                          | U                             | 5                           | U                           | 5.8                | U                           | 5.4          | U              | 6            | U              | 5.4          |
| 2-Chlorotoluene                          | U                             | 5                           | U                           | 5.8                | U                           | 5.4          | U              | 6            | U              | 5.4          |
| 1,3,5-Trimethylbenzene                   | U                             | 5                           | U                           | 5.8                | U                           | 5.4          | U              | 6            | U              | 5.4          |
| 4-Chlorotoluene                          | U                             | 5                           | U                           | 5.8                | U                           | 5.4          | U              | 6            | U              | 5.4          |
| t-Butylbenzene                           | U                             | 5                           | U                           | 5.8                | U                           | 5.4          | U              | 6            | U              | 5.4          |
| 1,2,4-Trimethylbenzene                   | U                             | 5                           | U                           | 5.8                | U                           | 5.4          | U              | 6            | U              | 5.4          |
| sec-Butylbenzene                         | U                             | 5                           | U                           | 5.8                | U                           | 5.4          | U              | 6            | U              | 5.4          |
| 1,3-Dichlorobenzene                      | U                             | 5                           | U                           | 5.8                | U                           | 5.4          | U              | 6            | U              | 5.4          |
| p-Isopropyltoluene                       | U                             | 5                           | U                           | 5.8                | U                           | 5.4          | U              | 6            | U              | 5.4          |
| 1,4-Dichlorobenzene                      | U                             | 5                           | U                           | 5.8                | U                           | 5.4          | U              | 6            | U              | 5.4          |
| n-Butylbenzene                           | U                             | 5                           | U                           | 5.8                | U                           | 5.4          | U              | 6            | U              | 5.4          |
| 1,2-Dichlorobenzene                      | U                             | 5                           | U                           | 5.8                | U                           | 5.4          | U              | 6            | U              | 5.4          |
| 1,2-Dibromo-3-chloropropane              | U                             | 10                          | U                           | 12                 | U                           | 11           | U              | 12           | U              | 11           |
| 1,2,4-Trichlorobenzene                   | U                             | 5                           | U                           | 5.8                | U                           | 5.4          | U              | 6            | U              | 5.4          |
| Hexachlorobutadiene                      | U                             | 5                           | U                           | 5.8                | U                           | 5.4          | U              | 6            | U              | 5.4          |
| Naphthalene                              | U                             | 5                           | U                           | 5.8                | U                           | 5.4          | U              | 6            | 0.9            | J 5.4        |
| 1,2,3-Trichlorobenzene                   | U                             | 5                           | U                           | 5.8                | U                           | 5.4          | U              | 6            | U              | 5.4          |

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Table 1.1 (Cont) Results of the Analysis for VOC in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 (Results are Based on Dry Weight)

| Sample No.<br>Sample Location<br>% Solid | B16669<br>Location 19<br>94 |              | B16670<br>Location 20<br>93 |              | B16671<br>Location 21<br>50 |              | B16675<br>Location 25<br>83 |              | B16676<br>Location 25-D<br>78 |              |
|--|-----------------------------|--------------|-----------------------------|--------------|-----------------------------|--------------|-----------------------------|--------------|-------------------------------|--------------|
|  | Conc.<br>µg/kg              | MDL<br>µg/kg | Conc.<br>µg/kg              | MDL<br>µg/kg | Conc.<br>µg/kg              | MDL<br>µg/kg | Conc.<br>µg/kg              | MDL<br>µg/kg | Conc.<br>µg/kg                | MDL<br>µg/kg |
| Dichlorodifluoromethane                  | U                           | 11           | U                           | 11           | U                           | 20           | U                           | 12           | U                             | 13           |
| Chloromethane                            | U                           | 11           | U                           | 11           | U                           | 20           | U                           | 12           | U                             | 13           |
| Vinyl chloride                           | U                           | 11           | U                           | 11           | U                           | 20           | U                           | 12           | U                             | 13           |
| Bromomethane                             | U                           | 11           | U                           | 11           | U                           | 20           | U                           | 12           | U                             | 13           |
| Chloroethane                             | U                           | 11           | U                           | 11           | U                           | 20           | U                           | 12           | U                             | 13           |
| Trichlorofluoromethane                   | U                           | 5.3          | U                           | 5.4          | U                           | 10           | U                           | 6.1          | U                             | 6.4          |
| Acrolein                                 | U                           | 53           | U                           | 54           | U                           | 100          | U                           | 61           | U                             | 64           |
| 1,1-Dichloroethylene                     | U                           | 5.3          | U                           | 5.4          | U                           | 10           | U                           | 6.1          | U                             | 6.4          |
| Trichlorotrifluoroethane                 | U                           | 5.3          | U                           | 5.4          | U                           | 10           | U                           | 6.1          | U                             | 6.4          |
| Acetone                                  | U                           | 21           | U                           | 22           | 37 J                        | 40           | U                           | 24           | 19 J                          | 26           |
| Iodomethane                              | U                           | 5.3          | U                           | 5.4          | U                           | 10           | U                           | 6.1          | U                             | 6.4          |
| Carbon disulfide                         | U                           | 5.3          | U                           | 5.4          | U                           | 10           | U                           | 6.1          | U                             | 6.4          |
| Methylene chloride                       | 6.6 B                       | 5.3          | 3.7 JB                      | 5.4          | 11 B                        | 10           | 18 B                        | 6.1          | 2.9 JB                        | 6.4          |
| trans-1,2-Dichloroethylene               | U                           | 5.3          | U                           | 5.4          | U                           | 10           | U                           | 6.1          | U                             | 6.4          |
| Methyl t-butyl ether                     | U                           | 5.3          | U                           | 5.4          | U                           | 10           | U                           | 6.1          | U                             | 6.4          |
| Acrylonitrile                            | U                           | 53           | U                           | 54           | U                           | 100          | U                           | 61           | U                             | 64           |
| 1,1-Dichloroethane                       | U                           | 5.3          | U                           | 5.4          | U                           | 10           | U                           | 6.1          | U                             | 6.4          |
| Vinyl acetate                            | U                           | 21           | U                           | 22           | U                           | 40           | U                           | 24           | U                             | 26           |
| cis-1,2-Dichloroethylene                 | U                           | 5.3          | U                           | 5.4          | U                           | 10           | U                           | 6.1          | U                             | 6.4          |
| 2-Butanone                               | U                           | 21           | U                           | 22           | U                           | 40           | U                           | 24           | U                             | 26           |
| Bromochloromethane                       | U                           | 5.3          | U                           | 5.4          | U                           | 10           | U                           | 6.1          | U                             | 6.4          |
| Chloroform                               | U                           | 5.3          | U                           | 5.4          | U                           | 10           | U                           | 6.1          | U                             | 6.4          |
| 1,1,1,-Trichloroethane                   | U                           | 5.3          | U                           | 5.4          | U                           | 10           | U                           | 6.1          | U                             | 6.4          |
| 2,2-Dichloropropane                      | U                           | 5.3          | U                           | 5.4          | U                           | 10           | U                           | 6.1          | U                             | 6.4          |
| Carbon tetrachloride                     | U                           | 5.3          | U                           | 5.4          | U                           | 10           | U                           | 6.1          | U                             | 6.4          |
| 1,1-Dichloropropene                      | U                           | 5.3          | U                           | 5.4          | U                           | 10           | U                           | 6.1          | U                             | 6.4          |
| 1,2-Dichloroethane                       | U                           | 5.3          | U                           | 5.4          | U                           | 10           | U                           | 6.1          | U                             | 6.4          |
| Benzene                                  | U                           | 5.3          | U                           | 5.4          | U                           | 10           | U                           | 6.1          | U                             | 6.4          |
| Trichloroethylene                        | U                           | 5.3          | U                           | 5.4          | U                           | 10           | U                           | 6.1          | U                             | 6.4          |
| 1,2-Dichloropropane                      | U                           | 5.3          | U                           | 5.4          | U                           | 10           | U                           | 6.1          | U                             | 6.4          |
| Dibromomethane                           | U                           | 5.3          | U                           | 5.4          | U                           | 10           | U                           | 6.1          | U                             | 6.4          |
| Bromochloromethane                       | U                           | 5.3          | U                           | 5.4          | U                           | 10           | U                           | 6.1          | U                             | 6.4          |
| 2-Chloroethyl vinyl ether                | U                           | 11           | U                           | 11           | U                           | 20           | U                           | 12           | U                             | 13           |
| cis-1,3-Dichloropropylene                | U                           | 5.3          | U                           | 5.4          | U                           | 10           | U                           | 6.1          | U                             | 6.4          |
| 4-Methyl-2-pentanone                     | U                           | 21           | U                           | 22           | U                           | 40           | U                           | 24           | U                             | 26           |
| Toluene                                  | U                           | 5.3          | U                           | 5.4          | U                           | 10           | U                           | 6.1          | U                             | 6.4          |



Table 1.1 (Cont) Results of the Analysis for VOC in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 (Results are Based on Dry Weight)

| Sample No.<br>Sample Location<br>% Solid | B16669         |              | B16670         |              | B16671         |              | B16675         |              | B16676         |              |
|--|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|
|  | Location 19    |              | Location 20    |              | Location 21    |              | Location 25    |              | Location 25-D  |              |
|  | 94             | 93           | 93             | 93           | 50             | 83           | 83             | 78           | 78             | 78           |
| Compound Name                            | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg |
| trans-1,3-Dichloropropylene              | U              | 5.3          | U              | 5.4          | U              | 10           | U              | 6.1          | U              | 6.4          |
| 1,1,2-Trichloroethane                    | U              | 5.3          | U              | 5.4          | U              | 10           | U              | 6.1          | U              | 6.4          |
| 2-Hexanone                               | U              | 21           | U              | 22           | U              | 40           | U              | 24           | U              | 26           |
| Tetrachloroethylene                      | U              | 5.3          | U              | 5.4          | U              | 10           | U              | 6.1          | U              | 6.4          |
| 1,3-Dichloropropane                      | U              | 5.3          | U              | 5.4          | U              | 10           | U              | 6.1          | U              | 6.4          |
| Dibromochloromethane                     | U              | 5.3          | U              | 5.4          | U              | 10           | U              | 6.1          | U              | 6.4          |
| 1,2-Dibromoethane                        | U              | 5.3          | U              | 5.4          | U              | 10           | U              | 6.1          | U              | 6.4          |
| 1-Chlorohexane                           | U              | 5.3          | U              | 5.4          | U              | 10           | U              | 6.1          | U              | 6.4          |
| Chlorobenzene                            | U              | 5.3          | U              | 5.4          | U              | 10           | U              | 6.1          | U              | 6.4          |
| 1,1,1,2-Tetrachloroethane                | U              | 5.3          | U              | 5.4          | U              | 10           | U              | 6.1          | U              | 6.4          |
| Ethyl benzene                            | U              | 5.3          | U              | 5.4          | U              | 10           | U              | 6.1          | U              | 6.4          |
| mp- Xylene                               | U              | 5.3          | U              | 5.4          | U              | 10           | U              | 6.1          | U              | 6.4          |
| o- Xylene                                | U              | 5.3          | U              | 5.4          | U              | 10           | U              | 6.1          | U              | 6.4          |
| Styrene                                  | U              | 5.3          | U              | 5.4          | U              | 10           | U              | 6.1          | U              | 6.4          |
| Bromoform                                | U              | 5.3          | U              | 5.4          | U              | 10           | U              | 6.1          | U              | 6.4          |
| Isopropylbenzene                         | U              | 5.3          | U              | 5.4          | U              | 10           | U              | 6.1          | U              | 6.4          |
| 1,2,3-Trichloropropane                   | U              | 5.3          | U              | 5.4          | U              | 10           | U              | 6.1          | U              | 6.4          |
| 1,1,1,2-Tetrachloroethane                | U              | 5.3          | U              | 5.4          | U              | 10           | U              | 6.1          | U              | 6.4          |
| Bromobenzene                             | U              | 5.3          | U              | 5.4          | U              | 10           | U              | 6.1          | U              | 6.4          |
| n-Propylbenzene                          | U              | 5.3          | U              | 5.4          | U              | 10           | U              | 6.1          | U              | 6.4          |
| 2-Chlorotoluene                          | U              | 5.3          | U              | 5.4          | U              | 10           | U              | 6.1          | U              | 6.4          |
| 1,3,5-Trimethylbenzene                   | U              | 5.3          | U              | 5.4          | U              | 10           | U              | 6.1          | U              | 6.4          |
| 4-Chlorotoluene                          | U              | 5.3          | U              | 5.4          | U              | 10           | U              | 6.1          | U              | 6.4          |
| t-Butylbenzene                           | U              | 5.3          | U              | 5.4          | U              | 10           | U              | 6.1          | U              | 6.4          |
| 1,2,4-Trimethylbenzene                   | U              | 5.3          | U              | 5.4          | U              | 10           | U              | 6.1          | U              | 6.4          |
| sec-Butylbenzene                         | U              | 5.3          | U              | 5.4          | U              | 10           | U              | 6.1          | U              | 6.4          |
| 1,3-Dichlorobenzene                      | U              | 5.3          | U              | 5.4          | U              | 10           | U              | 6.1          | U              | 6.4          |
| p-Isopropyltoluene                       | U              | 5.3          | U              | 5.4          | 100            | 10           | U              | 6.1          | U              | 6.4          |
| 1,4-Dichlorobenzene                      | U              | 5.3          | U              | 5.4          | U              | 10           | U              | 6.1          | U              | 6.4          |
| n-Butylbenzene                           | U              | 5.3          | U              | 5.4          | U              | 10           | U              | 6.1          | U              | 6.4          |
| 1,2-Dichlorobenzene                      | U              | 5.3          | U              | 5.4          | U              | 10           | U              | 6.1          | U              | 6.4          |
| 1,2-Dibromo-3-chloropropane              | U              | 11           | U              | 11           | U              | 20           | U              | 12           | U              | 13           |
| 1,2,4-Trichlorobenzene                   | U              | 5.3          | U              | 5.4          | U              | 10           | U              | 6.1          | U              | 6.4          |
| Hexachlorobutadiene                      | U              | 5.3          | U              | 5.4          | U              | 10           | U              | 6.1          | U              | 6.4          |
| Naphthalene                              | U              | 5.3          | 1.1            | 5.4          | 3.2            | 10           | 6.4            | 6.1          | U              | 6.4          |
| 1,2,3-Trichlorobenzene                   | U              | 5.3          | U              | 5.4          | U              | 10           | U              | 6.1          | U              | 6.4          |

00020

Table 1.1 (Cont) Results of the Analysis for VOC in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 (Results are Based on Dry Weight)

| Sample No.                 | B16677         |              | B16679         |              |
|----------------------------|----------------|--------------|----------------|--------------|
| Sample Location            | Location 26    |              | Location 27    |              |
| % Solid                    | 72             |              | 86             |              |
| Compound Name              | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg |
| Dichlorodifluoromethane    | U              | 14           | U              | 12           |
| Chloromethane              | U              | 14           | U              | 12           |
| Vinyl chloride             | U              | 14           | U              | 12           |
| Bromomethane               | U              | 14           | U              | 12           |
| Chloroethane               | U              | 14           | U              | 12           |
| Trichlorofluoromethane     | U              | 7            | U              | 5.8          |
| Acrolein                   | U              | 70           | U              | 58           |
| 1,1-Dichloroethylene       | U              | 7            | U              | 5.8          |
| Trichlorotrifluoroethane   | U              | 7            | U              | 5.8          |
| Acetone                    | U              | 28           | 20 J           | 23           |
| Iodomethane                | U              | 7            | U              | 5.8          |
| Carbon disulfide           | U              | 7            | U              | 5.8          |
| Methylene chloride         | 6.4 JB         | 7            | 3.1 JB         | 5.8          |
| trans-1,2-Dichloroethylene | U              | 7            | U              | 5.8          |
| Methyl t-butyl ether       | U              | 7            | U              | 5.8          |
| Acrylonitrile              | U              | 70           | U              | 58           |
| 1,1-Dichloroethane         | U              | 7            | U              | 5.8          |
| Vinyl acetate              | U              | 28           | U              | 23           |
| cis-1,2-Dichloroethylene   | U              | 7            | U              | 5.8          |
| 2-Butanone                 | U              | 28           | U              | 23           |
| Bromochloromethane         | U              | 7            | U              | 5.8          |
| Chloroform                 | U              | 7            | U              | 5.8          |
| 1,1,1,-Trichloroethane     | U              | 7            | U              | 5.8          |
| 2,2-Dichloropropane        | U              | 7            | U              | 5.8          |
| Carbon tetrachloride       | U              | 7            | U              | 5.8          |
| 1,1-Dichloropropene        | U              | 7            | U              | 5.8          |
| 1,2-Dichloroethane         | U              | 7            | U              | 5.8          |
| Benzene                    | U              | 7            | U              | 5.8          |
| Trichloroethylene          | U              | 7            | U              | 5.8          |
| 1,2-Dichloropropane        | U              | 7            | U              | 5.8          |
| Dibromomethane             | U              | 7            | U              | 5.8          |
| Bromochloromethane         | U              | 7            | U              | 5.8          |
| 2-Chloroethyl vinyl ether  | U              | 14           | U              | 12           |
| cis-1,3-Dichloropropylene  | U              | 7            | U              | 5.8          |
| 4-Methyl-2-pentanone       | U              | 28           | U              | 23           |
| Toluene                    | U              | 7            | U              | 5.8          |

00021

Table 1.1 (Cont) Results of the Analysis for VOC in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 (Results are Based on Dry Weight)

| Sample No.                  | B16677         |              | B16679         |              |
|-----------------------------|----------------|--------------|----------------|--------------|
| Sample Location             | Location 26    |              | Location 27    |              |
| % Solid                     | 72             |              | 86             |              |
| Compound Name               | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg |
| trans-1,3-Dichloropropylene | U              | 7            | U              | 5.8          |
| 1,1,2-Trichloroethane       | U              | 7            | U              | 5.8          |
| 2-Hexanone                  | U              | 28           | U              | 23           |
| Tetrachloroethylene         | U              | 7            | U              | 5.8          |
| 1,3-Dichloropropane         | U              | 7            | U              | 5.8          |
| Dibromochloromethane        | U              | 7            | U              | 5.8          |
| 1,2-Dibromoethane           | U              | 7            | U              | 5.8          |
| 1-Chlorohexane              | U              | 7            | U              | 5.8          |
| Chlorobenzene               | U              | 7            | U              | 5.8          |
| 1,1,1,2-Tetrachloroethane   | U              | 7            | U              | 5.8          |
| Ethyl benzene               | U              | 7            | U              | 5.8          |
| mp- Xylene                  | U              | 7            | U              | 5.8          |
| o- Xylene                   | U              | 7            | U              | 5.8          |
| Styrene                     | U              | 7            | U              | 5.8          |
| Bromoform                   | U              | 7            | U              | 5.8          |
| Isopropylbenzene            | U              | 7            | U              | 5.8          |
| 1,2,3-Trichloropropane      | U              | 7            | U              | 5.8          |
| 1,1,2,2-Tetrachloroethane   | U              | 7            | U              | 5.8          |
| Bromobenzene                | U              | 7            | U              | 5.8          |
| n-Propylbenzene             | U              | 7            | U              | 5.8          |
| 2-Chlorotoluene             | U              | 7            | U              | 5.8          |
| 1,3,5-Trimethylbenzene      | U              | 7            | U              | 5.8          |
| 4-Chlorotoluene             | U              | 7            | U              | 5.8          |
| t-Butylbenzene              | U              | 7            | U              | 5.8          |
| 1,2,4-Trimethylbenzene      | U              | 7            | U              | 5.8          |
| sec-Butylbenzene            | U              | 7            | U              | 5.8          |
| 1,3-Dichlorobenzene         | U              | 7            | U              | 5.8          |
| p-Isopropyltoluene          | U              | 7            | U              | 5.8          |
| 1,4-Dichlorobenzene         | U              | 7            | U              | 5.8          |
| n-Butylbenzene              | U              | 7            | U              | 5.8          |
| 1,2-Dichlorobenzene         | U              | 7            | U              | 5.8          |
| 1,2-Dibromo-3-chloropropane | U              | 14           | U              | 12           |
| 1,2,4-Trichlorobenzene      | U              | 7            | U              | 5.8          |
| Hexachlorobutadiene         | U              | 7            | U              | 5.8          |
| Naphthalene                 | U              | 7            | U              | 5.8          |
| 1,2,3-Trichlorobenzene      | U              | 7            | U              | 5.8          |

00022

Table 1.1 (Cont) Results of the Analysis for VOC in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 (Results are Based on Dry Weight)

| Sample No.<br>Sample Location<br>% Solid | B16680<br>Field Blank (VOC)<br>100 |              | B16678<br>Location 26D<br>55 |              | B16672<br>Location 22<br>80 |              | B16673<br>Location 23<br>84 |              |     |     |    |   |
|--|------------------------------------|--------------|------------------------------|--------------|-----------------------------|--------------|-----------------------------|--------------|-----|-----|----|---|
|  | Conc.<br>µg/kg                     | MDL<br>µg/kg | Conc.<br>µg/kg               | MDL<br>µg/kg | Conc.<br>µg/kg              | MDL<br>µg/kg | Conc.<br>µg/kg              | MDL<br>µg/kg |     |     |    |   |
| Dichlorodifluoromethane                  | U                                  | 10           | U                            | 18           | U                           | 12           | U                           | 12           |     |     |    |   |
| Chloromethane                            | U                                  | 10           | U                            | 18           | U                           | 12           | U                           | 12           |     |     |    |   |
| Vinyl chloride                           | U                                  | 10           | U                            | 18           | U                           | 12           | U                           | 12           |     |     |    |   |
| Bromomethane                             | U                                  | 10           | U                            | 18           | U                           | 12           | U                           | 12           |     |     |    |   |
| Chloroethane                             | U                                  | 10           | U                            | 18           | U                           | 12           | U                           | 12           |     |     |    |   |
| Trichlorofluoromethane                   | U                                  | 5            | U                            | 9            | U                           | 6.2          | U                           | 6            |     |     |    |   |
| Acrolein                                 | U                                  | 50           | U                            | 90           | U                           | 62           | U                           | 60           |     |     |    |   |
| 1,1-Dichloroethylene                     | U                                  | 5            | U                            | 9            | U                           | 6.2          | U                           | 6            |     |     |    |   |
| Trichlorotrifluoroethane                 | U                                  | 5            | U                            | 9            | U                           | 6.2          | U                           | 6            |     |     |    |   |
| Acetone                                  | U                                  | 20           | U                            | 36           | U                           | 25           | U                           | 24           |     |     |    |   |
| Iodomethane                              | U                                  | 5            | U                            | 9            | U                           | 6.2          | U                           | 6            |     |     |    |   |
| Carbon disulfide                         | U                                  | 5            | U                            | 9            | U                           | 6.2          | U                           | 6            |     |     |    |   |
| Methylene chloride                       | 2.4                                | JB           | 5                            | 4.1          | JB                          | 9            | 3.9                         | JB           | 6.2 | 3.6 | JB | 6 |
| trans-1,2-Dichloroethylene               | U                                  | 5            | U                            | 9            | U                           | 6.2          | U                           | 6            |     |     |    |   |
| Methyl t-butyl ether                     | U                                  | 5            | U                            | 9            | U                           | 6.2          | U                           | 6            |     |     |    |   |
| Acrylonitrile                            | U                                  | 50           | U                            | 90           | U                           | 62           | U                           | 60           |     |     |    |   |
| 1,1-Dichloroethane                       | U                                  | 5            | U                            | 9            | U                           | 6.2          | U                           | 6            |     |     |    |   |
| Vinyl acetate                            | U                                  | 20           | U                            | 36           | U                           | 25           | U                           | 24           |     |     |    |   |
| cis-1,2-Dichloroethylene                 | U                                  | 5            | U                            | 9            | U                           | 6.2          | U                           | 6            |     |     |    |   |
| 2-Butanone                               | U                                  | 20           | U                            | 36           | U                           | 25           | U                           | 24           |     |     |    |   |
| Bromochloromethane                       | U                                  | 5            | U                            | 9            | U                           | 6.2          | U                           | 6            |     |     |    |   |
| Chloroform                               | U                                  | 5            | U                            | 9            | U                           | 6.2          | U                           | 6            |     |     |    |   |
| 1,1,1,-Trichloroethane                   | U                                  | 5            | U                            | 9            | U                           | 6.2          | U                           | 6            |     |     |    |   |
| 2,2-Dichloropropane                      | U                                  | 5            | U                            | 9            | U                           | 6.2          | U                           | 6            |     |     |    |   |
| Carbon tetrachloride                     | U                                  | 5            | U                            | 9            | U                           | 6.2          | U                           | 6            |     |     |    |   |
| 1,1-Dichloropropene                      | U                                  | 5            | U                            | 9            | U                           | 6.2          | U                           | 6            |     |     |    |   |
| 1,2-Dichloroethane                       | U                                  | 5            | U                            | 9            | U                           | 6.2          | U                           | 6            |     |     |    |   |
| Benzene                                  | U                                  | 5            | U                            | 9            | U                           | 6.2          | U                           | 6            |     |     |    |   |
| Trichloroethylene                        | U                                  | 5            | U                            | 9            | U                           | 6.2          | U                           | 6            |     |     |    |   |
| 1,2-Dichloropropane                      | U                                  | 5            | U                            | 9            | U                           | 6.2          | U                           | 6            |     |     |    |   |
| Dibromomethane                           | U                                  | 5            | U                            | 9            | U                           | 6.2          | U                           | 6            |     |     |    |   |
| Bromochloromethane                       | U                                  | 5            | U                            | 9            | U                           | 6.2          | U                           | 6            |     |     |    |   |
| 2-Chloroethyl vinyl ether                | U                                  | 10           | U                            | 18           | U                           | 12           | U                           | 12           |     |     |    |   |
| cis-1,3-Dichloropropylene                | U                                  | 5            | U                            | 9            | U                           | 6.2          | U                           | 6            |     |     |    |   |
| 4-Methyl-2-pentanone                     | U                                  | 20           | U                            | 36           | U                           | 25           | U                           | 24           |     |     |    |   |
| Toluene                                  | U                                  | 5            | U                            | 9            | U                           | 6.2          | U                           | 6            |     |     |    |   |

Table 1.1 (Cont) Results of the Analysis for VOC in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 (Results are Based on Dry Weight)

| Sample No.                  | B16680            |              | B16678         |              | B16672         |              | B16673         |              |
|-----------------------------|-------------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|
| Sample Location             | Field Blank (VOC) |              | Location 26D   |              | Location 22    |              | Location 23    |              |
| % Solid                     | 100               |              | 55             |              | 80             |              | 84             |              |
| Compound Name               | Conc.<br>µg/kg    | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg |
| trans-1,3-Dichloropropylene | U                 | 5            | U              | 9            | U              | 6.2          | U              | 6            |
| 1,1,2-Trichloroethane       | U                 | 5            | U              | 9            | U              | 6.2          | U              | 6            |
| 2-Hexanone                  | U                 | 20           | U              | 36           | U              | 25           | U              | 24           |
| Tetrachloroethylene         | U                 | 5            | U              | 9            | U              | 6.2          | U              | 6            |
| 1,3-Dichloropropane         | U                 | 5            | U              | 9            | U              | 6.2          | U              | 6            |
| Dibromochloromethane        | U                 | 5            | U              | 9            | U              | 6.2          | U              | 6            |
| 1,2-Dibromoethane           | U                 | 5            | U              | 9            | U              | 6.2          | U              | 6            |
| 1-Chlorohexane              | U                 | 5            | U              | 9            | U              | 6.2          | U              | 6            |
| Chlorobenzene               | U                 | 5            | U              | 9            | U              | 6.2          | U              | 6            |
| 1,1,1,2-Tetrachloroethane   | U                 | 5            | U              | 9            | U              | 6.2          | U              | 6            |
| Ethyl benzene               | U                 | 5            | U              | 9            | U              | 6.2          | U              | 6            |
| mp- Xylene                  | U                 | 5            | U              | 9            | U              | 6.2          | U              | 6            |
| o- Xylene                   | U                 | 5            | U              | 9            | U              | 6.2          | U              | 6            |
| Styrene                     | U                 | 5            | U              | 9            | U              | 6.2          | U              | 6            |
| Bromoform                   | U                 | 5            | U              | 9            | U              | 6.2          | U              | 6            |
| Isopropylbenzene            | U                 | 5            | U              | 9            | U              | 6.2          | U              | 6            |
| 1,2,3-Trichloropropane      | U                 | 5            | U              | 9            | U              | 6.2          | U              | 6            |
| 1,1,2,2-Tetrachloroethane   | U                 | 5            | U              | 9            | U              | 6.2          | U              | 6            |
| Bromobenzene                | U                 | 5            | U              | 9            | U              | 6.2          | U              | 6            |
| n-Propylbenzene             | U                 | 5            | U              | 9            | U              | 6.2          | U              | 6            |
| 2-Chlorotoluene             | U                 | 5            | U              | 9            | U              | 6.2          | U              | 6            |
| 1,3,5-Trimethylbenzene      | U                 | 5            | U              | 9            | U              | 6.2          | U              | 6            |
| 4-Chlorotoluene             | U                 | 5            | U              | 9            | U              | 6.2          | U              | 6            |
| t-Butylbenzene              | U                 | 5            | U              | 9            | U              | 6.2          | U              | 6            |
| 1,2,4-Trimethylbenzene      | U                 | 5            | U              | 9            | U              | 6.2          | U              | 6            |
| sec-Butylbenzene            | U                 | 5            | U              | 9            | U              | 6.2          | U              | 6            |
| 1,3-Dichlorobenzene         | U                 | 5            | U              | 9            | U              | 6.2          | U              | 6            |
| p-Isopropyltoluene          | U                 | 5            | U              | 9            | U              | 6.2          | U              | 6            |
| 1,4-Dichlorobenzene         | U                 | 5            | U              | 9            | U              | 6.2          | U              | 6            |
| n-Butylbenzene              | U                 | 5            | U              | 9            | U              | 6.2          | U              | 6            |
| 1,2-Dichlorobenzene         | U                 | 5            | U              | 9            | U              | 6.2          | U              | 6            |
| 1,2-Dibromo-3-chloropropane | U                 | 10           | U              | 18           | U              | 12           | U              | 12           |
| 1,2,4-Trichlorobenzene      | U                 | 5            | U              | 9            | U              | 6.2          | U              | 6            |
| Hexachlorobutadiene         | U                 | 5            | U              | 9            | U              | 6.2          | U              | 6            |
| Naphthalene                 | U                 | 5            | U              | 9            | U              | 6.2          | U              | 6            |
| 1,2,3-Trichlorobenzene      | U                 | 5            | U              | 9            | U              | 6.2          | U              | 6            |

00024

Table 1.2 Results of the Analysis for Pesticide/PCBs in Soil  
 WA# 3-439 Lackawanna Foundry Site  
 Based on Dry Weight

| Client ID          | SBLK041599     |              | A16650         |              | A16651           |              | A16652         |              | A16653         |              |
|--------------------|----------------|--------------|----------------|--------------|------------------|--------------|----------------|--------------|----------------|--------------|
| Location           | -              |              | 1 West Fence   |              | 2 Surface Middle |              | 2D Middle      |              | 3 West Fence   |              |
| Percent Solid      | 100            |              | 57             |              | 94.7             |              | 90.5           |              | 73.2           |              |
| Analyte            | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg   | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg |
| a-BHC              | U              | 3.3          | U              | 5.8          | U                | 3.5          | U              | 3.7          | U              | 4.6          |
| g-BHC              | U              | 3.3          | U              | 5.8          | U                | 3.5          | U              | 3.7          | U              | 4.6          |
| b-BHC              | U              | 3.3          | U              | 5.8          | U                | 3.5          | U              | 3.7          | U              | 4.6          |
| Heptachlor         | U              | 3.3          | U              | 5.8          | U                | 3.5          | U              | 3.7          | U              | 4.6          |
| d-BHC              | U              | 3.3          | U              | 5.8          | U                | 3.5          | U              | 3.7          | U              | 4.6          |
| Aldnn              | U              | 3.3          | U              | 5.8          | U                | 3.5          | U              | 3.7          | U              | 4.6          |
| Heptachlor Epoxide | U              | 3.3          | U              | 5.8          | U                | 3.5          | U              | 3.7          | U              | 4.6          |
| g-Chlordane        | U              | 3.3          | U              | 5.8          | U                | 3.5          | U              | 3.7          | U              | 4.6          |
| a-Chlordane        | U              | 3.3          | U              | 5.8          | U                | 3.5          | U              | 3.7          | U              | 4.6          |
| Endosulfan (I)     | U              | 3.3          | U              | 5.8          | U                | 3.5          | U              | 3.7          | U              | 4.6          |
| p,p'-D D E         | U              | 3.3          | 12             | 5.8          | U                | 3.5          | U              | 3.7          | 30             | 4.6          |
| Dieldnn            | U              | 3.3          | U              | 5.8          | U                | 3.5          | U              | 3.7          | U              | 4.6          |
| Endnn              | U              | 3.3          | U              | 5.8          | U                | 3.5          | U              | 3.7          | U              | 4.6          |
| p,p'-D D D         | U              | 3.3          | U              | 5.8          | U                | 3.5          | U              | 3.7          | U              | 4.6          |
| Endosulfan (II)    | U              | 3.3          | U              | 5.8          | U                | 3.5          | U              | 3.7          | U              | 4.6          |
| p,p'-D D T         | U              | 3.3          | 25             | 5.8          | U                | 3.5          | U              | 3.7          | 14             | 4.6          |
| Endrin Aldehyde    | U              | 3.3          | U              | 5.8          | U                | 3.5          | U              | 3.7          | U              | 4.6          |
| Endosulfan Sulfate | U              | 3.3          | U              | 5.8          | U                | 3.5          | U              | 3.7          | U              | 4.6          |
| Methoxychlor       | U              | 3.3          | U              | 5.8          | U                | 3.5          | U              | 3.7          | U              | 4.6          |
| Endnn Ketone       | U              | 3.3          | U              | 5.8          | U                | 3.5          | U              | 3.7          | U              | 4.6          |
| Toxaphene          | U              | 83           | U              | 150          | U                | 88           | U              | 92           | U              | 110          |
| Aroclor 1016       | U              | 42           | U              | 73           | U                | 44           | U              | 46           | U              | 57           |
| Aroclor 1221       | U              | 83           | U              | 150          | U                | 88           | U              | 92           | U              | 110          |
| Aroclor 1232       | U              | 42           | U              | 73           | U                | 44           | U              | 46           | U              | 57           |
| Aroclor 1242       | U              | 42           | U              | 73           | U                | 44           | U              | 46           | U              | 57           |
| Aroclor 1248       | U              | 42           | U              | 73           | 560              | 44           | 2700           | 46           | U              | 57           |
| Aroclor 1254       | U              | 42           | U              | 73           | U                | 44           | U              | 46           | 25 J           | 57           |
| Aroclor 1260       | U              | 42           | U              | 73           | U                | 44           | U              | 46           | 19 J           | 57           |
| Aroclor 1268       | U              | 42           | U              | 73           | U                | 44           | U              | 46           | U              | 57           |

Table 1.2 (Cont) Results of the Analysis for Pesticide/PCBs in Soil  
 WA# 3-439 Lackawanna Foundry Site  
 Based on Dry Weight

| Client ID<br>Location<br>Percent Solid | A16654<br>4 West Fence<br>82.1 |              | A16655<br>5 West Fence<br>80.7 |              | A16656<br>6 Near Entrance<br>85 |              |
|--|--------------------------------|--------------|--------------------------------|--------------|---------------------------------|--------------|
|  | Conc.<br>µg/kg                 | MDL<br>µg/kg | Conc.<br>µg/kg                 | MDL<br>µg/kg | Conc.<br>µg/kg                  | MDL<br>µg/kg |
| a-BHC                                  | U                              | 4.1          | U                              | 4.1          | U                               | 3.9          |
| g-BHC                                  | U                              | 4.1          | U                              | 4.1          | U                               | 3.9          |
| b-BHC                                  | U                              | 4.1          | U                              | 4.1          | U                               | 3.9          |
| Heptachlor                             | U                              | 4.1          | U                              | 4.1          | U                               | 3.9          |
| d-BHC                                  | U                              | 4.1          | U                              | 4.1          | U                               | 3.9          |
| Aldnn                                  | U                              | 4.1          | U                              | 4.1          | U                               | 3.9          |
| Heptachlor Epoxide                     | U                              | 4.1          | U                              | 4.1          | U                               | 3.9          |
| g-Chlordane                            | U                              | 4.1          | U                              | 4.1          | U                               | 3.9          |
| a-Chlordane                            | U                              | 4.1          | U                              | 4.1          | U                               | 3.9          |
| Endosulfan (I)                         | U                              | 4.1          | U                              | 4.1          | U                               | 3.9          |
| p,p'-D D E                             | 19                             | 4.1          | 0.9 J                          | 4.1          | 12                              | 3.9          |
| Dieldnn                                | U                              | 4.1          | U                              | 4.1          | U                               | 3.9          |
| Endnn                                  | U                              | 4.1          | U                              | 4.1          | U                               | 3.9          |
| p,p'-D D D                             | 6.9                            | 4.1          | U                              | 4.1          | U                               | 3.9          |
| Endosulfan (II)                        | U                              | 4.1          | U                              | 4.1          | U                               | 3.9          |
| p,p'-D D T                             | 6.4                            | 4.1          | U                              | 4.1          | U                               | 3.9          |
| Endrin Aldehyde                        | U                              | 4.1          | U                              | 4.1          | U                               | 3.9          |
| Endosulfan Sulfate                     | U                              | 4.1          | U                              | 4.1          | U                               | 3.9          |
| Methoxychlor                           | U                              | 4.1          | U                              | 4.1          | U                               | 3.9          |
| Endrin Ketone                          | U                              | 4.1          | U                              | 4.1          | U                               | 3.9          |
| Toxaphene                              | U                              | 100          | U                              | 100          | U                               | 98           |
| Aroclor 1016                           | U                              | 51           | U                              | 52           | U                               | 49           |
| Aroclor 1221                           | U                              | 100          | U                              | 100          | U                               | 98           |
| Aroclor 1232                           | U                              | 51           | U                              | 52           | U                               | 49           |
| Aroclor 1242                           | U                              | 51           | U                              | 52           | U                               | 49           |
| Aroclor 1248                           | 120                            | 51           | 15 J                           | 52           | 770                             | 49           |
| Aroclor 1254                           | 81                             | 51           | 16 JW                          | 52           | 1100                            | 49           |
| Aroclor 1260                           | 31 J                           | 51           | 13 J                           | 52           | 680                             | 49           |
| Aroclor 1268                           | U                              | 51           | U                              | 52           | U                               | 49           |

Table 1.2 (Cont) Results of the Analysis for Pesticide/PCBs in Soil  
 WA# 3-439 Lackawanna Foundry Site  
 Based on Dry Weight

| Client ID          | SBLK041999     |              | A16657          |              | A16658          |              | A16659          |              | A16660           |              |
|--------------------|----------------|--------------|-----------------|--------------|-----------------|--------------|-----------------|--------------|------------------|--------------|
| Location           | -              |              | 7 Near Entrance |              | 8 Near Entrance |              | 9 Near Entrance |              | 10 Near Entrance |              |
| Percent Solid      | 100            |              | 76.4            |              | 86.8            |              | 91.2            |              | 98.9             |              |
| Analyte            | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg  | MDL<br>µg/kg | Conc.<br>µg/kg  | MDL<br>µg/kg | Conc.<br>µg/kg  | MDL<br>µg/kg | Conc.<br>µg/kg   | MDL<br>µg/kg |
| a-BHC              | U              | 3.3          | U               | 4.4          | U               | 3.8          | U               | 3.7          | U                | 3.4          |
| g-BHC              | U              | 3.3          | U               | 4.4          | U               | 3.8          | U               | 3.7          | U                | 3.4          |
| b-BHC              | U              | 3.3          | U               | 4.4          | U               | 3.8          | U               | 3.7          | U                | 3.4          |
| Heptachlor         | U              | 3.3          | U               | 4.4          | U               | 3.8          | U               | 3.7          | U                | 3.4          |
| d-BHC              | U              | 3.3          | U               | 4.4          | U               | 3.8          | U               | 3.7          | U                | 3.4          |
| Aldrin             | U              | 3.3          | U               | 4.4          | U               | 3.8          | U               | 3.7          | U                | 3.4          |
| Heptachlor Epoxide | U              | 3.3          | U               | 4.4          | U               | 3.8          | U               | 3.7          | U                | 3.4          |
| g-Chlordane        | U              | 3.3          | U               | 4.4          | U               | 3.8          | U               | 3.7          | U                | 3.4          |
| a-Chlordane        | U              | 3.3          | U               | 4.4          | U               | 3.8          | U               | 3.7          | U                | 3.4          |
| Endosulfan (I)     | U              | 3.3          | U               | 4.4          | U               | 3.8          | U               | 3.7          | U                | 3.4          |
| p,p'-D D E         | U              | 3.3          | 4.5             | 4.4          | 2.5 J           | 3.8          | 17              | 3.7          | 4.7              | 3.4          |
| Dieldrin           | U              | 3.3          | 2.9 J           | 4.4          | 2.0 J           | 3.8          | U               | 3.7          | 2.7 J            | 3.4          |
| Endrin             | U              | 3.3          | U               | 4.4          | U               | 3.8          | U               | 3.7          | U                | 3.4          |
| p,p'-D D D         | U              | 3.3          | U               | 4.4          | U               | 3.8          | U               | 3.7          | U                | 3.4          |
| Endosulfan (II)    | U              | 3.3          | U               | 4.4          | U               | 3.8          | U               | 3.7          | U                | 3.4          |
| p,p'-D D T         | U              | 3.3          | 9.6             | 4.4          | 5.3             | 3.8          | U               | 3.7          | U                | 3.4          |
| Endrin Aldehyde    | U              | 3.3          | U               | 4.4          | U               | 3.8          | U               | 3.7          | U                | 3.4          |
| Endosulfan Sulfate | U              | 3.3          | U               | 4.4          | U               | 3.8          | U               | 3.7          | U                | 3.4          |
| Methoxychlor       | U              | 3.3          | U               | 4.4          | U               | 3.8          | U               | 3.7          | U                | 3.4          |
| Endrin Ketone      | U              | 3.3          | U               | 4.4          | U               | 3.8          | U               | 3.7          | U                | 3.4          |
| Toxaphene          | U              | 83           | U               | 110          | U               | 96           | U               | 91           | U                | 84           |
| Aroclor 1016       | U              | 42           | U               | 55           | U               | 48           | U               | 46           | U                | 42           |
| Aroclor 1221       | U              | 83           | U               | 110          | U               | 96           | U               | 91           | U                | 84           |
| Aroclor 1232       | U              | 42           | U               | 55           | U               | 48           | U               | 46           | U                | 42           |
| Aroclor 1242       | U              | 42           | U               | 55           | U               | 48           | U               | 46           | 6500             | 42           |
| Aroclor 1248       | U              | 42           | 590             | 55           | 68              | 48           | 3500            | 46           | U                | 42           |
| Aroclor 1254       | U              | 42           | 200             | 55           | 68              | 48           | 1300            | 46           | U                | 42           |
| Aroclor 1260       | U              | 42           | 67              | 55           | 32 J            | 48           | 320             | 46           | U                | 42           |
| Aroclor 1268       | U              | 42           | U               | 55           | U               | 48           | U               | 46           | U                | 42           |



Table 1.2 (Cont) Results of the Analysis for Pesticide/PCBs in Soil  
 WA# 3-439 Lackawanna Foundry Site  
 Based on Dry Weight

| Client ID          | A16661           |              | A16662           |              | A16663           |              | A16664           |              | A16665           |              |
|--------------------|------------------|--------------|------------------|--------------|------------------|--------------|------------------|--------------|------------------|--------------|
| Location           | 11 Near Entrance |              | 12 Near Entrance |              | 13 Near Entrance |              | 14 Near Entrance |              | 15 Near Entrance |              |
| Percent Solid      | 84.5             |              | 79.6             |              | 91               |              | 73.5             |              | 62.4             |              |
| Analyte            | Conc.<br>µg/kg   | MDL<br>µg/kg | Conc.<br>µg/kg   | MDL<br>µg/kg | Conc.<br>µg/kg   | MDL<br>µg/kg | Conc.<br>µg/kg   | MDL<br>µg/kg | Conc.<br>µg/kg   | MDL<br>µg/kg |
| a-BHC              | U                | 3.9          | U                | 4.2          | U                | 3.7          | U                | 4.5          | U                | 5.3          |
| g-BHC              | U                | 3.9          | U                | 4.2          | U                | 3.7          | U                | 4.5          | U                | 5.3          |
| b-BHC              | U                | 3.9          | U                | 4.2          | U                | 3.7          | U                | 4.5          | U                | 5.3          |
| Heptachlor         | U                | 3.9          | U                | 4.2          | U                | 3.7          | U                | 4.5          | U                | 5.3          |
| d-BHC              | U                | 3.9          | U                | 4.2          | U                | 3.7          | U                | 4.5          | U                | 5.3          |
| Aldrin             | U                | 3.9          | U                | 4.2          | U                | 3.7          | U                | 4.5          | U                | 5.3          |
| Heptachlor Epoxide | U                | 3.9          | U                | 4.2          | U                | 3.7          | U                | 4.5          | U                | 5.3          |
| g-Chlordane        | U                | 3.9          | U                | 4.2          | U                | 3.7          | U                | 4.5          | U                | 5.3          |
| a-Chlordane        | U                | 3.9          | U                | 4.2          | U                | 3.7          | U                | 4.5          | U                | 5.3          |
| Endosulfan (I)     | U                | 3.9          | U                | 4.2          | U                | 3.7          | U                | 4.5          | U                | 5.3          |
| p,p'-DDE           | 3.9              | J 3.9        | 28               | 4.2          | 13               | 3.7          | 9.3              | 4.5          | U                | 5.3          |
| Dieldrin           | 1.1              | J 3.9        | 6.3              | 4.2          | 2.6              | J 3.7        | 2.4              | J 4.5        | U                | 5.3          |
| Endrin             | U                | 3.9          | U                | 4.2          | U                | 3.7          | U                | 4.5          | U                | 5.3          |
| p,p'-DDD           | U                | 3.9          | U                | 4.2          | U                | 3.7          | U                | 4.5          | U                | 5.3          |
| Endosulfan (II)    | U                | 3.9          | U                | 4.2          | U                | 3.7          | U                | 4.5          | U                | 5.3          |
| p,p'-DDT           | 5.2              | 3.9          | U                | 4.2          | U                | 3.7          | U                | 4.5          | U                | 5.3          |
| Endrin Aldehyde    | U                | 3.9          | U                | 4.2          | U                | 3.7          | U                | 4.5          | U                | 5.3          |
| Endosulfan Sulfate | U                | 3.9          | U                | 4.2          | U                | 3.7          | U                | 4.5          | U                | 5.3          |
| Methoxychlor       | U                | 3.9          | U                | 4.2          | U                | 3.7          | U                | 4.5          | U                | 5.3          |
| Endrin Ketone      | U                | 3.9          | U                | 4.2          | U                | 3.7          | U                | 4.5          | U                | 5.3          |
| Toxaphene          | U                | 99           | U                | 100          | U                | 92           | U                | 110          | U                | 130          |
| Aroclor 1016       | U                | 49           | U                | 52           | U                | 46           | U                | 57           | U                | 67           |
| Aroclor 1221       | U                | 99           | U                | 100          | U                | 92           | U                | 110          | U                | 130          |
| Aroclor 1232       | U                | 49           | U                | 52           | U                | 46           | U                | 57           | U                | 67           |
| Aroclor 1242       | U                | 49           | U                | 52           | U                | 46           | U                | 57           | U                | 67           |
| Aroclor 1248       | 1000             | 49           | 7400             | 52           | 3100             | 46           | 2800             | 57           | 42               | J 67         |
| Aroclor 1254       | U                | 49           | U                | 52           | 1300             | 46           | U                | 57           | U                | 67           |
| Aroclor 1260       | U                | 49           | U                | 52           | 240              | 46           | U                | 57           | U                | 67           |
| Aroclor 1268       | U                | 49           | U                | 52           | U                | 46           | U                | 57           | U                | 67           |

Table 1.2 (Cont) Results of the Analysis for Pesticide/PCBs in Soil  
 WA# 3-439 Lackawanna Foundry Site  
 Based on Dry Weight

| Client ID          | A16666         |              | A16667         |              | A16668         |              |
|--------------------|----------------|--------------|----------------|--------------|----------------|--------------|
| Location           | Location 16    |              | Location 17    |              | Location 18    |              |
| Percent Solid      | 89.6           |              | 87.8           |              | 92.6           |              |
| Analyte            | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg |
| a-BHC              | U              | 3.7          | U              | 3.8          | U              | 3.6          |
| g-BHC              | U              | 3.7          | U              | 3.8          | U              | 3.6          |
| b-BHC              | U              | 3.7          | U              | 3.8          | U              | 3.6          |
| Heptachlor         | U              | 3.7          | U              | 3.8          | U              | 3.6          |
| d-BHC              | U              | 3.7          | U              | 3.8          | U              | 3.6          |
| Aldrin             | U              | 3.7          | U              | 3.8          | U              | 3.6          |
| Heptachlor Epoxide | U              | 3.7          | U              | 3.8          | U              | 3.6          |
| g-Chlordane        | U              | 3.7          | U              | 3.8          | U              | 3.6          |
| a-Chlordane        | U              | 3.7          | U              | 3.8          | U              | 3.6          |
| Endosulfan (I)     | U              | 3.7          | U              | 3.8          | U              | 3.6          |
| p,p'-D D E         | 0.7 J          | 3.7          | 2.2 J          | 3.8          | U              | 3.6          |
| Dieldrin           | U              | 3.7          | 0.6 J          | 3.8          | U              | 3.6          |
| Endrin             | U              | 3.7          | U              | 3.8          | U              | 3.6          |
| p,p'-D D D         | U              | 3.7          | U              | 3.8          | U              | 3.6          |
| Endosulfan (II)    | U              | 3.7          | U              | 3.8          | U              | 3.6          |
| p,p'-D D T         | U              | 3.7          | U              | 3.8          | U              | 3.6          |
| Endrin Aldehyde    | U              | 3.7          | U              | 3.8          | U              | 3.6          |
| Endosulfan Sulfate | U              | 3.7          | U              | 3.8          | U              | 3.6          |
| Methoxychlor       | U              | 3.7          | U              | 3.8          | U              | 3.6          |
| Endrin Ketone      | U              | 3.7          | U              | 3.8          | U              | 3.6          |
| Toxaphene          | U              | 93           | U              | 95           | U              | 90           |
| Aroclor 1016       | U              | 47           | U              | 47           | U              | 45           |
| Aroclor 1221       | U              | 93           | U              | 95           | U              | 90           |
| Aroclor 1232       | U              | 47           | U              | 47           | U              | 45           |
| Aroclor 1242       | 340            | 47           | U              | 47           | U              | 45           |
| Aroclor 1248       | U              | 47           | 620            | 47           | 77             | 45           |
| Aroclor 1254       | U              | 47           | U              | 47           | 33 J           | 45           |
| Aroclor 1260       | U              | 47           | U              | 47           | U              | 45           |
| Aroclor 1268       | U              | 47           | U              | 47           | U              | 45           |

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Table 1.2 (Cont) Results of the Analysis for Pesticide/PCBs in Soil  
 WA# 3-439 Lackawanna Foundry Site  
 Based on Dry Weight

| Client ID          | SBLK042099     |              | A16669         |              | A16670         |              | A16671         |              | A16672         |              |
|--------------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|----------------|--------------|
| Location           | -              |              | Location 19    |              | Location 20    |              | Location 21    |              | Location 22    |              |
| Percent Solid      | 100            |              | 93.3           |              | 90.9           |              | 48.4           |              | 81.4           |              |
| Analyte            | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg |
| a-BHC              | U              | 3.3          | U              | 3.6          | U              | 3.7          | U              | 6.9          | U              | 4.1          |
| g-BHC              | U              | 3.3          | U              | 3.6          | 2.5 J          | 3.7          | U              | 6.9          | U              | 4.1          |
| b-BHC              | U              | 3.3          | U              | 3.6          | U              | 3.7          | U              | 6.9          | U              | 4.1          |
| Heptachlor         | U              | 3.3          | U              | 3.6          | U              | 3.7          | U              | 6.9          | U              | 4.1          |
| d-BHC              | U              | 3.3          | U              | 3.6          | U              | 3.7          | U              | 6.9          | U              | 4.1          |
| Aldrin             | U              | 3.3          | U              | 3.6          | U              | 3.7          | U              | 6.9          | U              | 4.1          |
| Heptachlor Epoxide | U              | 3.3          | U              | 3.6          | U              | 3.7          | U              | 6.9          | U              | 4.1          |
| g-Chlordane        | U              | 3.3          | U              | 3.6          | U              | 3.7          | U              | 6.9          | U              | 4.1          |
| a-Chlordane        | U              | 3.3          | U              | 3.6          | U              | 3.7          | U              | 6.9          | U              | 4.1          |
| Endosulfan (I)     | U              | 3.3          | U              | 3.6          | U              | 3.7          | U              | 6.9          | U              | 4.1          |
| p,p'-DDE           | U              | 3.3          | 5.1            | 3.6          | U              | 3.7          | 8.4            | 6.9          | 2.3 J          | 4.1          |
| Dieldrin           | U              | 3.3          | 2.2 J          | 3.6          | U              | 3.7          | 21             | 6.9          | U              | 4.1          |
| Endrin             | U              | 3.3          | U              | 3.6          | U              | 3.7          | U              | 6.9          | U              | 4.1          |
| p,p'-DDD           | U              | 3.3          | U              | 3.6          | U              | 3.7          | U              | 6.9          | 1.6 J          | 4.1          |
| Endosulfan (II)    | U              | 3.3          | U              | 3.6          | U              | 3.7          | U              | 6.9          | U              | 4.1          |
| p,p'-DDT           | U              | 3.3          | 4.9            | 3.6          | U              | 3.7          | U              | 6.9          | 1.9 J          | 4.1          |
| Endrin Aldehyde    | U              | 3.3          | U              | 3.6          | U              | 3.7          | U              | 6.9          | U              | 4.1          |
| Endosulfan Sulfate | U              | 3.3          | U              | 3.6          | U              | 3.7          | U              | 6.9          | U              | 4.1          |
| Methoxychlor       | U              | 3.3          | U              | 3.6          | U              | 3.7          | U              | 6.9          | U              | 4.1          |
| Endrin Ketone      | U              | 3.3          | U              | 3.6          | U              | 3.7          | U              | 6.9          | U              | 4.1          |
| Toxaphene          | U              | 83           | U              | 89           | U              | 92           | U              | 170          | U              | 100          |
| Aroclor 1016       | U              | 42           | U              | 45           | U              | 46           | U              | 86           | U              | 51           |
| Aroclor 1221       | U              | 83           | U              | 89           | U              | 92           | U              | 170          | U              | 100          |
| Aroclor 1232       | U              | 42           | U              | 45           | U              | 46           | U              | 86           | U              | 51           |
| Aroclor 1242       | U              | 42           | 2500           | 45           | U              | 46           | U              | 86           | U              | 51           |
| Aroclor 1248       | U              | 42           | U              | 45           | U              | 46           | 320            | 86           | 11 W           | 51           |
| Aroclor 1254       | U              | 42           | U              | 45           | U              | 46           | 300            | 86           | 16 W           | 51           |
| Aroclor 1260       | U              | 42           | U              | 45           | 3900           | 46           | 200            | 86           | U              | 51           |
| Aroclor 1268       | U              | 42           | U              | 45           | U              | 46           | U              | 86           | U              | 51           |

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Table 1.2 (Cont) Results of the Analysis for Pesticide/PCBs in Soil  
 WA# 3-433 Lackawanna Foundry Site  
 Based on Dry Weight

| Client ID<br>Location<br>Percent Solid | A16673<br>Location 23<br>82.2 |              | A16674<br>Location 24<br>75.4 |              | A16675<br>Location 25<br>82.9 |              | A16676<br>Location 25-D<br>77.2 |              | A16677<br>Location 26<br>69.5 |              |
|--|-------------------------------|--------------|-------------------------------|--------------|-------------------------------|--------------|---------------------------------|--------------|-------------------------------|--------------|
|  | Conc.<br>µg/kg                | MDL<br>µg/kg | Conc.<br>µg/kg                | MDL<br>µg/kg | Conc.<br>µg/kg                | MDL<br>µg/kg | Conc.<br>µg/kg                  | MDL<br>µg/kg | Conc.<br>µg/kg                | MDL<br>µg/kg |
| a-BHC                                  | U                             | 4.1          | U                             | 4.4          | U                             | 4.0          | U                               | 4.3          | U                             | 4.8          |
| g-BHC                                  | U                             | 4.1          | U                             | 4.4          | U                             | 4.0          | U                               | 4.3          | U                             | 4.8          |
| b-BHC                                  | U                             | 4.1          | U                             | 4.4          | U                             | 4.0          | U                               | 4.3          | U                             | 4.8          |
| Heptachlor                             | U                             | 4.1          | U                             | 4.4          | U                             | 4.0          | U                               | 4.3          | U                             | 4.8          |
| d-BHC                                  | U                             | 4.1          | U                             | 4.4          | U                             | 4.0          | U                               | 4.3          | U                             | 4.8          |
| Aldrn                                  | U                             | 4.1          | U                             | 4.4          | U                             | 4.0          | U                               | 4.3          | U                             | 4.8          |
| Heptachlor Epoxide                     | U                             | 4.1          | U                             | 4.4          | U                             | 4.0          | U                               | 4.3          | U                             | 4.8          |
| g-Chlordane                            | U                             | 4.1          | U                             | 4.4          | U                             | 4.0          | U                               | 4.3          | U                             | 4.8          |
| a-Chlordane                            | U                             | 4.1          | U                             | 4.4          | U                             | 4.0          | U                               | 4.3          | U                             | 4.8          |
| Endosulfan (I)                         | U                             | 4.1          | U                             | 4.4          | U                             | 4.0          | U                               | 4.3          | U                             | 4.8          |
| p,p'-D D E                             | 2.5 J                         | 4.1          | 1.9 J                         | 4.4          | 3.3 J                         | 4.0          | U                               | 4.3          | 11                            | 4.8          |
| Dieldrn                                | U                             | 4.1          | U                             | 4.4          | U                             | 4.0          | U                               | 4.3          | U                             | 4.8          |
| Endrn                                  | U                             | 4.1          | U                             | 4.4          | U                             | 4.0          | U                               | 4.3          | U                             | 4.8          |
| p,p'-D D D                             | U                             | 4.1          | U                             | 4.4          | U                             | 4.0          | U                               | 4.3          | U                             | 4.8          |
| Endosulfan (II)                        | U                             | 4.1          | U                             | 4.4          | U                             | 4.0          | U                               | 4.3          | U                             | 4.8          |
| p,p'-D D T                             | U                             | 4.1          | U                             | 4.4          | U                             | 4.0          | U                               | 4.3          | 18                            | 4.8          |
| Endrn Aldehyde                         | U                             | 4.1          | U                             | 4.4          | U                             | 4.0          | U                               | 4.3          | U                             | 4.8          |
| Endosulfan Sulfate                     | U                             | 4.1          | U                             | 4.4          | U                             | 4.0          | U                               | 4.3          | U                             | 4.8          |
| Methoxychlor                           | U                             | 4.1          | U                             | 4.4          | U                             | 4.0          | U                               | 4.3          | U                             | 4.8          |
| Endrn Ketone                           | U                             | 4.1          | U                             | 4.4          | U                             | 4.0          | U                               | 4.3          | U                             | 4.8          |
| Toxaphene                              | U                             | 100          | U                             | 110          | U                             | 100          | U                               | 110          | U                             | 120          |
| Aroclor 1016                           | U                             | 51           | U                             | 55           | U                             | 50           | U                               | 54           | U                             | 60           |
| Aroclor 1221                           | U                             | 100          | U                             | 110          | U                             | 100          | U                               | 110          | U                             | 120          |
| Aroclor 1232                           | U                             | 51           | U                             | 55           | U                             | 50           | U                               | 54           | U                             | 60           |
| Aroclor 1242                           | U                             | 51           | U                             | 55           | U                             | 50           | U                               | 54           | U                             | 60           |
| Aroclor 1248                           | 110                           | 51           | 120                           | 55           | 270                           | 50           | U                               | 54           | 170                           | 60           |
| Aroclor 1254                           | U                             | 51           | U                             | 55           | 83                            | 50           | U                               | 54           | U                             | 60           |
| Aroclor 1260                           | 14 J                          | 51           | 10 J                          | 55           | U                             | 50           | U                               | 54           | 42 J                          | 60           |
| Aroclor 1268                           | U                             | 51           | U                             | 55           | U                             | 50           | U                               | 54           | U                             | 60           |

00031

Table 1.2 (Cont) Results of the Analysis for Pesticide/PCBs in Soil  
 WA# 3-439 Lackawanna Foundry Site  
 Based on Dry Weight

| Client ID          | A16678        |              | A16679         |              |
|--------------------|---------------|--------------|----------------|--------------|
| Location           | Location 26-D |              | Location 27    |              |
| Percent Solid      | 76.5          |              | 83.2           |              |
| Analyte            | Conc<br>µg/kg | MDL<br>µg/kg | Conc.<br>µg/kg | MDL<br>µg/kg |
| a-BHC              | U             | 4.4          | U              | 4.0          |
| g-BHC              | U             | 4.4          | U              | 4.0          |
| b-BHC              | U             | 4.4          | U              | 4.0          |
| Heptachlor         | U             | 4.4          | U              | 4.0          |
| d-BHC              | U             | 4.4          | U              | 4.0          |
| Aldnn              | U             | 4.4          | U              | 4.0          |
| Heptachlor Epoxide | U             | 4.4          | U              | 4.0          |
| g-Chlordane        | U             | 4.4          | U              | 4.0          |
| a-Chlordane        | U             | 4.4          | U              | 4.0          |
| Endosulfan (I)     | U             | 4.4          | U              | 4.0          |
| p,p'-D D E         | U             | 4.4          | U              | 4.0          |
| Dieldnn            | U             | 4.4          | U              | 4.0          |
| Endrin             | U             | 4.4          | U              | 4.0          |
| p,p'-D D D         | U             | 4.4          | U              | 4.0          |
| Endosulfan (II)    | U             | 4.4          | U              | 4.0          |
| p,p'-D D T         | U             | 4.4          | U              | 4.0          |
| Endrin Aldehyde    | U             | 4.4          | U              | 4.0          |
| Endosulfan Sulfate | U             | 4.4          | U              | 4.0          |
| Methoxychlor       | U             | 4.4          | U              | 4.0          |
| Endrin Ketone      | U             | 4.4          | U              | 4.0          |
| Toxaphene          | U             | 110          | U              | 100          |
| Aroclor 1016       | U             | 54           | U              | 50           |
| Aroclor 1221       | U             | 110          | U              | 100          |
| Aroclor 1232       | U             | 54           | U              | 50           |
| Aroclor 1242       | U             | 54           | U              | 50           |
| Aroclor 1248       | U             | 54           | 39000          | 50           |
| Aroclor 1254       | U             | 54           | U              | 50           |
| Aroclor 1260       | U             | 54           | U              | 50           |
| Aroclor 1268       | U             | 54           | U              | 50           |

0003

## QA/QC for VOC

### Results of the Surrogate Percent Recoveries for VOC in Soil

Each sample was spiked with a three component mixture of CLP surrogate standards consisting of dibromofluoromethane, toluene-d<sub>8</sub> and 4-bromofluorobenzene. The surrogate percent recoveries, listed in Table 2.1, ranged from 65 to 156. One hundred and ten out of one hundred and twenty values were within the acceptable QC limits.

### Results of the LCS/LCSD Analysis for VOC in Soil

A laboratory control sample was analyzed in duplicate. The percent recoveries, ranging from 102 to 119, are listed in Table 2.2. All ten values were within the acceptable QC limits. The relative percent differences, also listed in Table 2.2, ranged from 5 to 12, and all five values were within the acceptable QC limits.

### Results of the LCS Analysis for VOC in Soil

Two laboratory control samples were analyzed. The percent recoveries, ranging from 92 to 112, are listed in Table 2.3. All ten values were within the acceptable QC limits.

### Results of the MS/MSD Analysis for VOC in Soil

Sample B 16665 was chosen for the matrix spike/matrix spike duplicate (MS/MSD) analyses. The percent recoveries, ranging from 65 to 100, are listed in Table 2.4. Eight out of ten values were within the acceptable QC limits. The relative percent differences, also listed in Table 2.4, ranged from 11 to 17, and all five were within the acceptable QC limits.

**Table 2.1 Results of the Surrogate Recoveries  
for VOC in Soil  
WA # 3-439 Lackawanna Foundry Site**

| Sample ID              | S1<br>(DBFM) | S2<br>(TOL) | S3<br>(BFB) | Total<br>Out |
|------------------------|--------------|-------------|-------------|--------------|
| B16650                 | 86           | 89          | 88          | 0            |
| B16651                 | 98           | 112         | 105         | 0            |
| B16652                 | 91           | 101         | 91          | 0            |
| B16653                 | 100          | 100         | 110         | 0            |
| B16655                 | 88           | 86          | 75          | 0            |
| B16656                 | 91           | 92          | 99          | 0            |
| B16657                 | 89           | 88          | 94          | 0            |
| B16658                 | 104          | 98          | 108         | 0            |
| Method Blank (4/23/99) | 92           | 110         | 90          | 0            |
| LCSD (4/23/99)         | 106          | 111         | 101         | 0            |
| LCS (4/23/99)          | 103          | 114         | 104         | 0            |
| B16659                 | 84           | 85          | 81          | 0            |
| B16660                 | 113          | 105         | 113         | 0            |
| B16661                 | 91           | 87          | 104         | 0            |
| B16662                 | 90           | 85          | 80          | 0            |
| B16663                 | 101          | 114         | 98          | 0            |
| B16664                 | 86           | 86          | 83          | 0            |
| B16654                 | 84           | 87          | 82          | 0            |
| Method Blank (4/26/99) | 102          | 111         | 100         | 0            |
| LCS (4/26/99)          | 90           | 107         | 92          | 0            |
| B16665 MSD             | 100          | 102         | 106         | 0            |
| B16665                 | 105          | 103         | 98          | 0            |
| B16665 MS              | 90           | 75          | 74          | 1            |
| B16666                 | 105          | 102         | 106         | 0            |
| B16667                 | 90           | 89          | 79          | 0            |
| B16668                 | 101          | 106         | 103         | 0            |
| B16669                 | 89           | 80          | 65          | 2            |
| B16670                 | 93           | 75          | 68          | 2            |
| B16671                 | 112          | 85          | 114         | 0            |
| B16672                 | 110          | 97          | 99          | 0            |
| B16673                 | 107          | 74          | 89          | 1            |
| B16674                 | 142          | 117         | 156         | 3            |
| B16675                 | 124          | 69          | 96          | 1            |
| B16676                 | 104          | 103         | 97          | 0            |
| B16677                 | 120          | 99          | 117         | 0            |
| B16678                 | 94           | 93          | 75          | 0            |
| B16679                 | 109          | 99          | 106         | 0            |
| B16680                 | 110          | 112         | 99          | 0            |
| Method Blank (4/27/99) | 98           | 104         | 86          | 0            |
| LCS (4/27/99)          | 99           | 106         | 94          | 0            |

QC Limits

S1 (DBFM) = Dibromofluoromethane      76-127  
 S3 (TOL) = Toluene-d8                      83-115  
 S2 (BFB) = Bromofluorobenzene        74-134

00034

Table 2.2 Results of the LCS/LCSD Analysis for VOC in Soil  
 WA # 3-439 Lackawanna Foundry Site

Date of Analysis 4/23/99

| Analyte            | Sample Conc<br>µg/kg | LCS Spike Added<br>µg/kg | LCS Conc<br>µg/kg | LCS %<br>Rec | LCSD Spike Added<br>µg/kg | LCSD Conc<br>µg/kg | LCSD %<br>Rec | RPD | Recommended QC Limits |     |
|--------------------|----------------------|--------------------------|-------------------|--------------|---------------------------|--------------------|---------------|-----|-----------------------|-----|
|                    |                      |                          |                   |              |                           |                    |               |     | %Rec                  | RPD |
| 1,1-Dichloroethene | U                    | 20                       | 21.2              | 106          | 20                        | 23.9               | 119           | 12  | 59-136                | 25  |
| Benzene            | U                    | 20                       | 20.9              | 105          | 20                        | 22.0               | 110           | 5   | 76-123                | 23  |
| Trichloroethene    | U                    | 20                       | 21.5              | 108          | 20                        | 22.7               | 114           | 5   | 74-127                | 25  |
| Toluene            | U                    | 20                       | 22.3              | 112          | 20                        | 23.4               | 117           | 5   | 75-124                | 24  |
| Chlorobenzene      | U                    | 20                       | 20.4              | 102          | 20                        | 21.5               | 108           | 5   | 75-124                | 24  |

00035



Table 2.3 Results of the LCS Analysis for VOC in Soil  
WA # 3-439 Lackawanna Foundry Site

Date of Analysis 4/26/99

| Analyte            | Spike<br>Conc<br>µg/kg | Rec<br>Conc<br>µg/kg | %<br>Rec | QC<br>Limits<br>% Rec |
|--------------------|------------------------|----------------------|----------|-----------------------|
| 1,1-Dichloroethene | 20                     | 19.2                 | 96       | 59-136                |
| Benzene            | 20                     | 19.9                 | 99       | 76-123                |
| Trichloroethene    | 20                     | 19.9                 | 99       | 74-127                |
| Toluene            | 20                     | 20.7                 | 103      | 75-124                |
| Chlorobenzene      | 20                     | 19.7                 | 99       | 75-124                |

Date of Analysis 4/27/99

| Analyte            | Spike<br>Conc<br>µg/kg | Rec<br>Conc<br>µg/kg | %<br>Rec | QC<br>Limits<br>% Rec |
|--------------------|------------------------|----------------------|----------|-----------------------|
| 1,1-Dichloroethene | 20                     | 21.4                 | 107      | 59-136                |
| Benzene            | 20                     | 20.7                 | 103      | 76-123                |
| Trichloroethene    | 20                     | 21.0                 | 105      | 74-127                |
| Toluene            | 20                     | 22.3                 | 112      | 75-124                |
| Chlorobenzene      | 20                     | 18.3                 | 92       | 75-124                |

00036

Table 2.4 Results of the MS/MSD Analysis for VOC in Soil  
 WA # 3-439 Lackawanna Foundry Site  
 Results Based on Dry Weight

Sample ID B 16665

| Analyte            | Sample Conc<br>µg/kg | MS                   |                  |             | MSD                  |                   |              | RPD | Recommended QC Limits |     |
|--------------------|----------------------|----------------------|------------------|-------------|----------------------|-------------------|--------------|-----|-----------------------|-----|
|                    |                      | Spike Added<br>µg/kg | MS Conc<br>µg/kg | MS %<br>Rec | Spike Added<br>µg/kg | MSD Conc<br>µg/kg | MSD %<br>Rec |     | %Rec                  | RPD |
| 1,1-Dichloroethene | U                    | 24.1                 | 24.2             | 100         | 24.1                 | 21.2              | 88           | 13  | 59-136                | 25  |
| Benzene            | U                    | 24.1                 | 24               | 100         | 24.1                 | 21.6              | 90           | 11  | 76-123                | 23  |
| Trichloroethene    | U                    | 24.1                 | 20.2             | 84          | 24.1                 | 17.4              | 72           | 15  | 74-127                | 25  |
| Toluene            | U                    | 24.1                 | 21.9             | 91          | 24.1                 | 19.7              | 82           | 11  | 75-124                | 24  |
| Chlorobenzene      | U                    | 24.1                 | 18.7             | 78          | 24.1                 | 15.7              | 65           | 17  | 75-124                | 24  |

00007

## QA/QC for Pesticides/PCBs

### Results of the Surrogate Recoveries for Pesticides/PCBs in Soil

Before extraction, each sample was spiked with a two component mixture of CLP surrogate standards consisting of tetrachloro-*m*-xylene and dichlorobiphenyl. The surrogate percent recoveries, listed in Table 2.5, ranged from 21 to 128. Forty-seven out of seventy-eight values were within the acceptable QC limits.

### Results of the MS/MSD Analysis for Pesticides/PCBs in Soil

Samples A 16653, A 16665 and A 16677 were chosen for the matrix spike/matrix spike duplicate (MS/MSD) analyses. The percent recoveries, ranging from 1.0 to 111, are listed in Table 2.6. Twenty-eight out of thirty-six values were within the acceptable QC limits. The relative percent differences, also listed in Table 2.6, ranged from 1 to 100 and fourteen out of eighteen values were within the acceptable QC limits.

Table 2.5 Results of the Surrogate Recoveries  
for Pesticides/PCBs in Soil  
WA# 3-439 Lackawanna Foundry Site

| Sample ID  | Percent Recovery |      |
|------------|------------------|------|
|            | TCMX             | DCBP |
| SBLK041599 | 89               | 86   |
| A16650     | 59 *             | 36 * |
| A16651     | 88               | 72   |
| A16652     | 71               | 21 * |
| A16653     | 83               | 49 * |
| A16653MS   | 91               | 43 * |
| A16653MSD  | 85               | 33 * |
| A16654     | 85               | 28 * |
| A16655     | 79               | 24 * |
| A16656     | 77               | 27 * |
| SBLK041999 | 99               | 95   |
| A16657     | 75               | 68   |
| A16658     | 81               | 52 * |
| A16659     | 101              | 77   |
| A16660     | 74               | 45 * |
| A16661     | 74               | 43 * |
| A16662     | 86               | 55 * |
| A16663     | 79               | 43 * |
| A16664     | 58 *             | 31 * |
| A16665     | 87               | 43 * |
| A16665MS   | 91               | 37 * |
| A16665MSD  | 96               | 53 * |
| A16666     | 67               | 31 * |
| A16667     | 81               | 39 * |
| A16668     | 70               | 42 * |
| SBLK042099 | 86               | 80   |
| A16669     | 90               | 80   |
| A16670     | 89               | 90   |
| A16671     | 84               | 76   |
| A16672     | 87               | 50 * |
| A16673     | 83               | 128  |
| A16674     | 68               | 40 * |
| A16675     | 78               | 54 * |
| A16676     | 75               | 46 * |
| A16677     | 84               | 52 * |
| A16677MS   | 72               | 30 * |
| A16677MSD  | 75               | 34 * |
| A16678     | 62               | 29 * |
| A16679     | 63               | 36 * |

TCMX denotes Tetrachloro-m-xylene  
DCBP denotes Decachlorobiphenyl

|      |          |
|------|----------|
|      | Advisory |
|      | QC       |
|      | Limits   |
| TCMX | 60-150   |
| DCBP | 60-150   |

03023

**Table 2.6 Results of the MS/MSD Analysis for Pesticide/PCB in Soil**  
**WA# 3-439 Lackawanna Foundry Site**  
**Results Based on Dry Weight**

Sample ID: A16653

| Compound   | Sample Conc<br>µg/kg | MS                   |                  |             | MSD                  |                   |              | RPD  | Advisory QC Limits |     |
|------------|----------------------|----------------------|------------------|-------------|----------------------|-------------------|--------------|------|--------------------|-----|
|            |                      | Spike Added<br>µg/kg | MS Conc<br>µg/kg | MS %<br>Rec | Spike Added<br>µg/kg | MSD Conc<br>µg/kg | MSD %<br>Rec |      | % Rec              | RPD |
| g-BHC      | U                    | 28.461               | 3.198            | 11          | * 28.461             | 8.142             | 29           | * 90 | * 46-127           | 50  |
| Heptachlor | U                    | 28.461               | 26.281           | 92          | 28.461               | 21.596            | 76           | 19   | 35-130             | 31  |
| Aldrin     | U                    | 28.461               | 25.157           | 88          | 28.461               | 24.236            | 85           | 3    | 34-132             | 43  |
| Dieldrin   | U                    | 56.922               | 54.380           | 96          | 56.922               | 49.607            | 87           | 10   | 31-134             | 38  |
| Endrin     | U                    | 56.922               | 58.717           | 103         | 56.922               | 52.604            | 92           | 11   | 42-139             | 45  |
| p,p'-DDT   | 13.718               | 56.922               | 15.537           | 3.0         | * 56.922             | 16.418            | 5.0          | * 50 | 23-134             | 50  |

Sample ID: A16665

| Compound   | Sample Conc<br>µg/kg | MS                   |                  |             | MSD                  |                   |              | RPD | Advisory QC Limits |     |
|------------|----------------------|----------------------|------------------|-------------|----------------------|-------------------|--------------|-----|--------------------|-----|
|            |                      | Spike Added<br>µg/kg | MS Conc<br>µg/kg | MS %<br>Rec | Spike Added<br>µg/kg | MSD Conc<br>µg/kg | MSD %<br>Rec |     | % Rec              | RPD |
| g-BHC      | U                    | 33.387               | 31.949           | 96          | 33.387               | 31.775            | 95           | 1   | 46-127             | 50  |
| Heptachlor | U                    | 33.387               | 28.996           | 87          | 33.387               | 32.089            | 96           | 10  | 35-130             | 31  |
| Aldrin     | U                    | 33.387               | 33.104           | 99          | 33.387               | 34.538            | 103          | 4   | 34-132             | 43  |
| Dieldrin   | U                    | 66.774               | 65.005           | 97          | 66.774               | 71.198            | 107          | 10  | 31-134             | 38  |
| Endrin     | U                    | 66.774               | 68.919           | 103         | 66.774               | 74.270            | 111          | 7   | 42-139             | 45  |
| p,p'-DDT   | U                    | 66.774               | 13.065           | 20          | * 66.774             | 23.275            | 35           | 55  | * 23-134           | 50  |

Sample ID: A16677

| Compound   | Sample Conc<br>µg/kg | MS                   |                  |             | MSD                  |                   |              | RPD   | Advisory QC Limits |     |
|------------|----------------------|----------------------|------------------|-------------|----------------------|-------------------|--------------|-------|--------------------|-----|
|            |                      | Spike Added<br>µg/kg | MS Conc<br>µg/kg | MS %<br>Rec | Spike Added<br>µg/kg | MSD Conc<br>µg/kg | MSD %<br>Rec |       | % Rec              | RPD |
| g-BHC      | U                    | 29.976               | 6.866            | 23          | * 29.976             | 14.024            | 47           | 69    | * 46-127           | 50  |
| Heptachlor | U                    | 29.976               | 15.589           | 52          | 29.976               | 15.338            | 51           | 2     | 35-130             | 31  |
| Aldrin     | U                    | 29.976               | 21.583           | 72          | 29.976               | 21.936            | 73           | 1     | 34-132             | 43  |
| Dieldrin   | U                    | 59.952               | 42.495           | 71          | 59.952               | 41.038            | 68           | 4     | 31-134             | 38  |
| Endrin     | U                    | 59.952               | 36.742           | 61          | 59.952               | 36.041            | 60           | 2     | 42-139             | 45  |
| p,p'-DDT   | 17.538               | 59.952               | 19.274           | 3.0         | * 59.952             | 17.846            | 1.0          | * 100 | * 23-134           | 50  |



Roy F. Weston, Inc.  
 GSA Rantan Depot  
 Bldg. 209 Annex (Bay F)  
 2890 Woodbridge Avenue  
 Edison, New Jersey 08837-3679  
 732-321-4200 • Fax 732-494-4021

**Paragon Analytics, Inc.**  
 225 Commerce Drive  
 Fort Collins, CO 80524

Attn: Claudia Clark

20 April 1999

Project # 3347-143-001-3439 Lackawanna Foundry Site

As per Weston REAC Purchase Order number 102755, please analyze samples according to the following parameters:

| Analysis/Method                                 | Matrix | # of samples |
|---|--------|--------------|
| VOA 5030\ SW-846-8260B                          | Soil   | 32           |
| Data package: Package with Diskette Deliverable |        |              |

Samples are expected to arrive at your laboratory on April 20, 1999. **All applicable QA/QC (MS/MSD) analysis as per method, will be performed on our sample matrix. Preliminary sample and QC result tables plus a signed copy of our Chain of Custody must be faxed to REAC 10 business days after receipt of the last samples.** The complete data package is due 21 business days after receipt of last batch of samples. The complete data package must include all items on the deliverables checklist. **Expect all samples to be difficult matrix and all raw data must be included in final analytical report.**

All sample and QC results (ie: MS/MSD, LCS, Duplicates, and Blanks) must be summarized in a ExCel diskette deliverable.

Please submit all reports and technical questions concerning this project to **John Johnson** at (732) 321-4248 or fax to (732) 494-4020. Any contractual question, please call Cynthia Lentini at (732) 321-4296.

Sincerely,

Misty Barkley  
 Data Validation and Report Writing Group Leader  
 Roy F. Weston, Inc. / REAC Project

MB:jj Attachments

cc. R. Singhvi  
 P. Campagna  
 3439\non\mem\9904\sub\3439Con2

V. Kansal  
 Subcontracting File  
 D. Angwenyi

C. Lentini  
 S. Fama  
 M. Barkley

000-11



Lackawanna Foundry VOA List

Dichlorodifluoromethane  
Chloromethane  
Vinyl Chloride  
Bromomethane  
Chloroethane  
Trichlorofluoromethane  
Acetone  
1,1-Dichloroethene  
Carbon Disulfide  
Methylene Chloride  
Methyl-tertiary-butylether  
trans-1,2-Dichloroethene  
1,1-Dichloroethane  
2-Butanone  
2,2-Dichloropropane  
cis-1,2-Dichloroethene  
Chloroform  
1,1-Dichloropropene  
1,2-Dichloroethane  
1,1,1-Trichloroethane  
Carbon Tetrachloride  
Benzene  
Trichloroethene  
1,2-Dichloropropane  
Dibromomethane  
Bromodichloromethane  
cis-1,3-Dichloropropene  
trans-1,3-Dichloropropene  
1,1,2-Trichloroethane  
1,3-Dichloropropane  
Dibromochloromethane  
1,2-Dibromoethane  
Bromoform  
4-Methyl-2-Pentanone  
Toluene  
2-Hexanone  
Tetrachloroethene  
Chlorobenzene  
1,1,1,2-Tetrachloroethane  
Ethylbenzene  
p & m-Xylene  
o-Xylene  
Styrene  
Isopropylbenzene  
1,1,2,2-Tetrachloroethane  
1,2,3-Trichloropropane  
Bromobenzene  
n-Propylbenzene  
2-Chlorotoluene  
4-Chlorotoluene  
1,3,5-Trimethylbenzene  
tert-Butylbenzene  
1,2,4-Trimethylbenzene  
sec-Butylbenzene  
1,3-Dichlorobenzene  
p-Isopropyltoluene  
1,4-Dichlorobenzene  
1,2-Dichlorobenzene  
n-Butylbenzene  
1,2-Dibromo-3-Chloropropane  
1,2,4-Trichlorobenzene  
Naphthalene  
Hexachlorobutadiene  
1,2,3-Trichlorobenzene

REAC, L. jon, NJ  
 (908) 321-4200  
 EPA Contract 68-C4-0022

**CHA OF CUSTODY RECORD**  
 Project Name: Lackawanna Foundry Site  
 Project Number: 03347-143-001-3439-01  
 RFW Contact: Johnson, Fama Phone: (732) 421-4200

No: 06505

SHEET NO. 1 OF 2

**Sample Identification**

**Analyses Requested**

| REAC # | Sample No. | Sampling Location | Matrix | Date Collected | # of Bottles | Container/Preservative | VOC |
|--------|------------|-------------------|--------|----------------|--------------|------------------------|-----|
| 755    | B16650     | 1 West Fence      | S      | 4/14/99        | 1            | 40mL / 4°C             | X   |
| 758    | B16651     | 2 Surface Middle  | S      |                |              |                        |     |
| 761    | B16652     | 2 Middle          | S      |                |              |                        |     |
| 764    | B16653     | 3 West Fence      | S      |                |              |                        |     |
| 767    | B16654     | 4 West Fence      | S      |                |              |                        |     |
| 770    | B16655     | 5 West Fence      | S      |                |              |                        |     |
| 773    | B16656     | 5 West Entrance   | S      |                |              |                        |     |
| 776    | B16657     | 7                 | S      |                |              |                        |     |
| 779    | B16658     | 8                 | S      |                |              |                        |     |
| 782    | B16659     | 9                 | S      |                |              |                        |     |
| 785    | B16660     | 10                | S      |                |              |                        |     |
| 788    | B16661     | 11                | S      |                |              |                        |     |
| 791    | B16662     | 12                | S      |                |              |                        |     |
| 794    | B16663     | 13                | S      |                |              |                        |     |
| 797    | B16664     | 14                | S      |                |              |                        |     |
| 800    | B16665     | 15                | S      |                |              |                        |     |

Special Instructions:

- Matrix: Sediment SD -
- Drum Solids DS -
- Drum Liquids DL -
- Other X -
- Potable Water PW -
- Groundwater GW -
- Surface Water SW -
- Sludge SL -
- Soil S -
- Water W -
- Oil O -
- Air A -

Lab to select Ms/MSD

**FOR SUBCONTRACTING USE ONLY**

**FROM CHAIN OF CUSTODY # 06908, 06909, 06910**

SUB OUT TO:

Perragon Analyticals, Inc  
 4/19/99

| Items/Reason | Relinquished By | Date    | Received By    | Date    | Relinquished By | Date    | Received By | Date | Time |
|--------------|-----------------|---------|----------------|---------|-----------------|---------|-------------|------|------|
| ALL ANALYSIS | John Johnson    | 4/19/99 | Daniel Johnson | 4/19/99 | Daniel Johnson  | 4/19/99 |             |      |      |
|              | John Johnson    |         |                |         |                 |         |             |      |      |
|              |                 |         |                |         |                 |         |             |      |      |
|              |                 |         |                |         |                 |         |             |      |      |



REAC, Edison, NJ  
 (908) 321-4200  
 EPA Contract 68-C4-0022

9906134

**CHAIN OF CUSTODY RECORD**

Project Name: Lackawanna Foundry Site  
 Project Number: 03347-143-001-01  
 RFW Contact: Johnson, Ferra Phone: (732) 421-4200

No: 06507

SHEET NO. 2 OF 2

**Sample Identification**

| REAC # | Sample No. | Sampling Location | Matrix | Date Collected | # of Bottles | Container/Preservative | VOC |
|--------|------------|-------------------|--------|----------------|--------------|------------------------|-----|
| 803    | B16666     | Location 16       | S      | 4/14/99        | 1            | 40ml VOC               | X   |
| 806    | B16667     | Location 17       |        |                |              |                        |     |
| 809    | B16668     | Location 18       |        |                |              |                        |     |
| 812    | B16669     | Location 19       |        |                |              |                        |     |
| 815    | B16670 PA  | Location 20       |        |                |              |                        |     |
| 818    | B16671     | Location 21       |        |                |              |                        |     |
| 821    | B16672     | Location 22       |        |                |              |                        |     |
| 824    | B16673     | Location 23       |        |                |              |                        |     |
| 827    | B16674     | Location 24       |        |                |              |                        |     |
| 830    | B16675     | Location 25       |        |                |              |                        |     |
| 833    | B16676     | Location 25-D     |        |                |              |                        |     |
| 836    | B16677     | Location 26       |        |                |              |                        |     |
| 839    | B16678     | Location 26D      |        |                |              |                        |     |
| 842    | B16679     | Location 27       |        |                |              |                        |     |
| 844    | B16680     | Field Blank (VOC) |        |                |              |                        |     |
| 845    | B16680     | Field Blank (VOC) |        |                |              |                        |     |

**Analyses Requested**

Special Instructions:

- Matrix: SD - Sediment, DS - Drum Solids, DL - Drum Liquids, X - Other  
 PW - Potable Water, GW - Groundwater, SW - Surface Water, SL - Sludge  
 S - Soil, W - Water, O - Oil, A - Air

Lab to Select  
 MS/MSD

**FOR SUBCONTRACTING USE ONLY**

**FROM CHAIN OF**

**CUSTODY # 06911, 06912, 06913**

| Items/Reason | Relinquished By    | Date    | Received By       | Date    | Time | Relinquished By   | Date    | Received By | Date | Time |
|--------------|--------------------|---------|-------------------|---------|------|-------------------|---------|-------------|------|------|
| ALL ANALYSIS | <i>[Signature]</i> | 4/19/99 | Daniel Angewinkel | 4/19/99 | 1350 | Daniel Angewinkel | 4/19/99 | FERRA       |      |      |
|              | <i>[Signature]</i> | 4/28/99 | G. Williams       | 4/28/99 | 900  |                   |         |             |      |      |



REAC, Edison, NJ  
 (908) 321-4200  
 EPA Contract 68-C4-0022

**CHAIN OF CUSTODY RECORD**

Project Name: Lacklawn Foundry Site  
 Project Number: 03347-143-001-3439-01  
 RFW Contact: John J. Fane Phone: 782-321-4200

No: 116103

SHEET NO. 2 OF 3

**Sample Identification**

| REAC # | Sample No. | Sampling Location | Matrix | Date Collected | # of Bottles | Container/Preservative | KNA, RES, AB | VOC | TAL Metals |
|--------|------------|-------------------|--------|----------------|--------------|------------------------|--------------|-----|------------|
| -774   | C16656     | EMR-2             | S      | 4/14/99        | 1            | 4oz/4°C                |              | ✓   | ✓          |
| -775   | A16657     |                   |        |                |              | 8oz/4°C                | ✓            |     |            |
| -776   | B16657     |                   |        |                |              | 40ml/4°C               |              | ✓   |            |
| -777   | C16657     |                   |        |                |              | 4oz/4°C                |              |     |            |
| -778   | A16658     |                   |        |                |              | 8oz/4°C                | ✓            |     |            |
| -779   | B16658     |                   |        |                |              | 40ml/4°C               |              | ✓   |            |
| -780   | C16658     |                   |        |                |              | 4oz/4°C                | ✓            |     |            |
| -781   | A16659     |                   |        |                |              | 8oz/4°C                |              |     |            |
| -782   | B16659     |                   |        |                |              | 40ml/4°C               | ✓            |     |            |
| -783   | C16659     |                   |        |                |              | 4oz/4°C                |              | ✓   |            |
| -784   | A16660     |                   |        |                |              | 8oz/4°C                | ✓            |     |            |
| -785   | B16660     |                   |        |                |              | 40ml/4°C               |              | ✓   |            |
| -786   | C16660     |                   |        |                |              | 4oz/4°C                |              |     | ✓          |
| -787   | A16661     |                   |        |                |              | 8oz/4°C                | ✓            |     |            |
| -788   | B16661     |                   |        |                |              | 40ml/4°C               |              | ✓   |            |
| -789   | C16661     |                   |        |                |              | 4oz/4°C                |              |     | ✓          |
| -790   | A16662     |                   |        |                |              | 8oz/4°C                | ✓            |     |            |
| -791   | B16662     |                   |        |                |              | 40ml/4°C               |              | ✓   |            |
| -792   | C16662     |                   |        |                |              | 4oz/4°C                |              |     | ✓          |
| -793   | A16663     |                   |        |                |              | 8oz/4°C                | ✓            |     |            |

Special Instructions:

Let to select MS/MSO  
 BNA - Bear new AmSpec on table  
 PCST - pesticides  
 PCST - Polychlorinated biphenyls  
 VOC - volatile organic compounds

**FOR SUBCONTRACTING USE ONLY**  
**FROM CHAIN OF CUSTODY #**

| Items/Reason | Relinquished By | Date    | Received By     | Date    | Time     |
|--------------|-----------------|---------|-----------------|---------|----------|
| all samples  | John J. Fane    | 4/15/99 | David Augustyni | 4/15/99 | 11:27    |
|              |                 |         | David Augustyni | 4/15/99 | 11:40 AM |

oz - ounces  
 ml - milliliters  
 oc - degrees celcius

REAC, Edison, NJ  
 (908) 321-4200  
 EPA Contract 68-C4-0022

**CHAIN OF CUSTODY RECORD**

Project Name: Loft Home Foundry Site  
 Project Number: 03347-143-001-3439-01  
 RFW Contact: Jim J. Fraz Phone: 732-321-4200

No: 06910

SHEET NO 3 OF 3

| Sample Identification |            |                   |        | Analyses Requested |              |                        |              |
|-----------------------|------------|-------------------|--------|--------------------|--------------|------------------------|--------------|
| REAC #                | Sample No. | Sampling Location | Matrix | Date Collected     | # of Bottles | Container/Preservative | ANALYSIS REQ |
| -794                  | B16663     | 13                | S      | 4/14/99            | 1            | 40ml/4°C               | VOC ✓        |
| -795                  | C16663     | 13                | S      |                    |              | 40ml/4°C               |              |
| -796                  | A16664     | 14                | S      |                    |              | 80ml/4°C               |              |
| -797                  | B16664     | 14                | S      |                    |              | 40ml/4°C               |              |
| -798                  | C16664     | 14                | S      |                    |              | 40ml/4°C               |              |
| -799                  | A16665     | 15                | S      |                    |              | 80ml/4°C               |              |
| -800                  | B16665     | 15                | S      |                    |              | 40ml/4°C               |              |
| -801                  | C16665     | 15                | S      |                    |              | 40ml/4°C               |              |

**Matrix:**  
 SD - Sediment  
 DS - Drum Solids  
 DL - Drum Liquids  
 X - Other

PW - Potable Water  
 GW - Groundwater  
 SW - Surface Water  
 SL - Sludge

VOC - volatile organic compounds

Soil  
 Water  
 Oil  
 Air

Special Instructions: id to select MS/MSD  
BNA - Benzene/Acid extractable  
PEST - pesticides  
PCB - polychlorinated biphenyls  
ml - milliliters  
02 - ounces

**FOR SUBCONTRACTING USE ONLY**  
**FROM CHAIN OF CUSTODY #**

| Items/Reason         | Relinquished By       | Date           | Received By           | Date           | Time           |
|----------------------|-----------------------|----------------|-----------------------|----------------|----------------|
| <u>cell analysis</u> | <u>David Provenyi</u> | <u>4/15/99</u> | <u>Y. Ex</u>          | <u>4/15/99</u> | <u>11:25</u>   |
|                      | <u>Jim Fraz</u>       | <u>4/14/99</u> | <u>David Provenyi</u> | <u>4/15/99</u> | <u>10:45am</u> |
|                      |                       |                | <u>David Provenyi</u> | <u>4/15/99</u> | <u>11:40am</u> |

REAC, Edison, NJ  
 (908) 321-4200  
 EPA Contract 68-C4-0022

**CHAIN OF CUSTODY RECORD**

Project Name 1-14 x ALUMINUM FOUNDRY  
 Project Number 03347-143-001-347-01  
 RFW Contact Substare Furna Phone: 732 311 4200

No: 06911

SHEET NO. 1 OF 3

**Sample Identification**

| REAC # | Sample No. | Sampling Location | Matrix | Date Collected | # of Bottles | Container/Preservative | BNA/PEST/ITS | VOCs | Metals |
|--------|------------|-------------------|--------|----------------|--------------|------------------------|--------------|------|--------|
| -802   | A166666    | Location 16       | S      | 14-APR-99      | 1            | 8oz glass / Ice        | ✓            | ✓    |        |
| -803   | B166666    | ↓                 |        |                | 1            | 40ml glass / Ice       |              |      |        |
| -804   | C166666    | ↓                 |        |                | 1            | 4oz glass / Ice        |              |      |        |
| -805   | A166667    | Location 17       |        |                | 1            | 8oz / Ice              | ✓            |      | ✓      |
| -806   | B166667    | ↓                 |        |                | 1            | 40ml / Ice             |              |      |        |
| -807   | C166667    | ↓                 |        |                | 1            | 40ml / Ice             |              |      |        |
| -808   | A166668    | Location 18       |        |                | 1            | 8oz / Ice              | ✓            |      | ✓      |
| -809   | B166668    | ↓                 |        |                | 1            | 40ml / Ice             |              |      |        |
| -810   | C166668    | ↓                 |        |                | 1            | 4oz / Ice              |              |      | ✓      |
| -811   | A166669    | Location 19       |        |                | 1            | 8oz / Ice              | ✓            |      | ✓      |
| -812   | B166669    | ↓                 |        |                | 1            | 40ml / Ice             |              |      |        |
| -813   | C166669    | ↓                 |        |                | 1            | 4oz / Ice              |              |      | ✓      |
| -814   | A166670    | Location 20       |        |                | 1            | 8oz / Ice              | ✓            |      |        |
| -815   | B166670    | ↓                 |        |                | 1            | 40ml / Ice             |              |      | ✓      |
| -816   | C166670    | ↓                 |        |                | 1            | 4oz / Ice              |              |      |        |
| -817   | A166671    | Location 21       |        |                | 1            | 8oz / Ice              | ✓            |      | ✓      |
| -818   | B166671    | ↓                 |        |                | 1            | 40ml / Ice             |              |      |        |
| -819   | C166671    | ↓                 |        |                | 1            | 4oz / Ice              |              |      | ✓      |
| -820   | A166672    | Location 22       |        |                | 1            | 8oz / Ice              | ✓            |      |        |
| -821   | B166672    | ↓                 |        |                | 1            | 40ml / Ice             |              |      | ✓      |

**Analyses Requested**

Special Instructions: LAB to choose MS/MSD

Matrix:  
 SD - Sediment  
 DS - Drum Solids  
 DL - Drum Liquids  
 X - Other

PW - Potable Water  
 GW - Groundwater  
 SW - Surface Water  
 SL - Sludge

S - Soil  
 W - Water  
 O - Oil  
 A - Air

02-ounce  
 ml - milliliters  
 BNA - Benzene not analyzed extractable  
 PEST - pesticides  
 PCB - polychlorinated biphenyls  
 VOC - volatile organic compounds

**FOR SUBCONTRACTING USE ONLY**  
**FROM CHAIN OF CUSTODY #**

| Items/Reason | Relinquished By | Date    | Received By    | Date    | Time    |
|--------------|-----------------|---------|----------------|---------|---------|
| all analysis | Substare        | 4/14/99 | David Angewsky | 4/15/99 | 11:00am |
|              |                 |         | David Angewsky | 4/15/99 | 11:40am |

REAC, Edison, NJ  
(908) 321-4200  
EPA Contract 68-C4-0022

**CHAIN OF CUSTODY RECORD**

Project Name: LAKAWANA FOUNDRY  
Project Number: 03347 143 001 3439 01  
RFW Contact: SALVATORE FAMA Phone: 732 301 4200

No: 06912

SHEET NO. OF 3

041519-

**Sample Identification**

| REAC # | Sample No. | Sampling Location | Matrix | Date Collected | # of Bottles | Container/Preservative | BVA | RES | RB | VOLS | METALS |
|--------|------------|-------------------|--------|----------------|--------------|------------------------|-----|-----|----|------|--------|
| -822   | C16672     | LOCATION 22       | S      | 14-APR-99      | 1            | 4oz glass / ICE        |     |     |    |      |        |
| -823   | A16673     | LOCATION 23       |        |                | 1            | 8oz glass /            |     |     |    |      |        |
| -824   | B16673     |                   |        |                | 1            | 40 ml glass /          |     |     |    | ✓    |        |
| -825   | C16673     |                   |        |                | 1            | 4 oz glass /           |     |     |    |      |        |
| -826   | A16674     | LOCATION 24       |        |                | 1            | 8 oz /                 |     |     |    |      |        |
| -827   | B16674     |                   |        |                | 1            | 40 ml /                |     |     |    | ✓    |        |
| -828   | C16674     |                   |        |                | 1            | 4 oz /                 |     |     |    |      |        |
| -829   | A16675     | LOCATION 25       |        |                | 1            | 8 oz /                 |     |     |    |      |        |
| -830   | B16675     |                   |        |                | 1            | 40 ml /                |     |     |    |      |        |
| -831   | C16675     |                   |        |                | 1            | 4 oz /                 |     |     |    |      |        |
| -832   | A16676     | LOCATION 25-D     |        |                | 1            | 8 oz /                 |     |     |    |      |        |
| -833   | B16676     |                   |        |                | 1            | 40 ml /                |     |     |    |      |        |
| -834   | C16676     |                   |        |                | 1            | 4 oz /                 |     |     |    |      |        |
| -835   | A16677     | LOCATION 26       |        |                | 1            | 8 oz /                 |     |     |    |      |        |
| -836   | B16677     |                   |        |                | 1            | 40 ml /                |     |     |    |      |        |
| -837   | C16677     |                   |        |                | 1            | 4 oz /                 |     |     |    |      |        |
| -838   | A16678     | LOCATION 26-D     |        |                | 1            | 8 oz /                 |     |     |    |      |        |
| -839   | B16678     |                   |        |                | 1            | 40 ml /                |     |     |    |      |        |
| -840   | C16678     |                   |        |                | 1            | 4 oz /                 |     |     |    |      |        |
| -841   | A16679     | LOCATION 27       |        |                | 1            | 8 oz                   |     |     |    |      |        |

**Analyses Requested**

Special Instructions: Lab to select MS/MSD

Soil: S.  
Water: W.  
Oil: O.  
Air: A.

PW - Potable Water  
GW - Groundwater  
SW - Surface Water  
SL - Sludge

Matrix:  
SD - Sediment  
DS - Drum Solids  
DL - Drum Liquids  
X - Other

**FOR SUBCONTRACTING USE ONLY**  
**FROM CHAIN OF CUSTODY #**

| Items/Reason  | Relinquished By | Date    | Received By            | Date    | Relinquished By        | Date    | Received By      | Date    | Time  |
|---------------|-----------------|---------|------------------------|---------|------------------------|---------|------------------|---------|-------|
| cell/analysis | <u>S. Fama</u>  | 4/14/99 | <u>David Anagnosty</u> | 4/15/99 | <u>David Anagnosty</u> | 4/15/99 | <u>Y. Chen</u>   | 4/15/99 | 11:25 |
|               |                 |         |                        |         | <u>David Anagnosty</u> | 4/15/99 | <u>M. Durkin</u> | 4/15/99 | 11:40 |

REFERENCE NO. 7

October 14, 1999

Mr. Kevin Matheis  
U.S. Environmental Protection Agency  
Response and Prevention Branch  
2890 Woodbridge Avenue  
Edison, NJ 08837

**EPA CONTRACT NO: 68-W5-0019**

**TDD NO: 02-99-09-0025**

**DOCUMENT CONTROL NO: START-02-F-**

**SUBJECT: QUALITY ASSURANCE PROJECT PLAN - LACKAWANNA FOUNDRY SITE**

Dear Mr. Matheis:

Enclosed please find the site Quality Assurance Project Plan (QAPP) for the Lackawanna Foundry Site located in Lackawanna, New York.

If you have any questions, please do not hesitate to call me at (908) 225-6116.

Very truly yours,

ROY F. WESTON, INC.

James Kearns  
Project Manager

Enclosure

cc: TDD File  
Joseph M. Soroka, QAO



**SITE QUALITY ASSURANCE PROJECT PLAN**

**LACKAWANNA FOUNDRY SITE  
Lackawanna, Erie County, New York**

Prepared by

Superfund Technical Assessment and Response Team  
Roy F. Weston, Inc.  
Federal Programs Division  
Edison, New Jersey 08837

Prepared for

U.S. Environmental Protection Agency  
Region II - Response and Prevention Branch  
Edison, New Jersey 08837

DCN #: START-02-F-  
TDD #: 02-99-09-0025  
EPA Contract No.: 68-W5-0019

**Approved by:**

START

\_\_\_\_\_  
James Kearns  
Project Manager

Date: \_\_\_\_\_

START

\_\_\_\_\_  
Joseph M. Soroka, START  
Quality Assurance Officer

Date: \_\_\_\_\_

EPA

\_\_\_\_\_  
Kevin Matheis  
Task Monitor

Date: \_\_\_\_\_

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The following elements are provided in the START Generic Quality Assurance Project Plan (QAPP) and are included by reference:

- QA REPORTS TO MANAGEMENT
- PREVENTIVE MAINTENANCE PROCEDURES AND SCHEDULES
- RECORDS MANAGEMENT SYSTEM
- LOGBOOK PROGRAM
- QUALITY-RELATED DOCUMENTS
- INSPECTION/ACCEPTANCE REQUIREMENTS FOR SUPPLIES AND CONSUMABLES

**LIST OF ATTACHMENTS**

ATTACHMENT A: Site Map

ATTACHMENT B: EPA/ERT Soil Sampling SOP #2012 and EPA/ERT Sampling  
Equipment Decontamination SOP#2006

## **1.0 INTRODUCTION**

Presented herein is the site Quality Assurance Project Plan (QAPP) for the sampling event to be conducted at the Lackawanna Foundry Site by the Region II Superfund Technical Assessment and Response Team (START). The site QAPP has been developed at the request of the United States Environmental Protection Agency (U.S. EPA) in accordance with the START Generic Quality Assurance Project Plan (QAPP).

This plan is based on information currently available and may be modified on site in light of field screening results and other acquired information. All deviations from the QAPP will be noted in the Sampling Trip Report.

## **2.0 PROJECT DESCRIPTION**

The Lackawanna Foundry Removal Action Site is located at 2 Elm Street, Lackawanna, Erie County, New York. The Site is located in a mixed residential and heavily industrialized area adjacent to Smokes Creek, a direct tributary to Lake Erie. The residential neighborhood immediately adjacent to the Site consists of twelve residences with over 100 residences being located within 1/8 mile to site. The population of the area within 1/2 mile of the Site consists of nearly 1,000 people.

The Site has been the location of a cast-iron foundry since the early 1900's. The current owner of the Site inherited the business from his father whom operated the foundry at this location for more than 30 years. Over the past five years, the current owner has operated a marginal business and has ceased operations since approximately 1997.

On December 22, 1998, the New York State Department of Environmental Conservation (NYSDEC) requested that U.S. Environmental Protection Agency (EPA) evaluate the site and perform a removal action, as appropriate, to address the threats at the Site from drums, PCB transformers, and PCB contaminated soil. Included in the referral was a drum inventory performed by the site owner that showed the presence of over 223 drums of various foundry raw materials and waste.

On March 19, 1999, the Acting Division Director of the Emergency and Remedial Response Division granted verbal authorization for the EPA to undertake emergency stabilization activities at the site to include fencing the perimeter of the Site, securing the buildings, moving drums into a secure location and stabilizing leaking containers. This stabilization action was initiated on March 22, 1999.

Two primary areas of contamination have been identified on Site to this date. One area being the On-Site Buildings which have PCB contamination, Laboratory chemicals, compressed gas cylinders, paint cans, and a couple of hundred drums.

The second area of contamination, the area of concern for this sampling event, is a recreational area owned by the City of Lackawanna and is located adjacent to the Site. The recreational area is a grassy area that Lackawanna maintains by frequent grass and weed cuttings. The arsenic contamination adjacent to the Site in the recreational area appears to be in the location of a bag-house dust collection system used during foundry operations. The contamination was caused by either airborne deposition or run-off from the baghouse collection area. Analytical data from an ERT mobilization in April of 1999 showed levels of arsenic contamination in this second area of contamination were as high as 77 parts per million (ppm).

### **3.0 PROJECT ORGANIZATION AND RESPONSIBILITIES**

The EPA Task Monitor (TM), Kevin Matheis, will provide overall direction to the staff concerning project sampling needs, objectives, and schedule. The Project Manager (PM), Jim Kearns, will be the primary point of contact with the TM. The PM is responsible for the development and completion of the Sampling QA/QC Plan, project team organization, and supervision of all project tasks, including reporting and deliverables. The Site QC Coordinator will be responsible for ensuring field adherence to the Sampling QA/QC Plan and recording of any deviations. The START Analytical Services Coordinator, Smita Sumbaly, will be the primary project team site contact with the subcontracted laboratory, if necessary.

START will arrange for the laboratory analyses. START personnel will transfer custody of the soil samples for shipment to the appropriate laboratory. The raw analytical data from the laboratory will be provided to the START Analytical Services Group for data validation.

The following sampling personnel will work on this project:

#### **Personnel**

James Kearns

#### **Responsibility**

Project Manager, Field Coordinator  
Sampling, Shipping  
Site QA/QC

The following laboratories will provide the following analyses:

| <u>Lab Name/Location</u> | <u>Sample Type</u> | <u>Parameters</u> |
|--------------------------|--------------------|-------------------|
| To be determined         | Soil               | Arsenic           |

A standard two week verbal and three week written results turnaround time has been requested by the TM.

#### **4.0 DATA USE OBJECTIVES**

The objective of this sampling event is to confirm and delineate the extent of arsenic contamination near the Lackawanna Foundry building. This sampling event will help in the remediation efforts at the site by helping delineate the extent of soil contamination. The data from this sampling event will be reviewed using the Site cleanup levels for the Lackawanna Foundry site of 20 ppm for arsenic.

#### **4.1 Quality Assurance Objectives**

The overall Quality Assurance (QA) objective for chemical measurement data associated with this sampling event is to provide analytical results that are legally defensible in a court of law. The QA program will incorporate Quality Control (QC) procedures for field sampling, chain of custody, laboratory analyses, and reporting to assure generation of sound analytical results.

The EPA Task Monitor (TM) has specified a critical level of QA of QA-2 for the Site. Details of this QA level are provided below.

The QA Protocols for a Level 2 QA objective sampling event are applicable to all sample matrices and include:

1. Sample documentation in the form of field logbooks, appropriate field data sheets, and chain of custody records (chain of custody records are optional for field screening locations).
2. Calibration of all monitoring and/or field-portable analytical equipment prior to collection and analyses of samples with results and/or performance check procedures/methods summarized and documented in a field, personal, and/or instrument

log notebook.

3. Field or laboratory determined method detection limits (MDLs) will be recorded along with corresponding analytical sample results, where appropriate.
4. Analytical holding times as determined from the time of sample collection through analysis. These will be documented in the field logbook or by the laboratory in the final data deliverable package.
5. Initial and continuous instrument calibration data.
6. QC blank results (rinsate, trip, method, preparation, instrument, etc.), as applicable.
7. Collection and analysis of blind field duplicate and MS/MSD QC samples to provide a quantitative measure of the analytical precision and accuracy, as applicable.
8. Use of the following QC procedure for QC analyses and data validation:

Definitive identification - confirm the identification of analytes on 100% of the "critical" samples, via an EPA-approved method; provide documentation such as gas chromatograms, mass spectra, etc.

The objective of this project/event applies to the following parameters:



**Table 1: Quality Assurance Objectives**

| <b>QA Parameters</b> | <b>Matrix</b> | <b>Intended Use of Data</b>            | <b>QA Objective</b> |
|----------------------|---------------|--|---------------------|
| TAL metals           | Soil          | Delineate extent of soil contamination | QA-2                |

A Field Sampling Summary is attached in Table 2 and a QA/QC Analysis and Objectives Summary is attached in Table 3. Section 5.1, Sampling Design, provides information on analyses to be performed on the individual soil samples.

**TABLE 2:**

**FIELD SAMPLING SUMMARY**

| lytical<br>meters | Matrix  | Container Size                       | Preservative                               | Holding Time <sup>1</sup>   | Subtotal<br>Samples | Rinsate<br>Blanks <sup>2</sup> | Duplicate<br>Samples <sup>3</sup> | MS/MSD<br>Samples <sup>3</sup> | Total<br>Field<br>Samples |
|-------------------|---------|--------------------------------------|--|-----------------------------|---------------------|--------------------------------|-----------------------------------|--------------------------------|---------------------------|
| ls-(arsenic)      | Soil    | 1 - 8 oz glass                       | Cool to 4°C                                | 6 months, 28 days for<br>Hg |                     |                                |                                   |                                |                           |
| ls-(arsenic)      | Aqueous | 1 - 1 Liter glass or<br>polyethylene | Cool to 4°C,<br>HNO <sub>3</sub> to pH < 2 | 6 months, 28 days for<br>Hg | 2                   | 2                              |                                   |                                | 2                         |

<sup>1</sup> Holding time from date of sampling.

<sup>2</sup> Only required if non-dedicated sampling equipment to be used. NR - not required, dedicated sampling equipment to be used.

<sup>3</sup> Not required for QA-1 (screening)

Please refer to Attachment B for a list of EPA/ERT SOPs to be used:

**TABLE 3**

**QA/QC Analysis and Objectives Summary**

| Analytical Parameters | Matrix | Analytical Method Reference       | QA/QC Quantitation Limits | QA<br>Objective |
|-----------------------|--------|-----------------------------------|---------------------------|-----------------|
| etals-(arsenic)       | Soil   | SW846 Method 5000 or CLP ILMO 4.0 | As per method             | QA-2            |

| Analytical Parameters | Matrix  | Analytical Method Reference       | QA/QC Quantitation Limits | QA Objective |
|-----------------------|---------|-----------------------------------|---------------------------|--------------|
| etals-(arsenic)       | Aqueous | SW846 Method 5000 or CLP ILMO 4.0 | As per method             | QA-2         |

Note: CLP-format deliverables required for all data packages.

## 5.0 APPROACH AND SAMPLING PROCEDURES

The following sampling activities will be conducted at the Lackawanna Foundry Site:

- Soil Sampling of area adjacent to recreational facility located near the Foundry building.

### 5.1 Sampling Design

The field program will include the collection of soil samples. All sampling activities will be performed by the Region II START, under the direction of the EPA Task Monitor (TM). All samples will be sent to the procured laboratory.

A maximum of 16 surface soil samples and 16 depth soil samples will be collected with dedicated scoops and analyzed for PCBs. Prior to the sample collection, gravel, debris, or foliage will be removed from the sampling point. Surface soil samples will be collected from 0-6 inches in depth. The depth samples will be collected from 18-24 inches in depth. A grid will be set up ahead of time so that all samples will be spaced out equally over the entire sampling area. Depth samples will be collected from the same locations as the surface samples and will be collected by first using a power auger to reach the proper depth then a decontaminated hand auger will be used to retrieve the sample.

### 5.2 Schedule of Activities

| <b>Proposed Start Date</b> | <b>Activity</b> | <b>End Date</b>  |
|----------------------------|-----------------|------------------|
| October 18, 1999           | Soil Sampling   | October 20, 1999 |

### 5.3 Sampling Equipment

In order to avoid cross-contamination, all samples will be collected with dedicated, disposable sampling equipment or decontaminated stainless steel scoops.

For subsurface samples, decontaminated stainless steel augers and trowels will be used.

#### **5.4 Sample Identification System**

Each sample collected by Region II START will be designated by a code which will identify the site. The code will be a site-specific project tracking number. The code for the Recreational Area of the Lackawanna Foundry Site is "RA". The media type will follow the numeric code. A hyphen will separate the site code and media type. Specific media types are as follows:

- S - Surface Soil
- SS - Subsurface Soil
- RIN - Rinsate Blank

After the media type, the sequential sample numbers will be listed; sample numbers will be identified as to their location area on the site location and/or the location on the x and y coordinates of the sampling grid. A duplicate sample will be identified in the same manner as other samples and will be distinguished and documented in the field logbook.

#### **5.5 Standard Operating Procedures (SOPs)**

##### **5.5.1 Sample Documentation**

All sample documents will be completed legibly, in ink. Any corrections or revisions will be made by lining through the incorrect entry and by initialing the error.

#### **FIELD LOGBOOK**

The field logbook is essentially a descriptive notebook detailing site activities and observations so that an accurate account of field procedures can be reconstructed in the writer's absence. All entries will be dated and signed by the individuals making the entries, and should include (at a minimum) the following:

1. Site name and project number.
2. Name(s) of personnel on site.
3. Dates and times of all entries (military time preferred).
4. Descriptions of all site activities, site entry and exit times.
5. Noteworthy events and discussions.
6. Weather conditions.

7. Site observations.
8. Sample and sample location identification and description\*.
9. Subcontractor information and names of on-site personnel.
10. Date and time of sample collections, along with chain of custody information.
11. Record of photographs.
12. Site sketches.

The description of the sample location will be noted in such a manner as to allow the reader to reproduce the location in the field at a later date.

### SAMPLE LABELS

Sample labels will clearly identify the particular sample, and should include the following:

1. Site/project number.
2. Sample identification number.
3. Sample collection date and time.
4. Designation of sample (grab or composite).
5. Sample preservation.
6. Analytical parameters.
7. Name of sampler.

Sample labels will be written in indelible ink and securely affixed to the sample container. Tie-on labels can be used if properly secured.

### CUSTODY SEALS

Custody seals demonstrate that a sample container has not been tampered with, or opened. The individual in possession of the sample(s) will sign and date the seal, affixing it in such a manner that the container cannot be opened without breaking the seal. The name of this individual, along with a description of the sample packaging, will be noted in the field logbook.

#### **5.5.2 Sampling SOPs**

The following Sampling SOPs will be used for this project:

- Soil Sampling EPA/ERT SOP# 2012
- Sampling Equipment Decontamination EPA/ERT SOP# 2006

Please refer to Attachment B for a list of EPA/ERT sampling SOPs.

### Decontamination

All stainless steel equipment involved in field sampling activities will be decontaminated in accordance to EPA/ERT SOP #2006 prior to and subsequent to sampling. Decontamination of sampling equipment will be kept to a minimum in the field; whenever possible dedicated sampling equipment, will be used. Decontamination of sampling equipment, including stainless steel augers and bowls, will be conducted as follows:

- 1) Alconox detergent and potable water scrub,
- 2) Potable water rinse,
- 3) A 10% nitric acid rinse (ultra pure grade) when sampling for inorganic parameters,
- 4) Distilled or potable water rinse,
- 5) Deionized water rinse,
- 6) Air dry (sufficient time will be allowed for the equipment to completely dry), and
- 7) Wrap or cover exposed ends of sampling equipment with aluminum foil (shiny side out) for transport and handling.

## **SOIL SAMPLING**

Soil sampling activities will be conducted in accordance with guidelines outlined in EPA/ERT Soil Sampling SOP #2012 (Attachment B).

### **5.5.3 Sample Handling and Shipment**

Each of the sample bottles will be sealed and labeled according to the following protocol. Caps will be secured with custody seals. Bottle labels will contain all required information including site/project code and sample number, time and date of collection, analyses requested, and preservative used. Sealed bottles will be placed in large metal or plastic coolers, and padded with an absorbent material such as vermiculite. All packaging will conform to IATA Transportation regulations for overnight carriers.

All sample documents will be sealed in a plastic bag and affixed to the underside of each cooler lid. The lid will be sealed and affixed on at least two sides with custody seals so that any sign of tampering is easily visible.

### **5.6 Sample Containers**

All sample containers will meet the QA/QC specifications in OSWER Directive 9240.0-05A, "Specifications and Guidance for Contaminant Free Sample Containers".

### **5.7 Disposal of PPE and contaminated sampling materials**

All used PPE and disposable sampling equipment will be left on site for disposal during the on-going removal action.

## **6.0 CHAIN OF CUSTODY RECORD**

A chain of custody record will be maintained from the time the sample is taken to its final deposition. Every transfer of custody must be noted and signed for, and a copy of this record kept by each individual who has signed. When samples (or groups of samples) are not under direct control of the individual responsible for them, they must be stored in a locked container sealed with a custody seal. Specific information regarding custody of the samples projected to be collected on the weekend will be noted in the field logbook.

The chain of custody record should include (at minimum) the following:

1. Sample identification number.
2. Sample information.
3. Sample location.
4. Sample date.
5. Name(s) and signature(s) of sampler(s).
6. Signature(s) of any individual(s) with control over samples.

A separate chain of custody form must accompany each cooler for each daily shipment. The chain of custody form must address all samples in that cooler, but not address samples in any other cooler. This practice maintains the chain of custody for all samples in case of mis-shipment.

## **7.0 FIELD INSTRUMENT CALIBRATION AND PREVENTIVE MAINTENANCE**

The sampling team is responsible for assuring that a calibration/maintenance log will be brought into the field and maintained for each measuring device. Each log will include at a minimum, where applicable:

- name of device and/or instrument calibrated
- device/instrument serial and/or ID number
- frequency of calibration
- date of calibration
- results of calibration
- name of person performing the calibration
- identification of the calibrant (PID, FID, pH meter)

Equipment to be used each day will be calibrated prior to the commencement of daily



activities.

## **8.0 ANALYTICAL METHODS**

Analytical methods to be utilized in the analyses of samples collected during this sampling event are detailed in Table 3.

## **9.0 DATA REDUCTION, VALIDATION AND REPORTING**

### **9.1 Deliverables**

The START PM, James Kearns, will maintain contact with the EPA TM, Kevin Matheis, to keep him informed about the technical and financial progress of this project. This communication will commence with the issuance of the work assignment and project scoping meeting. Activities under this project will be reported in status and trip reports and other deliverables (e.g., analytical reports, final reports) described herein. Activities will also be summarized in appropriate format for inclusion in monthly and annual reports.

The following deliverables will be provided under this project:

#### **TRIP REPORT**

A trip report will be prepared to provide a detailed accounting of what occurred during each sampling mobilization. The trip report will be prepared within one week of the last day of each sampling mobilization. Information will be provided on time of major events, dates, and personnel on site (including affiliations).

#### **MAPS/FIGURES**

Maps depicting site layout, contaminant source areas, and sample locations will be included in the trip report, as appropriate.

#### **ANALYTICAL REPORT**

An analytical report will be prepared for samples analyzed under this plan. Information regarding the analytical methods or procedures employed, sample results, QA/QC results, chain of custody documentation, laboratory correspondence, and raw data will be provided within this deliverable.

#### **DATA REVIEW**

A review of the data generated under this plan will be undertaken. The assessment of data acceptability or useability will be provided separately, or as part of the analytical report.

## **9.2 Data Validation**

Data generated under this QA/QC Sampling Plan will be evaluated according to criteria contained in the Removal Program Data Validation Procedures that accompany OSWER Directive number 9360.4-1 and in accordance with Region 2 guidelines.

Laboratory analytical results will be assessed by the data reviewer for compliance with required precision, accuracy, completeness, representativeness, and sensitivity.

## **10.0 FIELD QA/QC CHECK SAMPLES AND FREQUENCY**

This section details the Quality Assurance/Quality Control (QA/QC) requirements for field activities performed during the sampling effort.

QA/QC samples will include the collection of one field duplicate and one matrix spike/matrix spike duplicate sample for each matrix (soil/sediment) at a ratio of 1 per 20 samples. Extra sample volume will be submitted to allow the laboratory to perform matrix spike sample analysis. This analysis provides information about the effect of sample matrix on digestion and measurement methodology. Field duplicate samples provide an indication of analytical variability and analytical error and will not be identified to the laboratory.

Field Rinsate Blanks will be collected when non-dedicated sampling equipment is used. A field rinsate blank will consist of distilled deionized (DI), demonstrated analyte-free water that has been poured over decontaminated sampling equipment. The field rinsate blank analytical results will be utilized in evaluation of potential cross contamination resulting from inadequate decontamination only if non-dedicated sampling equipment is used. The frequency of field rinsate blank collection is one blank per decontamination event per type of equipment, not to exceed more than one per day. Blanks will be collected for all parameters of interest (excluding physical parameters) and shipped with the samples collected the same day. Field rinsate blanks will be collected by Region II START.

Field rinsate blanks will be collected in accordance with the procedure listed below:

- 1) Decontaminate sampling equipment using the procedure specified in Section 3.3 of this plan.
- 2) Pour DI water over the sampling device and collect the rinsate in the appropriate sample containers.

## **11.0 SYSTEM AUDIT**

The Field QA/QC Officer will observe sampling operations and review subsequent analytical results to ensure compliance with the QA/QC requirements of the project/sampling event.

## **12.0 CORRECTIVE ACTION**

All provisions will be taken in the field and laboratory to ensure that any problems that may develop will be dealt with as quickly as possible to ensure the continuity of the project/sampling events. Any deviations from this sampling plan will be noted in the final report.

**ATTACHMENT A**

**SITE MAPS**

**ATTACHMENT B**  
**EPA/ERT SAMPLING SOPS**

REFERENCE NO. 8

November 5, 1999

Mr. Kevin Matheis  
U.S. Environmental Protection Agency  
Removal Action Branch  
2890 Woodbridge Avenue  
Edison, NJ 08837

**EPA CONTRACT NO: 68-W5-0019**  
**TDD NO: 02-99-09-0025**  
**DOCUMENT CONTROL NO: START-02-F-03897**  
**SUBJECT: SAMPLING TRIP REPORT- LACKAWANNA FOUNDRY**

Dear Mr. Matheis:

Enclosed please find the Sampling Trip Report for the Lackawanna Foundry Site located at 2 Elm Street, Lackawanna, Erie County, New York. This Trip Report pertains to the soil sampling event for that occurred on October 18-20, 1999.

If you have any questions, do not hesitate to call me at (732) 225-6116.

Very truly yours,

ROY F. WESTON, INC.

James Kearns  
Project Manager

Enclosure

cc: TDD File  
Joseph M. Soroka

## SAMPLING TRIP REPORT

**SITE NAME:** Lackawanna Foundry Site

DCN #: START-02-F-03897

TDD #: 02-99-09-0025

### SAMPLING DATES:

1. Site Location: 2 Elm Street, Lackawanna, Erie County, New York
2. Sample Locations: Rear of Lackawanna Foundry building adjacent to recreational area owned by the City of Lackawanna (See Attachment A).
3. Sample Descriptions: Samples were taken from a grid that was set up to run along the south side of the Lackawanna Foundry building (the site) encompassing the area between the foundry building and the recreational area located south of the site. The recreational area which includes a baseball field and an indoor bocce ball court is owned by the City of Lackawanna. Samples were taken at 20 ft. northerly intervals and 35 ft. easterly intervals. All sample identification numbers reflect their distance from point N0-E0, the origin. Therefore, if a sample identification number was N20-E70, this meant that the location was 20 ft. north and 70 ft. east of the origin. The origin (N0-E0) was located 70 ft. from the Lackawanna Foundry building and 31 ft. from a utility pole located at the west end of the sampling grid.

4. Laboratories Receiving Samples:

| <u>Sample Type</u> | <u>Name and Address of Laboratory</u>                                  |
|--------------------|--|
| Soil               | STL Inc.<br>Suite 106<br>10 Hazelwood Drive<br>Amherst, New York 14228 |

5. Sample Dispatch Data:

On October 20, 1999, 35 samples (including 2 duplicate samples and 1 equipment blank) were hand delivered to the STL Inc. laboratory for arsenic analysis.

6. Sampling Personnel

| <u>Name</u>    | <u>Company</u>                  | <u>Duties on Site</u>     |
|----------------|---------------------------------|---------------------------|
| James Kearns   | Region II START                 | Project Manager           |
| Patrick Austin | Region II START                 | Sample Management Officer |
| Alex Steele    | Roy F. Weston- Corporate Office | Sampling                  |
| John Leeks     | Roy F. Weston- Corporate Office | Sampling                  |



7. Additional Comments:

On October 19, 1999, START performed soil sampling on the south side of the Lackawanna Foundry building near the recreational area owned by the City of Lackawanna which includes a baseball field and an indoor bocce ball court. Sampling included both surface and subsurface samples that were analyzed for arsenic contamination. Eleven samples were collected at various depths and 23 samples were taken at the surface. The original sampling plan called for 16 surface and 16 subsurface samples but refusal was encountered at several locations. The refusal area appeared to be a cement pad with slag beneath the pad. A gas powered two man auger was used in areas where the manual hand augers hit refusal. Even after using the gas powered auger there were still several locations that encountered refusal. Therefore, after discussing the situation with the EPA Task Monitor, the decision was made to collect the depth samples if a depth of one foot or greater could be reached before encountering refusal and if this depth could not be reached then extending the grid and increasing the amount of surface samples would occur. The sampling grid was extended and several additional surface samples were collected. Please refer to Table 1 and Table 2 for further sample information including information for sampling depths of each sample.

At sample location N20 E70 a hydrocarbon odor was detected between a depth of one to one and a half feet. Sampling was immediately halted and air monitoring with a Photovac Microtip showed readings at approximately 10 parts per million (ppm). Therefore, it was assumed that the soil was contaminated with volatiles. Breathing zone readings on the Microtip showed no readings. At the request of the EPA Task Monitor (TM) a sample of the contaminated soil was collected and given to the ERRS contractor for laboratory analysis.

An error occurred on the Chain of Custody form. Sample S20 E105 was incorrectly labeled as S20 135 and sample S20 E140 was incorrectly labeled as S20 140. The error was amended on the attached Chain of Custody form. Copies of the Chain of Custody records are presented in Attachment B.

Photo documentation is presented in Attachment C.

Table 1: Lackawanna Foundry Soil Sampling for Arsenic- Surface Samples

| <u>Sample #</u> | <u>Sample Date</u> | <u>Sample Time</u> | <u>Additional Information</u>   |
|-----------------|--------------------|--------------------|---|
| *N0 E0          | 10/19/99           | 0900               | Surface sample located at the Origin of the sampling grid (31' east of utility pole located on west end of sampling area and 70' south of Lackawanna Foundry building). |
| N0 E0-1         | 10/19/99           | 0905               | Duplicate sample of (N0 E0).  |
| *N20 E0         | 10/19/99           | 0907               | Surface sample located 20' north of origin.   |
| N20 E0-1        | 10/19/99           | 0908               | Duplicate sample of (N20 E0). Sample located 20' north of origin.   |
| N20 E35         | 10/19/99           | 0910               | Surface sample located 20' north and 35' east of origin.  |
| N0 E35          | 10/19/99           | 0915               | Surface sample located 35' east of origin.  |
| N40 E0          | 10/19/99           | 0915               | Surface sample located 40' north of origin.   |
| N40 E105        | 10/19/99           | 0925               | Surface sample located 40' north and 105' east of origin.   |
| N20 E70         | 10/19/99           | 0922               | Surface sample located 20' north and 70' east of the origin.  |
| N20 E105        | 10/19/99           | 0930               | Surface sample located 20' north and 105' east of the origin.   |
| N40 E35         | 10/19/99           | 0930               | Surface sample located 40' north and 35' east of the origin.  |
| N0 E70          | 10/19/99           | 0930               | Surface sample located 70' east of the origin.  |
| N20 E140        | 10/19/99           | 0938               | Sample located 20' north and 140' east of the origin.   |
| N0 E105         | 10/19/99           | 0940               | Sample located 105' east of the origin.   |
| N0 E140         | 10/19/99           | 0945               | Sample located 140' east of the origin.   |
| N40 E70         | 10/19/99           | 0950               | Sample located 40' north and 70' east of the origin.  |
| N0 E175         | 10/19/99           | 0955               | Sample collected 175' east of the origin.   |
| N20 E175        | 10/19/99           | 0935               | Sample collected 20' north and 175' east of the origin.   |
| N60 E0          | 10/19/99           | 1555               | Sample collected 60' north of origin.   |
| S20 E35         | 10/19/99           | 1205               | Sample collected 20' south and 35' east of the origin.  |
| S20 E70         | 10/19/99           | 1200               | Sample collected 20' south and 70' east of the origin.  |
| S20 E105        | 10/19/99           | 1215               | Sample collected 20' south and 105' east of the origin.   |
| S20 E140        | 10/19/99           | 1210               | Sample collected 20' south and 140' east of the origin.   |

\*- Extra volume for MS/MSD analysis was provided for these samples.

Table 2: Lackawanna Foundry Soil Sampling for Arsenic- Depth Samples

| <u>Sample #</u> | <u>Sample Date</u> | <u>Sample Time</u> | <u>Sample Depth</u> | <u>Additional Information</u>                              |
|-----------------|--------------------|--------------------|---------------------|--|
| N0 E0-D         | 10/19/99           | 11:40              | 12"                 | Sample location was at the origin of the sampling grid.    |
| N20 E0-D        | 10/19/99           | 11:45              | 12"                 | Sample location was 20' north of the origin.               |
| N40 E0-D        | 10/19/99           | 11:50              | 24"                 | Sample location was 40' north of the origin.               |
| N0 E35-D        | 10/19/99           | 12:50              | 12"-18"             | Sample location was 35' east of the origin.                |
| N20 E35-D       | 10/19/99           | 15:00              | 18"                 | Sample location was 20' north and 35' east of the origin.  |
| N40 E35-D       | 10/19/99           | 11:55              | 24"                 | Sample location was 40' north and 35' east of the origin.  |
| N40 E70-D       | 10/19/99           | 13:00              | 18"-24"             | Sample location was 40' north and 70' east of the origin.  |
| N20 E70-D       | 10/19/99           | 13:50              | 15"-18"             | Sample location was 20' north and 70' east of the origin.  |
| N40 E105-D      | 10/19/99           | 15:15              | 21"-24"             | Sample location was 40' north and 105' east of the origin. |
| N0 E140-D       | 10/19/99           | 15:15              | 12"-15"             | Sample location was 140' east of the origin.               |
| N0 E175-D       | 10/19/99           | 15:40              | 12"-18"             | Sample location was 175' east of the origin.               |

8. Report Prepared by: \_\_\_\_\_ Date: \_\_\_\_\_  
James Kearns, START PM
9. Report Approved by: \_\_\_\_\_ Date: \_\_\_\_\_  
Dan Crouse, START QC

**ATTACHMENT A  
SITE LOCATION AND SAMPLING  
LOCATION MAPS**

**ATTACHMENT C  
PHOTO DOCUMENTATION**

ghhg

REFERENCE NO. 9



Roy F. Weston, Inc.  
Federal Programs Division  
Suite 201  
1090 King Georges Post Road  
Edison, New Jersey 08837-3703  
732-225-6116 • Fax 732-225-7037

SUPERFUND TECHNICAL ASSESSMENT AND RESPONSE TEAM  
EPA CONTRACT 68-W5-0019

June 30, 2000

Mr. Kevin Matheis  
U.S. Environmental Protection Agency  
Removal Action Branch  
2890 Woodbridge Avenue  
Edison, NJ 08837

**EPA CONTRACT NO: 68-W5-0019**  
**TDD NO: 02-00-03-0032**  
**DOCUMENT CONTROL NO: START-02-F-04350**  
**SUBJECT: SAMPLING FINAL TRIP REPORT- LACKAWANNA FOUNDRY**

Dear Mr. Matheis:

Enclosed please find the Final PCB Sampling Trip Report for the Lackawanna Foundry Site located at 2 Elm Street, Lackawanna, Erie County, New York. This Trip Report pertains to the soil sampling event for that occurred between May 3-5, 2000.

If you have any questions, do not hesitate to call me at (732) 225-6116.

Very truly yours,

ROY F. WESTON, INC.

A handwritten signature in dark ink, appearing to read "D. Adams", is written over the typed name.

David L. Adams  
Project Manager

Enclosure

cc: TDD File  
Joseph M. Soroka



## SAMPLING TRIP REPORT

**SITE NAME:** Lackawanna Foundry Site  
DCN #: START-02-F-04350  
TDD #: 02-00-03-0032

### SAMPLING DATES:

1. Site Location: 2 Elm Street, Lackawanna, Erie County, New York
2. Sample Locations: Rear of Lackawanna Foundry building adjacent to recreational area owned by the City of Lackawanna (Attachment A).
3. Sample Descriptions: Two samples and a duplicate sample were collected for PCB analysis, on May 3, 2000. Sample LW-SS-03 is a duplicate of sample LW-SS-02. (Please refer to Table 1 for further sample information). Grab samples were collected from surface soil of locations selected by the OSC (Attachment A).

At the request of the EPA Task Monitor (TM) all samples were collected and given to the ERRS contractor for laboratory analysis. Copies of chains of custody can be found in Attachment B.

4. Laboratories Receiving Samples:

| <u>Sample Type</u> | <u>Name and Address of Laboratory</u>   |
|--------------------|---|
| Soil               | Upstate Laboratories<br>6034 Corporate Drive<br>East Syracuse, New York 13057 |

5. Sample Dispatch Data:

On May 3, 2000, 3 samples (including 1 duplicate sample) were hand delivered to Vernon Wilson, Earth Tech Response Manager, for delivery to Upstate Laboratories for PCB analysis.

6. Sampling Personnel

| <u>Name</u>    | <u>Company</u>  | <u>Duties on Site</u>    |
|----------------|-----------------|--------------------------|
| David L. Adams | Region II START | Project Manager/ Sampler |

7. Additional Comments:

On December 20, 1999, Earth Tech collected PCB confirmation samples in the excavation (Attachment A). Analytical data can be found in Attachment F.

On April 11, 2000, START performed soil sampling on the south side of the Lackawanna Foundry building near the recreational area owned by the City of Lackawanna which includes a baseball field and an indoor bocce ball court. Sampling included surface samples that were analyzed for arsenic contamination. Five samples were collected at the surface. START performed a PCB immunoassay test (Strategic Diagnostics Inc.) on suspect PCB contaminated soil, on the concrete pad located in the east end of the property (Attachment A). The screening was used to determine where the PCB samples would be collected. As a result, sample point X11 was used for PCB screening only (Attachment A). Please refer to Table 3 for further sample information.

On May 3, 2000, analytical results for the arsenic and PCB sampling event in April were received, all samples were below the cleanup levels for arsenic (Table 2). PCB results are detailed in Table 3. Analytical data can be found in Attachment C and D.

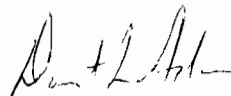
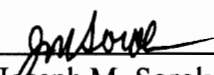
Analytical results for the backfill soil can be found in Attachment E.

Sample N20EO data (Attachment C), Attachment G details additional information obtained by the lab on the sample.

On May 18, 2000, analytical results for the PCB sampling event on May 3 were received, all samples were below the cleanup levels for PCBs (Table 1). Final analytical results can be found in Attachment H.

On May 18, 2000, analytical results for the topsoil were received. A copy of the report and the chain of custody can be found in Attachment I.

Results of the wetlands soil testing were received from Severn Trent Laboratories and can be found in Attachment J.

8. Report Prepared by:  Date: 6/30/00  
David L. Adams, START PM
9. Report Approved by:  Date: 7/3/00  
Joseph M. Soroka, START QCO

**Table 1: Lackawanna Foundry Soil Sampling for PCB**

| <b>Sample#</b> | <b>Sample Date</b> | <b>Sample Time</b> | <b>Analytical Results</b> | <b>Additional Information</b> |
|----------------|--------------------|--------------------|---------------------------|-------------------------------|
| LW-SS-01       | 5/3/00             | 0940               | .134 ppm                  | Brown sand/soil               |
| LW-SS-02       | 5/3/00             | 1000               | 5.8 ppm                   | Tan clay                      |
| LW-SS-03       | 5/3/00             | 1000               | 13 ppm                    | Tan clay                      |

**Table 2: Lackawanna Foundry Soil Sampling for Arsenic Results**

| <b>Sample #</b> | <b>Sample Date</b> | <b>Sample Time</b> | <b>Analytical Results</b> |
|-----------------|--------------------|--------------------|---------------------------|
| N20 E0          | 4/12/00            | 0955               | 14 ppm                    |
| N60 E0          | 4/12/00            | 0959               | 2.5 ppm                   |
| ARConf(3)       | 4/12/00            | 1000               | 0.3 ppm                   |
| ARConf(4)       | 4/12/00            | 1000               | 1.6 ppm                   |
| ARConf(6)<br>*  | 4/12/00            | 1010               | 1.0 ppm                   |

\* Triple volume taken for MS/MSD.

**Table 3: Lackawanna Foundry Soil Sampling for PCBs  
Screening and Analytical Results**

| <b>Sample #</b> | <b>Sample Date</b> | <b>Sample Time</b> | <b>Sample Depth</b> | <b>Immunoassay Screen Results</b>            | <b>Analytical Results</b> |
|-----------------|--------------------|--------------------|---------------------|--|---------------------------|
| X1*             | 4/11/00            | 0955               | 0-3"                | <10ppm on PCB screening                      | 12 ppm                    |
| X4              | 4/11/00            | 1010               | 0-3"                | >10ppm on PCB screening                      | 33 ppm                    |
| X5              | 4/11/00            | 1015               | 0-3"                | >10ppm on PCB screening                      | 62 ppm                    |
| X6              | 4/11/00            | 1020               | 0-3"                | >10ppm on PCB screening                      | 16 ppm                    |
| X7              | 4/11/00            | 1025               | 0-3"                | >10ppm on PCB screening                      | 11 ppm                    |
| X10             | 4/11/00            | 1035               | 0-3"                | >10ppm on PCB screening                      | 9 ppm                     |
| X99             | 4/11/00            | 1035               | 0-3"                | >10ppm on PCB screening,<br>Duplicate of X10 | 7 ppm                     |
| X12             | 4/11/00            | 1415               | 0-3"                | >10ppm on PCB screening                      | 19 ppm                    |

\* Triple volume taken for MS/MSD.

**ATTACHMENT A**

**ATTACHMENT B**

REF NO:  
 PO NO:



**SUPERFUND TECHNICAL ASSESSMENT AND RESPONSE TEAM**  
 EPA CONTRACT 68-W5-0019  
 Phone: 908-225-4116 Fax: 908-225-7037

| MATRIX BOX NO.     | PRESERVATIVE BOX NO. |
|--------------------|----------------------|
| 1. Surface Water   | 1. HCl               |
| 2. Ground Water    | 2. HNO3              |
| 3. Leachate        | 3. Na2SO4            |
| 4. Rinseate        | 4. H2SO4             |
| 5. Soil/Sediment   | 5. Other (Specify)   |
| 6. Oil             | 6. Ice Only          |
| 7. Waste           | N. Not Preserved     |
| 8. Other (Specify) |                      |

Send verbal and written results to: **Roy F. Weston, Inc., USEPA Region II START**  
 Suite 201, 1090 King Georges Post Road, Edison, New Jersey 08837-3703  
 Attention: **Smita Sumbaly, START Analytical Coordinator**

| Sample Number | Sample Collection<br>MM/DD/YY/Time | Sample<br>Matrix<br>(Water<br>box #) | Conc.<br>Low-L<br>Med-M<br>High-H | Sample<br>Type<br>Comp-C<br>Grab-G | Sample<br>Preserv.<br>(Enter<br>box #) | RAS ANALYSIS |     |      |     |     | ICRA ANALYSIS |     |     | Full<br>TCLP | OTHER  |
|---------------|------------------------------------|--------------------------------------|-----------------------------------|------------------------------------|--|--------------|-----|------|-----|-----|---------------|-----|-----|--------------|--|
|               |                                    |                                      |                                   |                                    |  | YCB          | HNA | PEST | PCB | TAL | INC           | KIN | COX |              |  |
| LW-SS-01      | 4/03/00 0940                       | 5                                    | L                                 | G                                  |  |              |     |      | X   |     |               |     |     |              | 24 hr TAT  |
| LW-SS-02      | 4/03/00 1000                       | 5                                    | L                                 | G                                  |  |              |     |      | X   |     |               |     |     |              |  |
| LW-SS-03      | 4/03/00 1005                       | 5                                    | L                                 | G                                  |  |              |     |      | X   |     |               |     |     |              | Fax Results<br>To<br>Kevin Matheis<br>@ 716-827-303<br>Mail Hard copy<br>to USEPA<br>Kevin Matheis<br>3 Elm St<br>LACKAWANNA, N<br>14218 |

|  |                  |      |        |              |                              |                 |
|--|------------------|------|--------|--------------|------------------------------|-----------------|
| Person Assuming Responsibility for Sample: |                  |      |        |              | Time                         | Date (MM/DD/YY) |
| DAVID L ADAMS                              |                  |      |        |              | 0940                         | 4/3/00          |
| Sample Number                              | Relinquished By: | Time | Date   | Received By: | Reason for Change of Custody |                 |
| All  | David L Adams    | 1030 | 4/3/00 | John U...    | shipment to<br>LAB.          |                 |
| Sample Number                              | Relinquished By: | Time | Date   | Received By: | Reason for Change of Custody |                 |
|  |                  |      |        |              |                              |                 |
| Sample Number                              | Relinquished By: | Time | Date   | Received By: | Reason for Change of Custody |                 |
|  |                  |      |        |              |                              |                 |

Roy F. Weston, Inc.  
 FEDERAL PROGRAMS DIVISION  
 In Association with Resource Applications, Inc., E.E. Sartira Associates, PRC Environmental Management,  
 C.C. Johnson & Malhotra, P.C., and GRB Environmental Services, Inc.

**ATTACHMENT C**



DATE: / /

Upstate Laboratories, Inc.  
 Analysis Results  
 Report Number: 10800063  
 Client I.D.: EARTH TECH

APPROVAL: \_\_\_\_\_  
 QC: \_\_\_\_\_  
 Lab I.D.: 10170  
 Sampled by: Client

ID:10800063 Mat:Soil N20EO 0955H 04/12/00 G

| PARAMETERS                      | RESULTS    | DATE ANAL. | KEY | FILE#  |
|---------------------------------|------------|------------|-----|--------|
| Percent Solids                  | 71%        | 04/18/00   |     | WC9891 |
| Total Arsenic by furnace method | 14mg/kg dw | 04/21/00   |     | ME2754 |

ID:10800064 Mat:Soil N60EO 0959H 04/12/00 G

| PARAMETERS                      | RESULTS     | DATE ANAL. | KEY | FILE#  |
|---------------------------------|-------------|------------|-----|--------|
| Percent Solids                  | 76%         | 04/18/00   |     | WC9891 |
| Total Arsenic by furnace method | 2.5mg/kg dw | 04/21/00   |     | ME2754 |

ID:10800065 Mat:Soil ARCONF(3) 1000H 04/12/00 G

| PARAMETERS                      | RESULTS      | DATE ANAL. | KEY | FILE#  |
|---------------------------------|--------------|------------|-----|--------|
| Percent Solids                  | 73%          | 04/18/00   |     | WC9891 |
| Total Arsenic by furnace method | 0.30mg/kg dw | 04/21/00   |     | ME2754 |

ID:10800066 Mat:Soil ARCONF(4) 1005H 04/12/00 G

| PARAMETERS                      | RESULTS     | DATE ANAL. | KEY | FILE#  |
|---------------------------------|-------------|------------|-----|--------|
| Percent Solids                  | 89%         | 04/18/00   |     | WC9891 |
| Total Arsenic by furnace method | 1.6mg/kg dw | 04/21/00   |     | ME2754 |

ID:10800067 Mat:Soil ARCOF(6) 1010H 04/12/00 G

| PARAMETERS                      | RESULTS     | DATE ANAL. | KEY | FILE#  |
|---------------------------------|-------------|------------|-----|--------|
| Percent Solids                  | 84%         | 04/18/00   |     | WC9891 |
| Total Arsenic by furnace method | 1.0mg/kg dw | 04/21/00   |     | ME2754 |

dw = Dry weight

**ATTACHMENT D**

DATE: / /

Upstate Laboratories, Inc.  
Analysis Results  
Report Number: 10800055  
Client I.D.: EARTH TECH

APPROVAL: \_ \_ \_ \_  
QC: \_ \_ \_ \_ Lab I.D.: 10170  
Sampled by: Client

ID:10800055 Mat:Soil X1 0955H 04/11/00 G

| PARAMETERS                        | RESULTS        | DATE ANAL. | KEY | FILE#  |
|-----------------------------------|----------------|------------|-----|--------|
| Percent Solids                    | 85%            | 04/18/00   |     | WC9881 |
| PCB (Aroclors) by EPA Method 8080 |                |            |     |        |
| Aroclor 1016                      | <170ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Aroclor 1221                      | <170ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Aroclor 1232                      | <170ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Aroclor 1242                      | 12,000ug/kg dw | 04/27/00   |     | GA0038 |
| Aroclor 1248                      | <170ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Aroclor 1254                      | <170ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Aroclor 1260                      | <170ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Total PCB                         | 12,000ug/kg dw | 04/27/00   |     | GA0038 |

ID:10800055 Mat:Soil X4 1016H 04/11/00 G

| PARAMETERS                        | RESULTS        | DATE ANAL. | KEY | FILE#  |
|-----------------------------------|----------------|------------|-----|--------|
| Percent Solids                    | 85%            | 04/18/00   |     | WC9881 |
| PCB (Aroclors) by EPA Method 8080 |                |            |     |        |
| Aroclor 1016                      | <170ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Aroclor 1221                      | <170ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Aroclor 1232                      | <170ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Aroclor 1242                      | 33,000ug/kg dw | 04/27/00   |     | GA0038 |
| Aroclor 1248                      | <170ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Aroclor 1254                      | <170ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Aroclor 1260                      | <170ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Total PCB                         | 33,000ug/kg dw | 04/27/00   |     | GA0038 |

ID:10800057 Mat:Soil X5 1015H 04/11/00 G

| PARAMETERS                        | RESULTS        | DATE ANAL. | KEY | FILE#  |
|-----------------------------------|----------------|------------|-----|--------|
| Percent Solids                    | 85%            | 04/18/00   |     | WC9881 |
| PCB (Aroclors) by EPA Method 8080 |                |            |     |        |
| Aroclor 1016                      | <850ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Aroclor 1221                      | <850ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Aroclor 1232                      | <850ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Aroclor 1242                      | 62,000ug/kg dw | 04/27/00   |     | GA0038 |
| Aroclor 1248                      | <850ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Aroclor 1254                      | <850ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Aroclor 1260                      | <830ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Total PCB                         | 62,000ug/kg dw | 04/27/00   |     | GA0038 |

dw = Dry weight

DATE: / /

Upstate Laboratories, Inc.  
Analysis Results  
Report Number: 10900055  
Client I.D.: EARTH TECH

APPROVAL: \_\_\_\_\_  
QC: \_\_\_\_\_  
Lab I.D.: 10170  
Sampled by: Client

ID:10900058 Mat:Soil X6 1020H 04/11/00 G

| PARAMETERS                        | RESULTS        | DATE ANAL. | KEY | FILE#  |
|-----------------------------------|----------------|------------|-----|--------|
| Percent Solids                    | 81%            | 04/18/00   |     | WC9891 |
| PCB (Aroclors) by EPA Method 8080 |                |            |     |        |
| Aroclor 1016                      | <170ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Aroclor 1221                      | <170ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Aroclor 1232                      | <170ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Aroclor 1242                      | 16,000ug/kg dw | 04/27/00   |     | GA0038 |
| Aroclor 1248                      | <170ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Aroclor 1254                      | <170ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Aroclor 1260                      | <170ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Total PCB                         | 16,000ug/kg dw | 04/27/00   |     | GA0038 |

ID:10900059 Mat:Soil X7 1025H 04/11/00 G

| PARAMETERS                        | RESULTS        | DATE ANAL. | KEY | FILE#  |
|-----------------------------------|----------------|------------|-----|--------|
| Percent Solids                    | 78%            | 04/18/00   |     | WC9891 |
| PCB (Aroclors) by EPA Method 8080 |                |            |     |        |
| Aroclor 1016                      | <170ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Aroclor 1221                      | <170ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Aroclor 1232                      | <170ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Aroclor 1242                      | 11,000ug/kg dw | 04/27/00   |     | GA0038 |
| Aroclor 1248                      | <170ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Aroclor 1254                      | <170ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Aroclor 1260                      | <170ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Total PCB                         | 11,000ug/kg dw | 04/27/00   |     | GA0038 |

ID:10900050 Mat:Soil X10 1035H 04/12/00 G

| PARAMETERS                        | RESULTS      | DATE ANAL. | KEY | FILE#  |
|-----------------------------------|--------------|------------|-----|--------|
| Percent Solids                    | 35%          | 04/18/00   |     | WC9891 |
| PCB (Aroclors) by EPA Method 8080 |              |            |     |        |
| Aroclor 1016                      | <170ug/kg dw | 04/27/00   | 05  | GA0038 |
| Aroclor 1221                      | <170ug/kg dw | 04/27/00   | 05  | GA0038 |
| Aroclor 1232                      | <170ug/kg dw | 04/27/00   | 05  | GA0038 |
| Aroclor 1242                      | 9000ug/kg dw | 04/27/00   |     | GA0038 |
| Aroclor 1248                      | <170ug/kg dw | 04/27/00   | 05  | GA0038 |
| Aroclor 1254                      | <170ug/kg dw | 04/27/00   | 05  | GA0038 |
| Aroclor 1260                      | <170ug/kg dw | 04/27/00   | 05  | GA0038 |
| Total PCB                         | 9000ug/kg dw | 04/27/00   |     | GA0038 |

dw = Dry weight

DATE: / /

Upstate Laboratories, Inc.  
Analysis Results  
Report Number: 10800055  
Client I.D.: EARTH TECH

APPROVAL: \_\_\_\_\_  
QC: \_\_\_\_\_  
Lab I.D.: 10170  
Sampled by: Client

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ID:10800061 Mat:Soil X99 1040H 04/12/00 G  
-----

| PARAMETERS                        | RESULTS      | DATE ANAL. | KEY | FILE#  |
|-----------------------------------|--------------|------------|-----|--------|
| Percent Solids                    | 98%          | 04/18/00   |     | WC9891 |
| PCB (Aroclors) by EPA Method 8080 |              |            |     |        |
| Aroclor 1016                      | <170ug/kg dw | 04/27/00   | 05  | GA0038 |
| Aroclor 1221                      | <170ug/kg dw | 04/27/00   | 05  | GA0038 |
| Aroclor 1232                      | <170ug/kg dw | 04/27/00   | 05  | GA0038 |
| Aroclor 1242                      | 7000ug/kg dw | 04/27/00   |     | GA0038 |
| Aroclor 1248                      | <170ug/kg dw | 04/27/00   | 05  | GA0038 |
| Aroclor 1254                      | <170ug/kg dw | 04/27/00   | 05  | GA0038 |
| Aroclor 1260                      | <170ug/kg dw | 04/27/00   | 05  | GA0038 |
| Total PCB                         | 7000ug/kg dw | 04/27/00   |     | GA0038 |

-----  
ID:10800062 Mat:Soil X12 1415H 04/12/00 G  
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| PARAMETERS                        | RESULTS        | DATE ANAL. | KEY | FILE#  |
|-----------------------------------|----------------|------------|-----|--------|
| Percent Solids                    | 90%            | 04/18/00   |     | WC9891 |
| PCB (Aroclors) by EPA Method 8080 |                |            |     |        |
| Aroclor 1016                      | <170ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Aroclor 1221                      | <170ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Aroclor 1232                      | <170ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Aroclor 1242                      | 19,000ug/kg dw | 04/27/00   |     | GA0038 |
| Aroclor 1248                      | <170ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Aroclor 1254                      | <170ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Aroclor 1260                      | <170ug/kg dw   | 04/27/00   | 05  | GA0038 |
| Total PCB                         | 19,000ug/kg dw | 04/27/00   |     | GA0038 |

dw = Dry weight

**ATTACHMENT E**

DATE: 11/19/99

Upstate Laboratories, Inc.  
Analysis Results  
Report Number: 30299001  
Client I.D.: EARTH TECH  
Sampled by: Client

APPROVAL: *ajs*  
QC: *js*  
Lab I.D.: 10170

3467401 LACKAWANNA  
PIT PILE-PILE 1 AM 10/28/99 C

*Rock fill Material + Sub topsoil*

ULI I.D.: 30299002

Matrix: Soil

| PARAMETERS                       | RESULTS       | DATE ANAL. | KEY | FILE#  |
|----------------------------------|---------------|------------|-----|--------|
| Corrosivity                      |               |            |     |        |
| pH                               | 7.9SU         | 11/03/99   |     | WC8007 |
| Density                          | 1.20g/ml      | 11/10/99   |     | WC8111 |
| Flash Point                      | >60degC       | 11/11/99   |     | WC8125 |
| Percent Solids                   | 97%           | 10/29/99   |     | WC7966 |
| TOC                              | 14,138mg/kg   | 11/04/99   |     | SC0001 |
| RCRA Reactivity                  |               |            |     |        |
| Reactive Sulfide                 | <50mg/kg      | 10/29/99   |     | WC7978 |
| Reactive Cyanide                 | <1.0mg/kg     | 11/10/99   |     | WC8048 |
| Total Aluminum                   | 1900mg/kg dw  | 11/03/99   |     | ME2404 |
| Total Antimony                   | <0.4mg/kg dw  | 11/03/99   |     | ME2404 |
| Total Arsenic by furnace method  | 5.5mg/kg dw   | 11/03/99   |     | ME2404 |
| Total Barium                     | 42mg/kg dw    | 11/03/99   |     | ME2404 |
| Total Beryllium                  | 0.62mg/kg dw  | 11/03/99   |     | ME2404 |
| Total Cadmium                    | 1.6mg/kg dw   | 11/03/99   |     | ME2404 |
| Total Calcium                    | 5000mg/kg dw  | 11/10/99   |     | ME2404 |
| Total Chromium                   | 20mg/kg dw    | 11/03/99   |     | ME2404 |
| Total Cobalt                     | <5.1mg/kg dw  | 11/03/99   |     | ME2404 |
| Total Copper                     | 83mg/kg dw    | 11/03/99   |     | ME2404 |
| Total Iron                       | 19000mg/kg dw | 11/03/99   |     | ME2404 |
| Total Lead                       | 82mg/kg dw    | 11/03/99   |     | ME2404 |
| Total Magnesium                  | 930mg/kg dw   | 11/10/99   |     | ME2413 |
| Total Manganese                  | 670mg/kg dw   | 11/03/99   |     | ME2404 |
| Total Mercury                    | 0.23mg/kg dw  | 11/09/99   |     | MB1662 |
| Total Nickel                     | 18mg/kg dw    | 11/03/99   |     | ME2404 |
| Total Potassium                  | 240mg/kg dw   | 11/10/99   |     | ME2413 |
| Total Selenium by furnace method | <0.2mg/kg dw  | 11/03/99   |     | ME2404 |
| Total Silver                     | <5.1mg/kg dw  | 11/03/99   |     | ME2404 |
| Total Sodium                     | 280mg/kg dw   | 11/10/99   |     | ME2413 |
| Total Thallium by furnace method | <0.4mg/kg dw  | 11/03/99   |     | ME2404 |
| Total Vanadium                   | <31mg/kg dw   | 11/03/99   |     | ME2404 |
| Total Zinc                       | 160mg/kg dw   | 11/03/99   |     | ME2404 |
| TCLP Arsenic                     | 0.26mg/l      | 11/10/99   |     | ME2415 |
| TCLP Barium                      | 0.4mg/l       | 11/10/99   |     | ME2415 |
| TCLP Cadmium                     | 0.014mg/l     | 11/10/99   |     | ME2415 |
| TCLP Chromium                    | <0.05mg/l     | 11/10/99   |     | ME2415 |
| TCLP Lead                        | <0.1mg/l      | 11/10/99   |     | ME2415 |
| TCLP Mercury                     | <0.0004mg/l   | 11/09/99   |     | MB1662 |

dw = Dry weight

DATE: 11/19/99

Upstate Laboratories, Inc.  
Analysis Results  
Report Number: 30299001  
Client I.D.: EARTH TECH  
Sampled by: Client

APPROVAL: *OJS*  
QC: *JP*  
Lab I.D.: 10170

3467401 LACKAWANNA  
PIT PILE-PILE 1 AM 10/28/99 C

ULI I.D.: 30299002

Matrix: Soil

| PARAMETERS    | RESULTS   | DATE ANAL. | KEY | FILE#  |
|---------------|-----------|------------|-----|--------|
| TCLP Selenium | <0.5mg/l  | 11/10/99   |     | ME2415 |
| TCLP Silver   | <0.05mg/l | 11/10/99   |     | ME2415 |

TCL Volatiles by EPA Method 8260

|                           |          |          |  |        |
|---------------------------|----------|----------|--|--------|
| Chloromethane             | <3ug/kg  | 11/08/99 |  | VM2659 |
| Bromomethane              | <3ug/kg  | 11/08/99 |  | VM2659 |
| Vinyl Chloride            | <2ug/kg  | 11/08/99 |  | VM2659 |
| Chloroethane              | <3ug/kg  | 11/08/99 |  | VM2659 |
| Methylene Chloride        | 100ug/kg | 11/08/99 |  | VM2659 |
| Acetone                   | <10ug/kg | 11/08/99 |  | VM2659 |
| Carbon Disulfide          | <3ug/kg  | 11/08/99 |  | VM2659 |
| 1,1-Dichloroethene        | <3ug/kg  | 11/08/99 |  | VM2659 |
| 1,1-Dichloroethane        | <3ug/kg  | 11/08/99 |  | VM2659 |
| trans-1,2-Dichloroethene  | <3ug/kg  | 11/08/99 |  | VM2659 |
| cis-1,2-Dichloroethene    | <3ug/kg  | 11/08/99 |  | VM2659 |
| Chloroform                | 3ug/kg   | 11/08/99 |  | VM2659 |
| 1,2-Dichloroethane        | <3ug/kg  | 11/08/99 |  | VM2659 |
| 2-Butanone                | <10ug/kg | 11/08/99 |  | VM2659 |
| 1,1,1-Trichloroethane     | <3ug/kg  | 11/08/99 |  | VM2659 |
| Carbon Tetrachloride      | <3ug/kg  | 11/08/99 |  | VM2659 |
| Bromodichloromethane      | <3ug/kg  | 11/08/99 |  | VM2659 |
| 1,2-Dichloropropane       | <3ug/kg  | 11/08/99 |  | VM2659 |
| cis-1,3-Dichloropropene   | <3ug/kg  | 11/08/99 |  | VM2659 |
| Trichloroethene           | <3ug/kg  | 11/08/99 |  | VM2659 |
| Dibromochloromethane      | <3ug/kg  | 11/08/99 |  | VM2659 |
| 1,1,2-Trichloroethane     | <3ug/kg  | 11/08/99 |  | VM2659 |
| Benzene                   | <3ug/kg  | 11/08/99 |  | VM2659 |
| trans-1,3-Dichloropropene | <3ug/kg  | 11/08/99 |  | VM2659 |
| Bromoform                 | <3ug/kg  | 11/08/99 |  | VM2659 |
| 4-Methyl-2-pentanone      | <10ug/kg | 11/08/99 |  | VM2659 |
| 2-Hexanone                | <10ug/kg | 11/08/99 |  | VM2659 |
| Tetrachloroethene         | <3ug/kg  | 11/08/99 |  | VM2659 |
| 1,1,2,2-Tetrachloroethane | <3ug/kg  | 11/08/99 |  | VM2659 |
| Toluene                   | <3ug/kg  | 11/08/99 |  | VM2659 |
| Chlorobenzene             | <3ug/kg  | 11/08/99 |  | VM2659 |
| Ethylbenzene              | <3ug/kg  | 11/08/99 |  | VM2659 |
| Styrene                   | <3ug/kg  | 11/08/99 |  | VM2659 |
| m-Xylene and p-Xylene     | <3ug/kg  | 11/08/99 |  | VM2659 |
| o-Xylene                  | <3ug/kg  | 11/08/99 |  | VM2659 |

dw = Dry weight



DATE: 11/19/99

Upstate Laboratories, Inc.  
Analysis Results  
Report Number: 30299001  
Client I.D.: EARTH TECH  
Sampled by: Client

APPROVAL: *ajs*  
QC: *sp*  
Lab I.D.: 10170

3467401 LACKAWANNA  
PIT PILE-PILE 1 AM 10/28/99 C

ULI I.D.: 30299002

Matrix: Soil

| PARAMETERS                              | RESULTS       | DATE ANAL. | KEY | FILE#  |
|---|---------------|------------|-----|--------|
| TCLP Volatile Organic Compounds by 8260 |               |            |     |        |
| TCLP Benzene                            | <0.03mg/l     | 11/08/99   |     | VM2659 |
| TCLP Carbon Tetrachloride               | <0.03mg/l     | 11/08/99   |     | VM2659 |
| TCLP Chlorobenzene                      | <0.03mg/l     | 11/08/99   |     | VM2659 |
| TCLP Chloroform                         | <0.03mg/l     | 11/08/99   |     | VM2659 |
| TCLP 1,4-Dichlorobenzene                | <0.03mg/l     | 11/08/99   |     | VM2659 |
| TCLP 1,2-Dichloroethane                 | <0.03mg/l     | 11/08/99   |     | VM2659 |
| TCLP 1,1-Dichloroethene                 | <0.03mg/l     | 11/08/99   |     | VM2659 |
| TCLP Methyl Ethyl Ketone                | <0.1mg/l      | 11/08/99   |     | VM2659 |
| TCLP Tetrachloroethene                  | <0.03mg/l     | 11/08/99   |     | VM2659 |
| TCLP Trichloroethene                    | <0.03mg/l     | 11/08/99   |     | VM2659 |
| TCLP Vinyl Chloride                     | <0.02mg/l     | 11/08/99   |     | VM2659 |
| TCL Semivolatiles by EPA Method 8270    |               |            |     |        |
| Phenol                                  | 8700ug/kg dw  | 11/11/99   |     | SA2200 |
| bis(2-Chloroethyl) ether                | <3400ug/kg dw | 11/11/99   |     | SA2200 |
| 2-Chlorophenol                          | <3400ug/kg dw | 11/11/99   |     | SA2200 |
| 1,3-Dichlorobenzene                     | <3400ug/kg dw | 11/11/99   |     | SA2200 |
| 1,4-Dichlorobenzene                     | <3400ug/kg dw | 11/11/99   |     | SA2200 |
| 1,2-Dichlorobenzene                     | <3400ug/kg dw | 11/11/99   |     | SA2200 |
| 2-Methylphenol                          | <3400ug/kg dw | 11/11/99   |     | SA2200 |
| 2,2'-Oxybis(1-Chloropropane)            | <3400ug/kg dw | 11/11/99   |     | SA2200 |
| 4-Methylphenol                          | <3400ug/kg dw | 11/11/99   |     | SA2200 |
| n-Nitrosodi-n-propylamine               | <3400ug/kg dw | 11/11/99   |     | SA2200 |
| Hexachloroethane                        | <3400ug/kg dw | 11/11/99   |     | SA2200 |
| Nitrobenzene                            | <3400ug/kg dw | 11/11/99   |     | SA2200 |
| Isophorone                              | <3400ug/kg dw | 11/11/99   |     | SA2200 |
| 2-Nitrophenol                           | <3400ug/kg dw | 11/11/99   |     | SA2200 |
| 2,4-Dimethylphenol                      | <3400ug/kg dw | 11/11/99   |     | SA2200 |
| bis(2-Chloroethoxy)methane              | <3400ug/kg dw | 11/11/99   |     | SA2200 |
| 2,4-Dichlorophenol                      | <3400ug/kg dw | 11/11/99   |     | SA2200 |
| 1,2,4-Trichlorobenzene                  | <3400ug/kg dw | 11/11/99   |     | SA2200 |
| Naphthalene                             | <3400ug/kg dw | 11/11/99   |     | SA2200 |
| 4-Chloroaniline                         | <3400ug/kg dw | 11/11/99   |     | SA2200 |
| Hexachlorobutadiene                     | <3400ug/kg dw | 11/11/99   |     | SA2200 |
| 4-Chloro-3-methylphenol                 | <3400ug/kg dw | 11/11/99   |     | SA2200 |
| 2-Methylnaphthalene                     | <3400ug/kg dw | 11/11/99   |     | SA2200 |
| Hexachlorocyclopentadiene               | <3400ug/kg dw | 11/11/99   |     | SA2200 |
| 2,4,6-Trichlorophenol                   | <3400ug/kg dw | 11/11/99   |     | SA2200 |

dw = Dry weight

DATE: 11/19/99

Upstate Laboratories, Inc.  
Analysis Results  
Report Number: 30299001  
Client I.D.: EARTH TECH  
Sampled by: Client

APPROVAL: *ajs*  
QC: *js*  
Lab I.D.: 10170

3467401 LACKAWANNA  
PIT PILE-PILE 1 AM 10/28/99 C

ULI I.D.: 30299002

Matrix: Soil

| PARAMETERS                   | RESULTS         | DATE ANAL. | KEY | FILE#  |
|------------------------------|-----------------|------------|-----|--------|
| 2,4,5-Trichlorophenol        | <3400ug/kg dw   | 11/11/99   |     | SA2200 |
| 2-Chloronaphthalene          | <3400ug/kg dw   | 11/11/99   |     | SA2200 |
| 2-Nitroaniline               | <34,000ug/kg dw | 11/11/99   |     | SA2200 |
| Dimethylphthalate            | <3400ug/kg dw   | 11/11/99   |     | SA2200 |
| Acenaphthylene               | <3400ug/kg dw   | 11/11/99   |     | SA2200 |
| 2,6-Dinitrotoluene           | <3400ug/kg dw   | 11/11/99   |     | SA2200 |
| 3-Nitroaniline               | <34,000ug/kg dw | 11/11/99   |     | SA2200 |
| Acenaphthene                 | <3400ug/kg dw   | 11/11/99   |     | SA2200 |
| 2,4-Dinitrophenol            | <34,000ug/kg dw | 11/11/99   |     | SA2200 |
| 4-Nitrophenol                | <34,000ug/kg dw | 11/11/99   |     | SA2200 |
| Dibenzofuran                 | <3400ug/kg dw   | 11/11/99   |     | SA2200 |
| 2,4-Dinitrotoluene           | <3400ug/kg dw   | 11/11/99   |     | SA2200 |
| Diethylphthalate             | <3400ug/kg dw   | 11/11/99   |     | SA2200 |
| 4-Chlorophenylphenylether    | <3400ug/kg dw   | 11/11/99   |     | SA2200 |
| Fluorene                     | <3400ug/kg dw   | 11/11/99   |     | SA2200 |
| 4-Nitroaniline               | <34,000ug/kg dw | 11/11/99   |     | SA2200 |
| 2-Methyl-4,6-dinitrophenol   | <34,000ug/kg dw | 11/11/99   |     | SA2200 |
| n-Nitrosodiphenylamine       | <3400ug/kg dw   | 11/11/99   |     | SA2200 |
| 4-Bromophenylphenylether     | <3400ug/kg dw   | 11/11/99   |     | SA2200 |
| Hexachlorobenzene            | <3400ug/kg dw   | 11/11/99   |     | SA2200 |
| Pentachlorophenol            | <690ug/kg dw    | 11/11/99   |     | SA2200 |
| Phenanthrene                 | <3400ug/kg dw   | 11/11/99   |     | SA2200 |
| Anthracene                   | <3400ug/kg dw   | 11/11/99   |     | SA2200 |
| Carbazole                    | <3400ug/kg dw   | 11/11/99   |     | SA2200 |
| di-n-butylphthalate          | <3400ug/kg dw   | 11/11/99   |     | SA2200 |
| Fluoranthene                 | <3400ug/kg dw   | 11/11/99   |     | SA2200 |
| Pyrene                       | <3400ug/kg dw   | 11/11/99   |     | SA2200 |
| Butylbenzylphthalate         | <3400ug/kg dw   | 11/11/99   |     | SA2200 |
| 3,3'-Dichlorobenzidine       | <3400ug/kg dw   | 11/11/99   |     | SA2200 |
| Benzo (a) anthracene         | <3400ug/kg dw   | 11/11/99   |     | SA2200 |
| Chrysene                     | <3400ug/kg dw   | 11/11/99   |     | SA2200 |
| bis (2-Ethylhexyl) phthalate | <3400ug/kg dw   | 11/11/99   |     | SA2200 |
| di-n-octylphthalate          | <3400ug/kg dw   | 11/11/99   |     | SA2200 |
| Benzo (b) fluoranthene       | <3400ug/kg dw   | 11/11/99   |     | SA2200 |
| Benzo (k) fluoranthene       | <3400ug/kg dw   | 11/11/99   |     | SA2200 |
| Benzo (a) pyrene             | <3400ug/kg dw   | 11/11/99   |     | SA2200 |
| Indeno (1,2,3-cd) pyrene     | <690ug/kg dw    | 11/11/99   |     | SA2200 |
| Dibenzo (a, h) anthracene    | <690ug/kg dw    | 11/11/99   |     | SA2200 |
| Benzo (ghi) perylene         | <690ug/kg dw    | 11/11/99   |     | SA2200 |

dw = Dry weight

DATE: 11/19/99

Upstate Laboratories, Inc.  
Analysis Results  
Report Number: 30299001  
Client I.D.: EARTH TECH  
Sampled by: Client

APPROVAL: *ajs*  
QC: *jo*  
Lab I.D.: 10170

3467401 LACKAWANNA  
PIT PILE-PILE 1 AM 10/28/99 C

ULI I.D.: 30299002

Matrix: Soil

| PARAMETERS                          | RESULTS       | DATE ANAL. | KEY | FILE#  |
|-------------------------------------|---------------|------------|-----|--------|
| TCLP Semivolatile Compounds by 8270 |               |            |     |        |
| TCLP Cresol, Total                  | <0.05mg/l     | 11/11/99   |     | SA2207 |
| TCLP 2,4-Dinitrotoluene             | <0.05mg/l     | 11/11/99   |     | SA2207 |
| TCLP Hexachlorobenzene              | <0.05mg/l     | 11/11/99   |     | SA2207 |
| TCLP Hexachlorobutadiene            | <0.05mg/l     | 11/11/99   |     | SA2207 |
| TCLP Hexachloroethane               | <0.05mg/l     | 11/11/99   |     | SA2207 |
| TCLP Nitrobenzene                   | <0.05mg/l     | 11/11/99   |     | SA2207 |
| TCLP Pentachlorophenol              | <0.10mg/l     | 11/11/99   |     | SA2207 |
| TCLP Pyridine                       | <0.05mg/l     | 11/11/99   |     | SA2207 |
| TCLP 2,4,5-Trichlorophenol          | <0.05mg/l     | 11/11/99   |     | SA2207 |
| TCLP 2,4,6-Trichlorophenol          | <0.05mg/l     | 11/11/99   |     | SA2207 |
| PCB (Aroclors) by EPA Method 8082   |               |            |     |        |
| Aroclor 1016                        | <0.08mg/kg dw | 11/17/99   |     | PA5354 |
| Aroclor 1221                        | <0.08mg/kg dw | 11/17/99   |     | PA5354 |
| Aroclor 1232                        | <0.08mg/kg dw | 11/17/99   |     | PA5354 |
| Aroclor 1242                        | <0.08mg/kg dw | 11/17/99   |     | PA 4   |
| Aroclor 1248                        | <0.08mg/kg dw | 11/17/99   |     | PA 4   |
| Aroclor 1254                        | <0.08mg/kg dw | 11/17/99   |     | PA5354 |
| Aroclor 1260                        | <0.08mg/kg dw | 11/17/99   |     | PA5354 |
| Total PCB                           | <0.08mg/kg dw | 11/17/99   |     | PA5354 |

dw = Dry weight

| PARAMETER                        | SO27        | PILE 1        |
|----------------------------------|-------------|---------------|
| Arsenic                          | <0.5mg/l    | 0.26mg/l      |
| Barium                           | <0.3mg/l    | 0.4mg/l       |
| Cadmium                          | 0.018mg/l   | 0.014mg/l     |
| Chromium                         | <0.05mg/l   | <0.05mg/l     |
| Copper                           | 0.21mg/l    |               |
| Lead                             | <0.1mg/l    | <0.1mg/l      |
| Mercury                          | <0.0004mg/l | <0.0004mg/l   |
| Selenium                         | <0.5mg/l    | <0.5mg/l      |
| Silver                           | <0.05mg/l   | <0.05mg/l     |
| Zinc                             | 1.3mg/l     |               |
| pH                               |             | 7.9SU         |
| Density                          |             | 1.20g/ml      |
| Flash Point                      |             | >60degC       |
| Percent Solids                   |             | 97%           |
| TOC                              |             | 14,138mg/kg   |
| Reactive Sulfide                 |             | <50mg/kg      |
| Reactive Cyanide                 |             | <1.0mg/kg     |
| Total Aluminum                   |             | 1900mg/kg dw  |
| Total Antimony                   |             | <0.4mg/kg dw  |
| Total Arsenic by furnace method  |             | 5.5mg/kg dw   |
| Total Barium                     |             | 42mg/kg dw    |
| Total Beryllium                  |             | 0.62mg/kg dw  |
| Total Cadmium                    |             | 1.6mg/kg dw   |
| Total Calcium                    |             | 5000mg/kg dw  |
| Total Chromium                   |             | 20mg/kg dw    |
| Total Cobalt                     |             | <5.1mg/kg dw  |
| Total Copper                     |             | 83mg/kg dw    |
| Total Iron                       |             | 19000mg/kg dw |
| Total Lead                       |             | 82mg/kg dw    |
| Total Magnesium                  |             | 930mg/kg dw   |
| Total Manganese                  |             | 670mg/kg dw   |
| Total Mercury                    |             | 0.23mg/kg dw  |
| Total Nickel                     |             | 18mg/kg dw    |
| Total Potassium                  |             | 240mg/kg dw   |
| Total Selenium by furnace method |             | <0.2mg/kg dw  |
| Total Silver                     |             | <5.1mg/kg dw  |
| Total Sodium                     |             | 280mg/kg dw   |



| PARAMETER                        | SO27 | PILE 1         |
|----------------------------------|------|----------------|
| o - Xylene                       |      | < 3ug/kg       |
| Benzene                          |      | < 0.03mg/l     |
| Carbon Tetrachloride             |      | < 0.03mg/l     |
| Chlorobenzene                    |      | < 0.03mg/l     |
| Chloroform                       |      | < 0.03mg/l     |
| 1,4 - Dichlorobenzene            |      | < 0.03mg/l     |
| 1,2 - Dichloroethane             |      | < 0.03mg/l     |
| 1,1 - Dichloroethene             |      | < 0.03mg/l     |
| Methyl Ethyl Ketone              |      | < 0.1mg/l      |
| Tetrachloroethene                |      | < 0.03mg/l     |
| Trichloroethene                  |      | < 0.03mg/l     |
| Vinyl Chloride                   |      | < 0.02mg/l     |
| Phenol                           |      | 8700ug/kg dw   |
| bis(2 - Chloroethyl) ether       |      | < 3400ug/kg dw |
| 2 - Chlorophenol                 |      | < 3400ug/kg dw |
| 1,3 - Dichlorobenzene            |      | < 3400ug/kg dw |
| 1,4 - Dichlorobenzene            |      | < 3400ug/kg dw |
| 1,2 - Dichlorobenzene            |      | < 3400ug/kg dw |
| 2 - Methylphenol                 |      | < 3400ug/kg dw |
| 2,2' - Oxybis(1 - Chloropropane) |      | < 3400ug/kg dw |
| 4 - Methylphenol                 |      | < 3400ug/kg dw |
| n - Nitrosodi - n - propylamine  |      | < 3400ug/kg dw |
| Hexachloroethane                 |      | < 3400ug/kg dw |
| Nitrobenzene                     |      | < 3400ug/kg dw |
| Isophorone                       |      | < 3400ug/kg dw |
| 2 - Nitrophenol                  |      | < 3400ug/kg dw |
| 2,4 - Dimethylphenol             |      | < 3400ug/kg dw |
| bis(2 - Chloroethoxy)methane     |      | < 3400ug/kg dw |
| 2,4 - Dichlorophenol             |      | < 3400ug/kg dw |
| 1,2,4 - Trichlorobenzene         |      | < 3400ug/kg dw |
| Naphthalene                      |      | < 3400ug/kg dw |
| 4 - Chloroaniiline               |      | < 3400ug/kg dw |
| Hexachlorobutadiene              |      | < 3400ug/kg dw |
| 4 - Chloro - 3 - methylphenol    |      | < 3400ug/kg dw |
| 2 - Methylinaphthalene           |      | < 3400ug/kg dw |
| Hexachlorocyclopentadiene        |      | < 3400ug/kg dw |
| 2,4,6 - Trichlorophenol          |      | < 3400ug/kg dw |

## SO27

## PILE 1

|                               |                 |
|-------------------------------|-----------------|
| PARAMETER                     |                 |
| 2,4,5- Trichlorophenol        | <3400ug/kg dw   |
| 2- Chloronaphthalene          | <3400ug/kg dw   |
| 2- Nitroaniline               | <34,000ug/kg dw |
| Dimethylphthalate             | <3400ug/kg dw   |
| Acenaphthylene                | <3400ug/kg dw   |
| 2,6- Dinitrotoluene           | <34,000ug/kg dw |
| 3- Nitroaniline               | <3400ug/kg dw   |
| Acenaphthene                  | <34,000ug/kg dw |
| 2,4- Dinitrophenol            | <34,000ug/kg dw |
| 4- Nitrophenol                | <34,000ug/kg dw |
| Dibenzofuran                  | <3400ug/kg dw   |
| 2,4- Dinitrotoluene           | <3400ug/kg dw   |
| Diethylphthalate              | <3400ug/kg dw   |
| 4- Chlorophenylphenylether    | <3400ug/kg dw   |
| Fluorene                      | <3400ug/kg dw   |
| 4- Nitroaniline               | <34,000ug/kg dw |
| 2- Methyl- 4,6- dinitrophenol | <34,000ug/kg dw |
| n- Nitrosodiphenylamine       | <3400ug/kg dw   |
| 4- Bromophenylphenylether     | <3400ug/kg dw   |
| Hexachlorobenzene             | <3400ug/kg dw   |
| Pentachlorophenol             | <690ug/kg dw    |
| Phenanthrene                  | <3400ug/kg dw   |
| Anthracene                    | <3400ug/kg dw   |
| Carbazole                     | <3400ug/kg dw   |
| di- n- butylphthalate         | <3400ug/kg dw   |
| Fluoranthene                  | <3400ug/kg dw   |
| Pyrene                        | <3400ug/kg dw   |
| Butylbenzylphthalate          | <3400ug/kg dw   |
| 3,3'- Dichlorobenzidine       | <3400ug/kg dw   |
| Benzo(a)anthracene            | <3400ug/kg dw   |
| Chrysene                      | <3400ug/kg dw   |
| bis(2- Ethylhexyl)phthalate   | <3400ug/kg dw   |
| di- n- octylphthalate         | <3400ug/kg dw   |
| Benzo(b)fluoranthene          | <3400ug/kg dw   |
| Benzo(k)fluoranthene          | <3400ug/kg dw   |
| Benzo(a)pyrene                | <3400ug/kg dw   |
| Indeno(1,2,3- cd)pyrene       | <690ug/kg dw    |

| PARAMETER              | SO27 | PILE 1        |
|------------------------|------|---------------|
| Dibenzo(a,h)anthracene |      | <690ug/kg dw  |
| Benzo(ghi)perylene     |      | <690ug/kg dw  |
| Cresol, Total          |      | <0.05mg/l     |
| 2,4-Dinitrotoluene     |      | <0.05mg/l     |
| Hexachlorobenzene      |      | <0.05mg/l     |
| Hexachlorobutadiene    |      | <0.05mg/l     |
| Hexachloroethane       |      | <0.05mg/l     |
| Nitrobenzene           |      | <0.05mg/l     |
| Pentachlorophenol      |      | <0.10mg/l     |
| Pyridine               |      | <0.05mg/l     |
| 2,4,5-Trichlorophenol  |      | <0.05mg/l     |
| 2,4,6-Trichlorophenol  |      | <0.05mg/l     |
| Aroclor 1016           |      | <0.08mg/kg dw |
| Aroclor 1221           |      | <0.08mg/kg dw |
| Aroclor 1232           |      | <0.08mg/kg dw |
| Aroclor 1242           |      | <0.08mg/kg dw |
| Aroclor 1248           |      | <0.08mg/kg dw |
| Aroclor 1254           |      | <0.08mg/kg dw |
| Aroclor 1260           |      | <0.08mg/kg dw |
| Total PCB              |      | <0.08mg/kg dw |





**ATTACHMENT F**

# Upstate Laboratories Inc.

Shipping: 6034 Corporate Dr. • E. Syracuse, NY 13067-1017 • (315) 437-0255 • Fax (315) 437-1209  
Mailing: Box 289 • Syracuse, NY 13206  
Albany (518) 459-3134  
Binghamton (607) 724-0478

Buffalo (716) 649-2533  
Rochester (716) 436-9070  
New Jersey (201) 703-1324

February 9, 2000

Ms. Lane Aulick  
Earth Tech, Inc.  
7870 Villa Park Dr.  
Suite 400  
Richmond, VA 23228

Re: Analysis Report #35599012 - 38340 Lackawanna Foundry

Dear Ms. Aulick:

Please find enclosed the results for your samples which were picked up by ULI personnel on December 20, 1999.

We have included the Chain of Custody Record as part of your report. You may need to reference this form for a more detailed explanation of your sample. Samples will be disposed of approximately one month from final report date.

Should you have any questions, please feel free to give us a call.

Thank you for your patronage.

Sincerely,

UPSTATE LABORATORIES, INC.

*Anthony J. Scala*  
Anthony J. Scala  
Director

AJS/lw

Enclosures: report, spreadsheets, disk

cc/encs: N. Scala, ULI  
file  
M. Kromis, Earth Tech (category B deliverables to follow)

Note: Faxed results were given to your office on 12/20/99. AJS

Disclaimer: The test results and procedures utilized, and laboratory interpretations of data obtained by ULI as contained in this report are believed by ULI to be accurate and reliable for sample(s) tested. In accepting this report, the customer agrees that the full extent of any and all liability for actual and consequential damages of ULI for the services performed shall be equal to the fee charged to the customer for the services as liquidated damages.

DATE: 02/09/00

Upstate Laboratories, Inc.  
 Analysis Results  
 Report Number: 35599012  
 Client I.D.: EARTH TECH

APPROVAL: *gjs*  
 QC: *gjs*  
 Lab I.D.: 10170  
 Sampled by: Client

ID:35599012 Mat:Soil 38340 LACKAWANNA FOUNDRY CONF 7 SUB-SOIL PCB/ 1000H 12/20/99 G  
 TRANSFORMER AREA

| PARAMETERS                        | RESULTS       | DATE ANAL. | KEY | FILE#  |
|-----------------------------------|---------------|------------|-----|--------|
| Percent Solids                    | 59%           | 12/22/99   |     | WC8615 |
| PCB (Aroclors) by EPA Method 8080 |               |            |     |        |
| Aroclor 1016                      | <0.14ug/kg dw | 01/06/00   |     | PA5467 |
| Aroclor 1221                      | <0.14ug/kg dw | 01/06/00   |     | PA5467 |
| Aroclor 1232                      | <0.14ug/kg dw | 01/06/00   |     | PA5467 |
| Aroclor 1242                      | <0.14ug/kg dw | 01/06/00   |     | PA5467 |
| Aroclor 1248                      | <0.14ug/kg dw | 01/06/00   |     | PA5467 |
| Aroclor 1254                      | <0.14ug/kg dw | 01/06/00   |     | PA5467 |
| Aroclor 1260                      | <0.14ug/kg dw | 01/06/00   |     | PA5467 |
| Total PCB                         | <0.14ug/kg dw | 01/06/00   |     | PA5467 |

ID:35599013 Mat:Soil 38340 LACKAWANNA FOUNDRY CONF 8 SUB-SOIL PCB/ 1000H 12/20/99 G  
 TRANSFORMER AREA

| PARAMETERS                        | RESULTS       | DATE ANAL. | KEY | FILE#  |
|-----------------------------------|---------------|------------|-----|--------|
| Percent Solids                    | 70%           | 12/22/99   |     | WC8615 |
| PCB (Aroclors) by EPA Method 8080 |               |            |     |        |
| Aroclor 1016                      | <0.11ug/kg dw | 01/06/00   |     | PA5467 |
| Aroclor 1221                      | <0.11ug/kg dw | 01/06/00   |     | PA5467 |
| Aroclor 1232                      | <0.11ug/kg dw | 01/06/00   |     | PA5467 |
| Aroclor 1242                      | 1400ug/kg dw  | 01/06/00   |     | PA5467 |
| Aroclor 1248                      | <0.11ug/kg dw | 01/06/00   |     | PA5467 |
| Aroclor 1254                      | <0.11ug/kg dw | 01/06/00   |     | PA5467 |
| Aroclor 1260                      | <0.11ug/kg dw | 01/06/00   |     | PA5467 |
| Total PCB                         | 1400ug/kg dw  | 01/06/00   |     | PA5467 |

ID:35599014 Mat:Soil 38340 LACKAWANNA FOUNDRY CONF 9 SUB-SOIL PCB/ 1000H 12/20/99 G  
 TRANSFORMER AREA

| PARAMETERS                        | RESULTS       | DATE ANAL. | KEY | FILE#  |
|-----------------------------------|---------------|------------|-----|--------|
| Percent Solids                    | 66%           | 12/22/99   |     | WC8615 |
| PCB (Aroclors) by EPA Method 8080 |               |            |     |        |
| Aroclor 1016                      | <0.12ug/kg dw | 01/06/00   |     | PA5467 |
| Aroclor 1221                      | <0.12ug/kg dw | 01/06/00   |     | PA5467 |
| Aroclor 1232                      | <0.12ug/kg dw | 01/06/00   |     | PA5467 |
| Aroclor 1242                      | <0.12ug/kg dw | 01/06/00   |     | PA5467 |
| Aroclor 1248                      | <0.12ug/kg dw | 01/06/00   |     | PA5467 |
| Aroclor 1254                      | <0.12ug/kg dw | 01/06/00   |     | PA5467 |
| Aroclor 1260                      | <0.12ug/kg dw | 01/06/00   |     | PA5467 |
| Total PCB                         | <0.12ug/kg dw | 01/06/00   |     | PA5467 |

dw = Dry weight

DATE: 02/09/00

Upstate Laboratories, Inc.  
 Analysis Results  
 Report Number: 35599012  
 Client I.D.: EARTH TECH

APPROVAL: *[Signature]*  
 QC: *[Signature]* Lab I.D.: 10170  
 Sampled by: Client

ID: 35599015 Mat: Soil 38340 LACKAWANNA FOUNDRY CONF 10 SUB-SOIL PCB/ 1000H 12/20/99 G  
 TRANSFORMER AREA

| PARAMETERS                        | RESULTS     | DATE ANAL. | KEY | FILE#  |
|-----------------------------------|-------------|------------|-----|--------|
| Percent Solids                    | 74%         | 12/22/99   |     | WC8615 |
| PCB (Aroclors) by EPA Method 8080 |             |            |     |        |
| Aroclor 1016                      | <54ug/kg dw | 01/06/00   | 05  | PA5467 |
| Aroclor 1221                      | <54ug/kg dw | 01/06/00   | 05  | PA5467 |
| Aroclor 1232                      | <54ug/kg dw | 01/06/00   | 05  | PA5467 |
| Aroclor 1242                      | <54ug/kg dw | 01/06/00   | 05  | PA5467 |
| Aroclor 1248                      | <54ug/kg dw | 01/06/00   | 05  | PA5467 |
| Aroclor 1254                      | <54ug/kg dw | 01/06/00   | 05  | PA5467 |
| Aroclor 1260                      | <54ug/kg dw | 01/06/00   | 05  | PA5467 |
| Total PCB                         | <54ug/kg dw | 01/06/00   | 05  | PA5467 |

dw = Dry weight

## KEY PAGE

1 MATRIX INTERFERENCE PRECLUDES LOWER DETECTION LIMITS  
 2 MATRIX INTERFERENCE  
 3 PRESENT IN BLANK  
 4 ANALYSIS NOT PERFORMED BECAUSE OF INSUFFICIENT SAMPLE  
 5 THE PRESENCE OF OTHER TARGET ANALYTE(S) PRECLUDES LOWER DETECTION LIMITS  
 6 BLANK CORRECTED  
 7 HEAD SPACE PRESENT IN SAMPLE  
 8 QUANTITATION LIMIT IS GREATER THAN THE CALCULATED REGULATORY LEVEL. THE  
 9 QUANTITATION LIMIT THEREFORE BECOMES THE REGULATORY LEVEL.  
 10 THE OIL WAS TREATED AS A SOLID AND LEACHED WITH EXTRACTION FLUID  
 11 ADL(AVERAGE DETECTION LIMITS)  
 12 PQL(PRACTICAL QUANTITATION LIMITS)  
 13 SAMPLE ANALYZED OVER HOLDING TIME  
 14 DISSOLVED VALUE MAY BE HIGHER THAN TOTAL DUE TO CONTAMINATION FROM  
 15 THE FILTERING PROCEDURE  
 16 SAMPLED BY ULI  
 17 DISSOLVED VALUE MAY BE HIGHER THAN TOTAL; HOWEVER, THE VALUES ARE  
 18 WITHIN EXPERIMENTAL ERROR  
 19 AN INHIBITORY FACTOR WAS OBSERVED IN THIS ANALYSIS  
 20 PARAMETER NOT ANALYZED WITHIN 15 MINUTES OF SAMPLING  
 21 THE SERIAL DILUTION OF THIS SAMPLE SUGGESTS A POSSIBLE PHYSICAL AND/OR CHEMICAL  
 22 INTERFERENCE IN THIS DETERMINATION. THE DATA MAY BE BIASED EITHER HIGH OR LOW.  
 23 CALCULATION BASED ON DRY WEIGHT  
 24 INDICATES AN ESTIMATED VALUE, DETECTED BUT BELOW THE PRACTICAL QUANTITATION  
 25 LIMITS  
 26 UG/KG AS REC.D / UG/KG DRY WT  
 27 MG/KG AS REC.D / MG/KG DRY WT  
 28 INSUFFICIENT SAMPLE PRECLUDES LOWER DETECTION LIMITS  
 29 SAMPLE DILUTED/BLANK CORRECTED  
 30 ND(NON-DETECTED)  
 31 MATRIX INTERFERENCE PRECLUDES LOWER DETECTION LIMITS/BLANK CORRECTED  
 32 SPIKE RECOVERY ABNORMALLY HIGH/LOW DUE TO MATRIX INTERFERENCE  
 33 POST-DIGESTION SPIKE FOR FURNACE AA ANALYSIS IS OUTSIDE OF THE CONTROL  
 34 LIMITS (85-115%); HOWEVER, THE SAMPLE CONCENTRATION IS BELOW THE PQL  
 35 ANALYZED BY METHOD OF STANDARD ADDITIONS  
 36 METHOD PERFORMANCE STUDY HAS NOT BEEN COMPLETED/ND(NON-DETECTED)  
 37 FIELD MEASURED PARAMETER TAKEN BY CLIENT  
 38 TARGET ANALYTE IS BIODEGRADED AND/OR ENVIRONMENTALLY WEATHERED  
 39 NON-POTABLE WATER SOURCE  
 40 VOLATILE ASP CODES  
 -----  
 (B) POSSIBLE/PROBABLE BLANK CONTAMINATION (D) ALL COMPOUNDS IDENTIFIED  
 AT A SECONDARY DILUTION FACTOR (J) ESTIMATED VALUE  
 35 THE HYDROCARBONS DETECTED IN THE SAMPLE DID NOT CROSS-MATCH WITH COMMON  
 PETROLEUM DISTILLATES  
 36 MATRIX INTERFERENCE CAUSING SPIKES TO RESULT IN LESS THAN 50.0% RECOVERY  
 37 MILLIGRAMS PER LITER (MG/L) / POUNDS (LBS) PER DAY  
 38 MILLIGRAMS PER LITER (MG/L) OF RESIDUAL CHLORINE (CL2) / POUNDS (LBS)  
 PER DAY OF CL2  
 39 MICROGRAMS PER LITER (UG/L) / POUNDS (LBS) PER DAY  
 40 MILLIGRAMS PER LITER (MG/L) LINEAR ALKYL SULFONATE (LAS) / POUNDS (LBS)  
 PER DAY LAS  
 41 RESULTS ARE REPORTED ON AN AS REC.D BASIS  
 42 THE SAMPLE WAS ANALYZED ON A TOTAL BASIS; THE TEST RESULT CAN BE COMPARED  
 TO THE TCLP REGULATORY CRITERIA BY DIVIDING THE TEST RESULT BY 20,  
 CREATING A THEORETICAL TCLP VALUE  
 43 METAL BY CONCENTRATION PROCEDURE  
 44 POSSIBLE CONTAMINATION FROM FIELD/LABORATORY

| PARAMETER      | CONF-7        | CONF-8        | CONF-9        | CONF-10     |
|----------------|---------------|---------------|---------------|-------------|
| Percent Solids | 59%           | 70%           | 66%           | 74%         |
| Aroclor 1016   | <0.14ug/kg dw | <0.11ug/kg dw | <0.12ug/kg dw | <54ug/kg dw |
| Aroclor 1221   | <0.14ug/kg dw | <0.11ug/kg dw | <0.12ug/kg dw | <54ug/kg dw |
| Aroclor 1232   | <0.14ug/kg dw | <0.11ug/kg dw | <0.12ug/kg dw | <54ug/kg dw |
| Aroclor 1242   | <0.14ug/kg dw | 1400ug/kg dw  | <0.12ug/kg dw | <54ug/kg dw |
| Aroclor 1248   | <0.14ug/kg dw | <0.11ug/kg dw | <0.12ug/kg dw | <54ug/kg dw |
| Aroclor 1254   | <0.14ug/kg dw | <0.11ug/kg dw | <0.12ug/kg dw | <54ug/kg dw |
| Aroclor 1260   | <0.14ug/kg dw | <0.11ug/kg dw | <0.12ug/kg dw | <54ug/kg dw |
| Total PCB      | <0.14ug/kg dw | 1400ug/kg dw  | <0.12ug/kg dw | <54ug/kg dw |



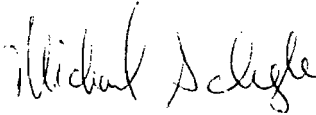


**ATTACHMENT G**

May 15, 2000

To: Kevin Matheis, OSC  
USEPA-Region II  
Lackawanna Foundry Site  
Lackawanna, New York

From: Mike Ingle, Field Chemist  
Earth Tech



Cc: Lane Aulick, Disposal Coordinator  
Vernon Wilson, Response Manager  
Mark Kromis, QA/QC Manager  
Earth Tech

Re: Suspect Arsenic Data at the Lackawanna Foundry Site

Various samples received from The Lackawanna Foundry Site at Upstate Laboratories on April 14, 2000 were analyzed for Trace Arsenic. One of these results was 14 mg/kg, which was near the site established action level of 20 mg/kg. The quality control (QC) parameters for the samples was low, therefore suggesting that the actual samples result might be biased high. After discussions with site T&D coordinator Lane Aulick, it was decided that a re-digestion and analysis of this sample, which is site sample designated as N20E0, would be appropriate to determine whether further site action is necessary for this sample designation. A sample aliquot was digested and analyzed on May 9, 2000 and produced results of 34 mg/kg and 27 mg/kg, respectively. Additional reruns of the digested aliquot on May 10, 2000 produced results of 34 mg/kg, 30 mg/kg, 26 mg/kg, respectively. After a review of the data with Mark Kromis, Earth Tech's QA/QC Manager, it was decided that the laboratory re-digest 3 separate sample aliquots in order to determine some statistical comparisons. These samples were analyzed on May 12, 2000 and produced results of 20 mg/kg, 27 mg/kg, and 11 mg/kg, respectively. Using all the results listed with the exception of the original value of 14 mg/kg, an average value of 26.1 mg/kg was calculated with 29.2 %RSD between all the numbers.

**ATTACHMENT H**

# Upstate Laboratories inc.

Shipping: 6034 Corporate Dr. • E. Syracuse, NY 13057-1017 • (315) 437-0255 • Fax (315) 437-1209

Mailing: Box 289 • Syracuse, NY 13206

Albany (518) 459-3134

Binghamton (607) 724-0478

Buffalo (716) 649-2533

Rochester (716) 436-9070

New Jersey (201) 703-1324

May 18, 2000

Ms. Lane Aulick  
Earth Tech  
7870 Villa Park Dr.  
Suite 400  
Richmond, VA 23228

Re: Analysis Report #12500017 - Lackawanna Foundry

Dear Ms. Aulick:

Please find enclosed the results for your samples which were picked up by ULI personnel on May 3, 2000.

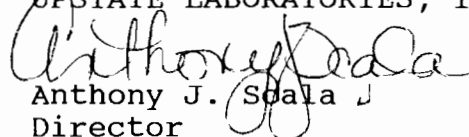
We have included the Chain of Custody Record as part of your report. You may need to reference this form for a more detailed explanation of your sample. Samples will be disposed of approximately one month from final report date.

Should you have any questions, please feel free to give us a call.

Thank you for your patronage.

Sincerely,

UPSTATE LABORATORIES, INC.

  
Anthony J. Scala  
Director

AJS/lw

Enclosures: report, spreadsheets, disk

cc/encs: N. Scala, ULI  
file

M. Kromis, Earth Tech (category B deliverables sent on  
5/10/00)

Note: Faxed results were given to your office and K. Matheis on  
5/5/00. AJS

**Disclaimer:** The test results and procedures utilized, and laboratory interpretations of data obtained by ULI as contained in this report are believed by ULI to be accurate and reliable for sample(s) tested. In accepting this report, the customer agrees that the full extent of any and all liability for actual and consequential damages of ULI for the services performed shall be equal to the fee charged to the customer for the services as liquidated damages.

DATE: 05/18/00

Upstate Laboratories, Inc.  
Analysis Results  
Report Number: 12500017  
Client I.D.: EARTH TECH

APPROVAL: *QSS*  
QC: *JD* Lab I.D.: 10170  
Sampled by: Client

ID:12500017 Mat:Soil LACKAWANNA FOUNDRY LW-SS-01 0940H 05/03/00 G

| PARAMETERS                        | RESULTS      | DATE ANAL. | KEY | FILE#  |
|-----------------------------------|--------------|------------|-----|--------|
| Percent Solids                    | 95%          | 05/04/00   |     | WD0102 |
| PCB (Aroclors) by EPA Method 8080 |              |            |     |        |
| Aroclor 1016                      | <1.7ug/kg dw | 05/04/00   |     | GA0057 |
| Aroclor 1221                      | <1.7ug/kg dw | 05/04/00   |     | GA0057 |
| Aroclor 1232                      | <1.7ug/kg dw | 05/04/00   |     | GA0057 |
| Aroclor 1242                      | 134ug/kg dw  | 05/04/00   |     | GA0057 |
| Aroclor 1248                      | <1.7ug/kg dw | 05/04/00   |     | GA0057 |
| Aroclor 1254                      | <1.7ug/kg dw | 05/04/00   |     | GA0057 |
| Aroclor 1260                      | <1.7ug/kg dw | 05/04/00   |     | GA0057 |
| Total PCB                         | 134ug/kg dw  | 05/04/00   |     | GA0057 |

ID:12500018 Mat:Soil LACKAWANNA FOUNDRY LW-SS-02 1000H 05/03/00 G

| PARAMETERS                        | RESULTS      | DATE ANAL. | KEY | FILE#  |
|-----------------------------------|--------------|------------|-----|--------|
| Percent Solids                    | 69%          | 05/04/00   |     | WD0102 |
| PCB (Aroclors) by EPA Method 8080 |              |            |     |        |
| Aroclor 1016                      | <170ug/kg dw | 05/05/00   | 05  | GA0057 |
| Aroclor 1221                      | <170ug/kg dw | 05/05/00   | 05  | GA0057 |
| Aroclor 1232                      | <170ug/kg dw | 05/05/00   | 05  | GA0057 |
| Aroclor 1242                      | 5800ug/kg dw | 05/05/00   |     | GA0057 |
| Aroclor 1248                      | <170ug/kg dw | 05/05/00   | 05  | GA0057 |
| Aroclor 1254                      | <170ug/kg dw | 05/05/00   | 05  | GA0057 |
| Aroclor 1260                      | <170ug/kg dw | 05/05/00   | 05  | GA0057 |
| Total PCB                         | 5800ug/kg dw | 05/05/00   |     | GA0057 |

ID:12500019 Mat:Soil LACKAWANNA FOUNDRY LW-SS-03 1005H 05/03/00 G

| PARAMETERS                        | RESULTS        | DATE ANAL. | KEY | FILE#  |
|-----------------------------------|----------------|------------|-----|--------|
| Percent Solids                    | 68%            | 05/04/00   |     | WD0102 |
| PCB (Aroclors) by EPA Method 8080 |                |            |     |        |
| Aroclor 1016                      | <170ug/kg dw   | 05/04/00   | 05  | GA0057 |
| Aroclor 1221                      | <170ug/kg dw   | 05/04/00   | 05  | GA0057 |
| Aroclor 1232                      | <170ug/kg dw   | 05/04/00   | 05  | GA0057 |
| Aroclor 1242                      | 13,000ug/kg dw | 05/04/00   |     | GA0057 |
| Aroclor 1248                      | <170ug/kg dw   | 05/04/00   | 05  | GA0057 |
| Aroclor 1254                      | <170ug/kg dw   | 05/04/00   | 05  | GA0057 |
| Aroclor 1260                      | <170ug/kg dw   | 05/04/00   | 05  | GA0057 |
| Total PCB                         | 13,000ug/kg dw | 05/04/00   |     | GA0057 |

iw = Dry weight

KEY PAGE

1 MATRIX INTERFERENCE PRECLUDES LOWER DETECTION LIMITS  
2 MATRIX INTERFERENCE  
3 PRESENT IN BLANK  
4 ANALYSIS NOT PERFORMED BECAUSE OF INSUFFICIENT SAMPLE  
5 THE PRESENCE OF OTHER TARGET ANALYTE(S) PRECLUDES LOWER DETECTION LIMITS  
6 BLANK CORRECTED  
7 HEAD SPACE PRESENT IN SAMPLE  
8 QUANTITATION LIMIT IS GREATER THAN THE CALCULATED REGULATORY LEVEL. THE  
9 QUANTITATION LIMIT THEREFORE BECOMES THE REGULATORY LEVEL.  
10 THE OIL WAS TREATED AS A SOLID AND LEACHED WITH EXTRACTION FLUID  
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12 PQL(PRACTICAL QUANTITATION LIMITS)  
13 SAMPLE ANALYZED OVER HOLDING TIME  
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15 THE FILTERING PROCEDURE  
16 SAMPLED BY ULI  
17 DISSOLVED VALUE MAY BE HIGHER THAN TOTAL; HOWEVER, THE VALUES ARE  
18 WITHIN EXPERIMENTAL ERROR  
19 AN INHIBITORY FACTOR WAS OBSERVED IN THIS ANALYSIS  
20 PARAMETER NOT ANALYZED WITHIN 15 MINUTES OF SAMPLING  
21 THE SERIAL DILUTION OF THIS SAMPLE SUGGESTS A POSSIBLE PHYSICAL AND/OR CHEMICAL  
22 INTERFERENT IN THIS DETERMINATION. THE DATA MAY BE BIASED EITHER HIGH OR LOW.  
23 CALCULATION BASED ON DRY WEIGHT  
24 INDICATES AN ESTIMATED VALUE, DETECTED BUT BELOW THE PRACTICAL QUANTITATION  
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29 SAMPLE DILUTED/BLANK CORRECTED  
30 ND(NON-DETECTED)  
31 MATRIX INTERFERENCE PRECLUDES LOWER DETECTION LIMITS/BLANK CORRECTED  
32 SPIKE RECOVERY ABNORMALLY HIGH/LOW DUE TO MATRIX INTERFERENCE  
33 POST-DIGESTION SPIKE FOR FURNACE AA ANALYSIS IS OUTSIDE OF THE CONTROL  
34 LIMITS (85-115%); HOWEVER, THE SAMPLE CONCENTRATION IS BELOW THE PQL  
35 ANALYZED BY METHOD OF STANDARD ADDITIONS  
36 METHOD PERFORMANCE STUDY HAS NOT BEEN COMPLETED/ND(NON-DETECTED)  
37 FIELD MEASURED PARAMETER TAKEN BY CLIENT  
38 TARGET ANALYTE IS BIODEGRADED AND/OR ENVIRONMENTALLY WEATHERED  
39 NON-POTABLE WATER SOURCE  
40 VOLATILE ASP CODES  
-----  
41 (B) POSSIBLE/PROBABLE BLANK CONTAMINATION (D) ALL COMPOUNDS IDENTIFIED AT A  
42 SECONDARY DILUTION FACTOR (J) DETECTED BELOW THE CRQL  
43 THE HYDROCARBONS DETECTED IN THE SAMPLE DID NOT CROSS-MATCH WITH COMMON  
44 PETROLEUM DISTILLATES  
45 MATRIX INTERFERENCE CAUSING SPIKES TO RESULT IN LESS THAN 50.0% RECOVERY  
46 MILLIGRAMS PER LITER (MG/L) / POUNDS (LBS) PER DAY  
47 MILLIGRAMS PER LITER (MG/L) OF RESIDUAL CHLORINE (CL2) / POUNDS (LBS)  
48 PER DAY OF CL2  
49 MICROGRAMS PER LITER (UG/L) / POUNDS (LBS) PER DAY  
50 MILLIGRAMS PER LITER (MG/L) LINEAR ALKYL SULFONATE (LAS) / POUNDS (LBS)  
51 PER DAY LAS  
52 RESULTS ARE REPORTED ON AN AS REC.D BASIS  
53 THE SAMPLE WAS ANALYZED ON A TOTAL BASIS; THE TEST RESULT CAN BE COMPARED  
54 TO THE TCLP REGULATORY CRITERIA BY DIVIDING THE TEST RESULT BY 20,  
55 CREATING A THEORETICAL TCLP VALUE  
56 METAL BY CONCENTRATION PROCEDURE  
57 POSSIBLE CONTAMINATION FROM FIELD/LABORATORY

| PARAMETER      | LW-SS-01     | LW-SS-02     | LW-SS-03       |
|----------------|--------------|--------------|----------------|
| Percent Solids | 95%          | 69%          | 68%            |
| Aroclor 1016   | <1.7ug/kg dw | <170ug/kg dw | <170ug/kg dw   |
| Aroclor 1221   | <1.7ug/kg dw | <170ug/kg dw | <170ug/kg dw   |
| Aroclor 1232   | <1.7ug/kg dw | <170ug/kg dw | <170ug/kg dw   |
| Aroclor 1242   | 134ug/kg dw  | 5800ug/kg dw | 13,000ug/kg dw |
| Aroclor 1248   | <1.7ug/kg dw | <170ug/kg dw | <170ug/kg dw   |
| Aroclor 1254   | <1.7ug/kg dw | <170ug/kg dw | <170ug/kg dw   |
| Aroclor 1260   | <1.7ug/kg dw | <170ug/kg dw | <170ug/kg dw   |
| Total PCB      | 134ug/kg dw  | 5800ug/kg dw | 13,000ug/kg dw |

ACE NO. \_\_\_\_\_  
 PO No. \_\_\_\_\_



**SUPERFUND TECHNICAL ASSESSMENT AND RESPONSE TEAM**  
 EPA CONTRACT 68-W5-0019  
 Phone: 908-225-6116 Fax: 908-225-7037

- MATRIX BOX NO.:
1. Surface Water
  2. Ground Water
  3. Leachate
  4. Rinse
  5. Soil/Sediment
  6. Oil
  7. Waste
  8. Other (Specify)

- PRESERVATIVE BOX NO.:
1. HCl
  2. HNO3
  3. Na2SO4
  4. H2SO4
  5. Other (Specify)
  6. Ice Only
  7. Not Preserved
- 5/5  
HDD

Send verbal and written results to: Roy F. Weston, Inc., USEPA Region II START  
 Suite 201, 1090 King Georges Post Road, Edison, New Jersey 08837-3703  
 Attention: Smita Sumbaly, START Analytical Coordinator

| Sample Number | Sample Collection<br>MM/DD/YY/Time | Sample Matrix<br>(Enter box #) | Conc. Low-L<br>Med-M<br>High-H | Sample Type<br>Comp-C<br>Grab-G | Sample Preserv.<br>(Enter box #) | RAS ANALYSIS |     |      |      |     |     | BCRA ANALYSIS |     |      | Full<br>TCLP<br>%<br>SPL/CS | OTHER |                |
|---------------|------------------------------------|--------------------------------|--------------------------------|---------------------------------|----------------------------------|--------------|-----|------|------|-----|-----|---------------|-----|------|-----------------------------|-------|----------------|
|               |                                    |                                |                                |                                 |                                  | VOL          | HNA | PEST | PCBs | TAL | END | MIN           | CON | REAC |                             |       |                |
| LW-SS-01      | 5/03/00 0940                       | 5                              | L                              | G                               |                                  |              |     |      |      |     |     |               |     |      |                             | (x)   | 24 hr TAT      |
| LW-SS-02      | 5/03/00 1000                       | 5                              | L                              | G                               |                                  |              |     |      |      |     |     |               |     |      |                             | (x)   |                |
| LW-SS-03      | 5/03/00 1005                       | 5                              | L                              | G                               |                                  |              |     |      |      |     |     |               |     |      |                             | (x)   | Fax Results    |
|               |                                    |                                |                                |                                 |                                  |              |     |      |      |     |     |               |     |      |                             |       | To             |
|               |                                    |                                |                                |                                 |                                  |              |     |      |      |     |     |               |     |      |                             |       | Kevin Matheis  |
|               |                                    |                                |                                |                                 |                                  |              |     |      |      |     |     |               |     |      |                             |       | @ 716-827-3036 |
|               |                                    |                                |                                |                                 |                                  |              |     |      |      |     |     |               |     |      |                             |       | Mail Hard copy |
|               |                                    |                                |                                |                                 |                                  |              |     |      |      |     |     |               |     |      |                             |       | To USEPA       |
|               |                                    |                                |                                |                                 |                                  |              |     |      |      |     |     |               |     |      |                             |       | Kevin Matheis  |
|               |                                    |                                |                                |                                 |                                  |              |     |      |      |     |     |               |     |      |                             |       | 3 Elm St       |
|               |                                    |                                |                                |                                 |                                  |              |     |      |      |     |     |               |     |      |                             |       | LACKAWANNA, NY |
|               |                                    |                                |                                |                                 |                                  |              |     |      |      |     |     |               |     |      |                             |       | 14218          |

Person Assuming Responsibility for Sample: DAVID L ADAMS Time: 0940 Date (MM/DD/YY): 4/3/00

Sample Number: All Relinquished By: David L Adams Time: 1030 Date: 4/3/00 Received By: [Signature] Reason for Change of Custody: Shipment to Lab.

Sample Number: \_\_\_\_\_ Relinquished By: [Signature] Time: \_\_\_\_\_ Date: \_\_\_\_\_ Received By: [Signature] Reason for Change of Custody: \_\_\_\_\_

Sample Number: \_\_\_\_\_ Relinquished By: [Signature] Time: 1730 Date: 5/3/00 Received By: [Signature] Reason for Change of Custody: \_\_\_\_\_



**ATTACHMENT I**

# Upstate Laboratories inc.

Shipping: 6034 Corporate Dr. • E. Syracuse, NY 13057-1017 • (315) 437-0255 • Fax (315) 437-1209

Mailing: Box 289 • Syracuse, NY 13206

Albany (518) 459-3134

Binghamton (607) 724-0478

Buffalo (716) 649-2533

Rochester (716) 436-9070

New Jersey (201) 703-1324

May 18, 2000

Ms. Lane Aulick  
Earth Tech  
7870 Villa Park Dr.  
Suite 400  
Richmond, VA 23228

1st Topsoil  
sample

Re: Analysis Report #13100037 - 38340 Lackawanna Foundry

Dear Ms. Aulick:

Please find enclosed the results for your sample which was received on May 10, 2000.

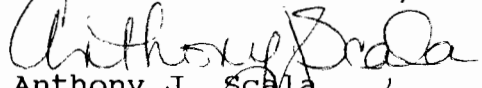
We have included the Chain of Custody Record as part of your report. You may need to reference this form for a more detailed explanation of your sample. Samples will be disposed of approximately one month from final report date.

Should you have any questions, please feel free to give us a call.

Thank you for your patronage.

Sincerely,

UPSTATE LABORATORIES, INC.

  
Anthony J. Scala  
Director

AJS/lw

Enclosures: report, spreadsheets, disk

cc/encs: N. Scala, ULI  
file

M. Kromis, Earth Tech (category B deliverables to follow)

Note: Faxed results were given to your office and K. Matheis on 5/16 and 5/17/00. AJS

Disclaimer: The test results and procedures utilized, and laboratory interpretations of data obtained by ULI as contained in this report are believed by ULI to be accurate and reliable for sample(s) tested. In accepting this report, the customer agrees that the full extent of any and all liability for actual and consequential damages of ULI for the services performed shall be equal to the fee charged to the customer for the services as liquidated damages.

DATE: 05/18/00

Upstate Laboratories, Inc.  
Analysis Results  
Report Number: 13100037  
Client I.D.: EARTH TECH  
Sampled by:

APPROVAL: *AS*  
QC: *AS*  
Lab I.D.: 10170

38340 LACKAWANNA  
FOUNDRY 1ST STA MAIN ENTRANCE 05/09/00 G

ULI I.D.: 13100037

Matrix: Soil

| PARAMETERS                       | RESULTS        | DATE ANAL. | KEY | FILE#  |
|----------------------------------|----------------|------------|-----|--------|
| Percent Solids                   | 87%            | 05/10/00   |     | WD0174 |
| Total Aluminum                   | 5400mg/kg dw   | 05/15/00   |     | ME2805 |
| Total Antimony                   | <2mg/kg dw     | 05/15/00   |     | ME2805 |
| Total Arsenic by furnace method  | <2mg/kg dw     | 05/15/00   |     | ME2805 |
| Total Barium                     | 56mg/kg dw     | 05/15/00   |     | ME2805 |
| Total Beryllium                  | <0.4mg/kg dw   | 05/15/00   |     | ME2805 |
| Total Cadmium                    | 1.3mg/kg dw    | 05/15/00   |     | ME2805 |
| Total Calcium                    | 2900mg/kg dw   | 05/15/00   |     | ME2805 |
| Total Chromium                   | 8.8mg/kg dw    | 05/15/00   |     | ME2805 |
| Total Cobalt                     | 3.0mg/kg dw    | 05/15/00   |     | ME2805 |
| Total Copper                     | 5.3mg/kg dw    | 05/15/00   |     | ME2805 |
| Total Iron                       | 10,000mg/kg dw | 05/15/00   |     | ME2805 |
| Total Lead                       | 8.8mg/kg dw    | 05/15/00   |     | ME2805 |
| Total Magnesium                  | 1200mg/kg dw   | 05/15/00   |     | ME2805 |
| Total Manganese                  | 93mg/kg dw     | 05/15/00   |     | ME2805 |
| Total Mercury                    | <0.3mg/kg dw   | 05/11/00   |     | MB2123 |
| Total Nickel                     | 7.8mg/kg dw    | 05/15/00   |     | ME2805 |
| Total Potassium                  | 490mg/kg dw    | 05/15/00   |     | ME2805 |
| Total Selenium by furnace method | <0.6mg/kg dw   | 05/16/00   |     | ME2807 |
| Total Silver                     | <2mg/kg dw     | 05/16/00   |     | ME2807 |
| Total Sodium                     | 230mg/kg dw    | 05/15/00   |     | ME2805 |
| Total Thallium by furnace method | <2mg/kg dw     | 05/15/00   |     | ME2805 |
| Total Vanadium                   | 14mg/kg dw     | 05/15/00   |     | ME2805 |
| Total Zinc                       | 35mg/kg dw     | 05/15/00   |     | ME2805 |

TCL Volatiles by EPA Method 8260

|                          |             |          |    |        |
|--------------------------|-------------|----------|----|--------|
| Chloromethane            | <3ug/kg dw  | 05/12/00 |    | VM2880 |
| Bromomethane             | <3ug/kg dw  | 05/12/00 |    | VM2880 |
| Vinyl Chloride           | <2ug/kg dw  | 05/12/00 |    | VM2880 |
| Chloroethane             | <3ug/kg dw  | 05/12/00 |    | VM2880 |
| Methylene Chloride       | 19ug/kg dw  | 05/12/00 | 44 | VM2880 |
| Acetone                  | <11ug/kg dw | 05/12/00 |    | VM2880 |
| Carbon Disulfide         | <3ug/kg dw  | 05/12/00 |    | VM2880 |
| 1,1-Dichloroethene       | <3ug/kg dw  | 05/12/00 |    | VM2880 |
| 1,1-Dichloroethane       | <3ug/kg dw  | 05/12/00 |    | VM2880 |
| trans-1,2-Dichloroethene | <3ug/kg dw  | 05/12/00 |    | VM2880 |
| cis-1,2-Dichloroethene   | <3ug/kg dw  | 05/12/00 |    | VM2880 |
| Chloroform               | <3ug/kg dw  | 05/12/00 |    | VM2880 |
| 1,2-Dichloroethane       | <3ug/kg dw  | 05/12/00 |    | VM2880 |
| 2-Butanone               | <11ug/kg dw | 05/12/00 |    | VM2880 |

dw = Dry weight

DATE: 05/18/00

Upstate Laboratories, Inc.  
Analysis Results  
Report Number: 13100037  
Client I.D.: EARTH TECH  
Sampled by:

APPROVAL: *AS*  
QC: *JD*  
Lab I.D.: 10170

38340 LACKAWANNA  
FOUNDRY 1ST STA MAIN ENTRANCE 05/09/00 G

ULI I.D.: 13100037

Matrix: Soil

| PARAMETERS                | RESULTS     | DATE ANAL. | KEY | FILE#  |
|---------------------------|-------------|------------|-----|--------|
| 1,1,1-Trichloroethane     | <3ug/kg dw  | 05/12/00   |     | VM2880 |
| Carbon Tetrachloride      | <3ug/kg dw  | 05/12/00   |     | VM2880 |
| Bromodichloromethane      | <3ug/kg dw  | 05/12/00   |     | VM2880 |
| 1,2-Dichloropropane       | <3ug/kg dw  | 05/12/00   |     | VM2880 |
| cis-1,3-Dichloropropene   | <3ug/kg dw  | 05/12/00   |     | VM2880 |
| Trichloroethene           | <3ug/kg dw  | 05/12/00   |     | VM2880 |
| Dibromochloromethane      | <3ug/kg dw  | 05/12/00   |     | VM2880 |
| 1,1,2-Trichloroethane     | <3ug/kg dw  | 05/12/00   |     | VM2880 |
| Benzene                   | <3ug/kg dw  | 05/12/00   |     | VM2880 |
| trans-1,3-Dichloropropene | <3ug/kg dw  | 05/12/00   |     | VM2880 |
| Bromoform                 | <3ug/kg dw  | 05/12/00   |     | VM2880 |
| 4-Methyl-2-pentanone      | <11ug/kg dw | 05/12/00   |     | VM2880 |
| 2-Hexanone                | <11ug/kg dw | 05/12/00   |     | VM2880 |
| Tetrachloroethene         | <3ug/kg dw  | 05/12/00   |     | VM2880 |
| 1,1,2,2-Tetrachloroethane | <3ug/kg dw  | 05/12/00   |     | VM2880 |
| Toluene                   | <3ug/kg dw  | 05/12/00   |     | VM2880 |
| Chlorobenzene             | <3ug/kg dw  | 05/12/00   |     | VM2880 |
| Ethylbenzene              | <3ug/kg dw  | 05/12/00   |     | VM2880 |
| Styrene                   | <3ug/kg dw  | 05/12/00   |     | VM2880 |
| m-Xylene and p-Xylene     | <3ug/kg dw  | 05/12/00   |     | VM2880 |
| o-Xylene                  | <3ug/kg dw  | 05/12/00   |     | VM2880 |

TCL Semivolatiles by EPA Method 8270

|                              |              |          |  |        |
|------------------------------|--------------|----------|--|--------|
| Phenol                       | <380ug/kg dw | 05/16/00 |  | SA2409 |
| bis(2-Chloroethyl) ether     | <380ug/kg dw | 05/16/00 |  | SA2409 |
| 2-Chlorophenol               | <380ug/kg dw | 05/16/00 |  | SA2409 |
| 1,3-Dichlorobenzene          | <380ug/kg dw | 05/16/00 |  | SA2409 |
| 1,4-Dichlorobenzene          | <380ug/kg dw | 05/16/00 |  | SA2409 |
| 1,2-Dichlorobenzene          | <380ug/kg dw | 05/16/00 |  | SA2409 |
| 2-Methylphenol               | <380ug/kg dw | 05/16/00 |  | SA2409 |
| 2,2'-Oxybis(1-Chloropropane) | <380ug/kg dw | 05/16/00 |  | SA2409 |
| 4-Methylphenol               | <380ug/kg dw | 05/16/00 |  | SA2409 |
| n-Nitrosodi-n-propylamine    | <380ug/kg dw | 05/16/00 |  | SA2409 |
| Hexachloroethane             | <380ug/kg dw | 05/16/00 |  | SA2409 |
| Nitrobenzene                 | <380ug/kg dw | 05/16/00 |  | SA2409 |
| Isophorone                   | <380ug/kg dw | 05/16/00 |  | SA2409 |
| 2-Nitrophenol                | <380ug/kg dw | 05/16/00 |  | SA2409 |
| 2,4-Dimethylphenol           | <380ug/kg dw | 05/16/00 |  | SA2409 |
| bis(2-Chloroethoxy)methane   | <380ug/kg dw | 05/16/00 |  | SA2409 |
| 2,4-Dichlorophenol           | <380ug/kg dw | 05/16/00 |  | SA2409 |

dw = Dry weight

DATE: 05/18/00

Upstate Laboratories, Inc.  
Analysis Results  
Report Number: 13100037  
Client I.D.: EARTH TECH  
Sampled by:

APPROVAL: *AS*  
QC: *JD*  
Lab I.D.: 10170

38340 LACKAWANNA  
FOUNDRY 1ST STA MAIN ENTRANCE 05/09/00 G

ULI I.D.: 13100037

Matrix: Soil

| PARAMETERS                 | RESULTS       | DATE ANAL. | KEY | FILE#  |
|----------------------------|---------------|------------|-----|--------|
| 1,2,4-Trichlorobenzene     | <380ug/kg dw  | 05/16/00   |     | SA2409 |
| Naphthalene                | <380ug/kg dw  | 05/16/00   |     | SA2409 |
| 4-Chloroaniline            | <380ug/kg dw  | 05/16/00   |     | SA2409 |
| Hexachlorobutadiene        | <380ug/kg dw  | 05/16/00   |     | SA2409 |
| 4-Chloro-3-methylphenol    | <380ug/kg dw  | 05/16/00   |     | SA2409 |
| 2-Methylnaphthalene        | <380ug/kg dw  | 05/16/00   |     | SA2409 |
| Hexachlorocyclopentadiene  | <380ug/kg dw  | 05/16/00   |     | SA2409 |
| 2,4,6-Trichlorophenol      | <380ug/kg dw  | 05/16/00   |     | SA2409 |
| 2,4,5-Trichlorophenol      | <380ug/kg dw  | 05/16/00   |     | SA2409 |
| 2-Chloronaphthalene        | <380ug/kg dw  | 05/16/00   |     | SA2409 |
| 2-Nitroaniline             | <3800ug/kg dw | 05/16/00   |     | SA2409 |
| Dimethylphthalate          | <380ug/kg dw  | 05/16/00   |     | SA2409 |
| Acenaphthylene             | <380ug/kg dw  | 05/16/00   |     | SA2409 |
| 2,6-Dinitrotoluene         | <380ug/kg dw  | 05/16/00   |     | SA2409 |
| 3-Nitroaniline             | <3800ug/kg dw | 05/16/00   |     | SA2409 |
| Acenaphthene               | <380ug/kg dw  | 05/16/00   |     | SA2409 |
| 2,4-Dinitrophenol          | <3800ug/kg dw | 05/16/00   |     | SA2409 |
| 4-Nitrophenol              | <3800ug/kg dw | 05/16/00   |     | SA2409 |
| Dibenzofuran               | <380ug/kg dw  | 05/16/00   |     | SA2409 |
| 2,4-Dinitrotoluene         | <380ug/kg dw  | 05/16/00   |     | SA2    |
| Diethylphthalate           | <380ug/kg dw  | 05/16/00   |     | SA2409 |
| 4-Chlorophenylphenylether  | <380ug/kg dw  | 05/16/00   |     | SA2409 |
| Fluorene                   | <380ug/kg dw  | 05/16/00   |     | SA2409 |
| 4-Nitroaniline             | <3800ug/kg dw | 05/16/00   |     | SA2409 |
| 2-Methyl-4,6-dinitrophenol | <3800ug/kg dw | 05/16/00   |     | SA2409 |
| n-Nitrosodiphenylamine     | <380ug/kg dw  | 05/16/00   |     | SA2409 |
| 4-Bromophenylphenylether   | <380ug/kg dw  | 05/16/00   |     | SA2409 |
| Hexachlorobenzene          | <380ug/kg dw  | 05/16/00   |     | SA2409 |
| Pentachlorophenol          | <760ug/kg dw  | 05/16/00   |     | SA2409 |
| Phenanthrene               | <380ug/kg dw  | 05/16/00   |     | SA2409 |
| Anthracene                 | <380ug/kg dw  | 05/16/00   |     | SA2409 |
| Carbazole                  | <380ug/kg dw  | 05/16/00   |     | SA2409 |
| di-n-butylphthalate        | <380ug/kg dw  | 05/16/00   |     | SA2409 |
| Fluoranthene               | <380ug/kg dw  | 05/16/00   |     | SA2409 |
| Pyrene                     | <380ug/kg dw  | 05/16/00   |     | SA2409 |
| Butylbenzylphthalate       | <380ug/kg dw  | 05/16/00   |     | SA2409 |
| 3,3'-Dichlorobenzidine     | <380ug/kg dw  | 05/16/00   |     | SA2409 |
| Benzo(a)anthracene         | <380ug/kg dw  | 05/16/00   |     | SA2409 |
| Chrysene                   | <380ug/kg dw  | 05/16/00   |     | SA2409 |
| bis(2-Ethylhexyl)phthalate | <380ug/kg dw  | 05/16/00   |     | SA2409 |

dw = Dry weight

DATE: 05/18/00

Upstate Laboratories, Inc.

Analysis Results

Report Number: 13100037

Client I.D.: EARTH TECH

Sampled by:

APPROVAL: *AS*

QC: *JD*

Lab I.D.: 10170

38340 LACKAWANNA

FOUNDRY 1ST STA MAIN ENTRANCE 05/09/00 G

ULI I.D.: 13100037

Matrix: Soil

| PARAMETERS             | RESULTS      | DATE ANAL. | KEY | FILE#  |
|------------------------|--------------|------------|-----|--------|
| di-n-octylphthalate    | <380ug/kg dw | 05/16/00   |     | SA2409 |
| Benzo(b)fluoranthene   | <380ug/kg dw | 05/16/00   |     | SA2409 |
| Benzo(k)fluoranthene   | <380ug/kg dw | 05/16/00   |     | SA2409 |
| Benzo(a)pyrene         | <380ug/kg dw | 05/16/00   |     | SA2409 |
| Indeno(1,2,3-cd)pyrene | <380ug/kg dw | 05/16/00   |     | SA2409 |
| Dibenzo(a,h)anthracene | <380ug/kg dw | 05/16/00   |     | SA2409 |
| Benzo(ghi)perylene     | <380ug/kg dw | 05/16/00   |     | SA2409 |

TCL Pesticides by EPA Method 8080

|                    |              |          |  |        |
|--------------------|--------------|----------|--|--------|
| BHC (a-isomer)     | <2.0ug/kg dw | 05/17/00 |  | GA0077 |
| BHC (b-isomer)     | <2.0ug/kg dw | 05/17/00 |  | GA0077 |
| BHC (d-isomer)     | <2.0ug/kg dw | 05/17/00 |  | GA0077 |
| BHC (g-isomer)     | <2.0ug/kg dw | 05/17/00 |  | GA0077 |
| Heptachlor         | <2.0ug/kg dw | 05/17/00 |  | GA0077 |
| Aldrin             | <2.0ug/kg dw | 05/17/00 |  | GA0077 |
| Heptachlor Epoxide | <2.0ug/kg dw | 05/17/00 |  | GA0077 |
| Endosulfan I       | <2.0ug/kg dw | 05/17/00 |  | GA0077 |
| Dieldrin           | <3.8ug/kg dw | 05/17/00 |  | GA0077 |
| 4,4'-DDE           | <3.8ug/kg dw | 05/17/00 |  | GA0077 |
| Endrin             | <3.8ug/kg dw | 05/17/00 |  | GA0077 |
| Endosulfan II      | <3.8ug/kg dw | 05/17/00 |  | GA0077 |
| 4,4'-DDD           | <3.8ug/kg dw | 05/17/00 |  | GA0077 |
| Endosulfan Sulfate | <3.8ug/kg dw | 05/17/00 |  | GA0077 |
| 4,4'-DDT           | <3.8ug/kg dw | 05/17/00 |  | GA0077 |
| Methoxychlor       | <20ug/kg dw  | 05/17/00 |  | GA0077 |
| Endrin Ketone      | <3.8ug/kg dw | 05/17/00 |  | GA0077 |
| Endrin Aldehyde    | <3.8ug/kg dw | 05/17/00 |  | GA0077 |
| alpha-Chlordane    | <2.0ug/kg dw | 05/17/00 |  | GA0077 |
| gamma-Chlordane    | <2.0ug/kg dw | 05/17/00 |  | GA0077 |
| Toxaphene          | <195ug/kg dw | 05/17/00 |  | GA0077 |

dw = Dry weight

KEY PAGE

1 MATRIX INTERFERENCE PRECLUDES LOWER DETECTION LIMITS  
2 MATRIX INTERFERENCE  
3 PRESENT IN BLANK  
4 ANALYSIS NOT PERFORMED BECAUSE OF INSUFFICIENT SAMPLE  
5 THE PRESENCE OF OTHER TARGET ANALYTE(S) PRECLUDES LOWER DETECTION LIMITS  
6 BLANK CORRECTED  
7 HEAD SPACE PRESENT IN SAMPLE  
8 QUANTITATION LIMIT IS GREATER THAN THE CALCULATED REGULATORY LEVEL. THE  
9 QUANTITATION LIMIT THEREFORE BECOMES THE REGULATORY LEVEL.  
10 THE OIL WAS TREATED AS A SOLID AND LEACHED WITH EXTRACTION FLUID  
11 ADL(AVERAGE DETECTION LIMITS)  
12 PQL(PRACTICAL QUANTITATION LIMITS)  
13 SAMPLE ANALYZED OVER HOLDING TIME  
14 DISSOLVED VALUE MAY BE HIGHER THAN TOTAL DUE TO CONTAMINATION FROM  
15 THE FILTERING PROCEDURE  
16 SAMPLED BY ULI  
17 DISSOLVED VALUE MAY BE HIGHER THAN TOTAL; HOWEVER, THE VALUES ARE  
18 WITHIN EXPERIMENTAL ERROR  
19 AN INHIBITORY FACTOR WAS OBSERVED IN THIS ANALYSIS  
20 PARAMETER NOT ANALYZED WITHIN 15 MINUTES OF SAMPLING  
21 THE SERIAL DILUTION OF THIS SAMPLE SUGGESTS A POSSIBLE PHYSICAL AND/OR CHEMICAL  
22 INTERFERENT IN THIS DETERMINATION. THE DATA MAY BE BIASED EITHER HIGH OR LOW.  
23 CALCULATION BASED ON DRY WEIGHT  
24 INDICATES AN ESTIMATED VALUE, DETECTED BUT BELOW THE PRACTICAL QUANTITATION  
25 LIMITS  
26 UG/KG AS REC.D / UG/KG DRY WT  
27 MG/KG AS REC.D / MG/KG DRY WT  
28 INSUFFICIENT SAMPLE PRECLUDES LOWER DETECTION LIMITS  
29 SAMPLE DILUTED/BLANK CORRECTED  
30 ND(NON-DETECTED)  
31 MATRIX INTERFERENCE PRECLUDES LOWER DETECTION LIMITS/BLANK CORRECTED  
32 SPIKE RECOVERY ABNORMALLY HIGH/LOW DUE TO MATRIX INTERFERENCE  
33 POST-DIGESTION SPIKE FOR FURNACE AA ANALYSIS IS OUTSIDE OF THE CONTROL  
34 LIMITS (85-115%); HOWEVER, THE SAMPLE CONCENTRATION IS BELOW THE PQL  
35 ANALYZED BY METHOD OF STANDARD ADDITIONS  
36 METHOD PERFORMANCE STUDY HAS NOT BEEN COMPLETED/ND(NON-DETECTED)  
37 FIELD MEASURED PARAMETER TAKEN BY CLIENT  
38 TARGET ANALYTE IS BIODEGRADED AND/OR ENVIRONMENTALLY WEATHERED  
39 NON-POTABLE WATER SOURCE  
40 VOLATILE ASP CODES  
-----  
41 (B)POSSIBLE/PROBABLE BLANK CONTAMINATION (D)ALL COMPOUNDS IDENTIFIED AT A  
42 SECONDARY DILUTION FACTOR (J)DETECTED BELOW THE CRQL  
43 THE HYDROCARBONS DETECTED IN THE SAMPLE DID NOT CROSS-MATCH WITH COMMON  
44 PETROLEUM DISTILLATES  
45 MATRIX INTERFERENCE CAUSING SPIKES TO RESULT IN LESS THAN 50.0% RECOVERY  
46 MILLIGRAMS PER LITER (MG/L) / POUNDS (LBS) PER DAY  
47 MILLIGRAMS PER LITER (MG/L) OF RESIDUAL CHLORINE (CL2) / POUNDS (LBS)  
48 PER DAY OF CL2  
49 MICROGRAMS PER LITER (UG/L) / POUNDS (LBS) PER DAY  
50 MILLIGRAMS PER LITER (MG/L) LINEAR ALKYL SULFONATE (LAS) / POUNDS (LBS)  
51 PER DAY LAS  
52 RESULTS ARE REPORTED ON AN AS REC.D BASIS  
53 THE SAMPLE WAS ANALYZED ON A TOTAL BASIS; THE TEST RESULT CAN BE COMPARED  
54 TO THE TCLP REGULATORY CRITERIA BY DIVIDING THE TEST RESULT BY 20,  
55 CREATING A THEORETICAL TCLP VALUE  
56 METAL BY CONCENTRATION PROCEDURE  
57 POSSIBLE CONTAMINATION FROM FIELD/LABORATORY

1st Station Main Entrance

| PARAMETER                        | 1st Station Main Entrance |
|----------------------------------|---------------------------|
| Percent Solids                   | 87%                       |
| Total Aluminum                   | 5400mg/kg dw              |
| Total Antimony                   | <2mg/kg dw                |
| Total Arsenic by furnace method  | <2mg/kg dw                |
| Total Barium                     | 56mg/kg dw                |
| Total Beryllium                  | <0.4mg/kg dw              |
| Total Cadmium                    | 1.3mg/kg dw               |
| Total Calcium                    | 2900mg/kg dw              |
| Total Chromium                   | 8.8mg/kg dw               |
| Total Cobalt                     | 3.0mg/kg dw               |
| Total Copper                     | 5.3mg/kg dw               |
| Total Iron                       | 10,000mg/kg dw            |
| Total Lead                       | 8.8mg/kg dw               |
| Total Magnesium                  | 1200mg/kg dw              |
| Total Manganese                  | 93mg/kg dw                |
| Total Mercury                    | <0.3mg/kg dw              |
| Total Nickel                     | 7.8mg/kg dw               |
| Total Potassium                  | 490mg/kg dw               |
| Total Selenium by furnace method | <0.6mg/kg dw              |
| Total Silver                     | <2mg/kg dw                |
| Total Sodium                     | 230mg/kg dw               |
| Total Thallium by furnace method | <2mg/kg dw                |
| Total Vanadium                   | 14mg/kg dw                |
| Total Zinc                       | 35mg/kg dw                |
| Chloromethane                    | <3ug/kg dw                |
| Bromomethane                     | <3ug/kg dw                |
| Vinyl Chloride                   | <2ug/kg dw                |
| Chloroethane                     | <3ug/kg dw                |
| Methylene Chloride               | 19ug/kg dw                |
| Acetone                          | <11ug/kg dw               |
| Carbon Disulfide                 | <3ug/kg dw                |
| 1,1 – Dichloroethene             | <3ug/kg dw                |
| 1,1 – Dichloroethane             | <3ug/kg dw                |
| trans – 1,2 – Dichloroethene     | <3ug/kg dw                |
| cis – 1,2 – Dichloroethene       | <3ug/kg dw                |
| Chloroform                       | <3ug/kg dw                |
| 1,2 – Dichloroethane             | <3ug/kg dw                |



1st Station Main Entrance

| PARAMETER                    | 1st Station Main Entrance |
|------------------------------|---------------------------|
| 2-Butanone                   | < 11ug/kg dw              |
| 1,1,1-Trichloroethane        | < 3ug/kg dw               |
| Carbon Tetrachloride         | < 3ug/kg dw               |
| Bromodichloromethane         | < 3ug/kg dw               |
| 1,2-Dichloropropane          | < 3ug/kg dw               |
| cis-1,3-Dichloropropene      | < 3ug/kg dw               |
| Trichloroethene              | < 3ug/kg dw               |
| Dibromochloromethane         | < 3ug/kg dw               |
| 1,1,2-Trichloroethane        | < 3ug/kg dw               |
| Benzene                      | < 3ug/kg dw               |
| trans-1,3-Dichloropropene    | < 3ug/kg dw               |
| Bromoform                    | < 3ug/kg dw               |
| 4-Methyl-2-pentanone         | < 11ug/kg dw              |
| 2-Hexanone                   | < 11ug/kg dw              |
| Tetrachloroethene            | < 3ug/kg dw               |
| 1,1,2,2-Tetrachloroethane    | < 3ug/kg dw               |
| Toluene                      | < 3ug/kg dw               |
| Chlorobenzene                | < 3ug/kg dw               |
| Ethylbenzene                 | < 3ug/kg dw               |
| Styrene                      | < 3ug/kg dw               |
| m-Xylene and p-Xylene        | < 3ug/kg dw               |
| o-Xylene                     | < 3ug/kg dw               |
| Phenol                       | < 380ug/kg dw             |
| bis(2-Chloroethyl)ether      | < 380ug/kg dw             |
| 2-Chlorophenol               | < 380ug/kg dw             |
| 1,3-Dichlorobenzene          | < 380ug/kg dw             |
| 1,4-Dichlorobenzene          | < 380ug/kg dw             |
| 1,2-Dichlorobenzene          | < 380ug/kg dw             |
| 2-Methylphenol               | < 380ug/kg dw             |
| 2,2'-Oxybis(1-Chloropropane) | < 380ug/kg dw             |
| 4-Methylphenol               | < 380ug/kg dw             |
| n-Nitrosodi-n-propylamine    | < 380ug/kg dw             |
| Hexachloroethane             | < 380ug/kg dw             |
| Nitrobenzene                 | < 380ug/kg dw             |
| Isophorone                   | < 380ug/kg dw             |
| 2-Nitrophenol                | < 380ug/kg dw             |
| 2,4-Diethylphenol            | < 380ug/kg dw             |

1st Station Main Entrance

| PARAMETER                        |                |
|----------------------------------|----------------|
| bis(2--Chloroethoxy)methane      | < 380ug/kg dw  |
| 2,4 - Dichlorophenol             | < 380ug/kg dw  |
| 1,2,4 - Trichlorobenzene         | < 380ug/kg dw  |
| Naphthalene                      | < 380ug/kg dw  |
| 4 - Chloroaniline                | < 380ug/kg dw  |
| Hexachlorobutadiene              | < 380ug/kg dw  |
| 4 - Chloro - 3 - methylphenol    | < 380ug/kg dw  |
| 2 - Methylnaphthalene            | < 380ug/kg dw  |
| Hexachlorocyclopentadiene        | < 380ug/kg dw  |
| 2,4,6 - Trichlorophenol          | < 380ug/kg dw  |
| 2,4,5 - Trichlorophenol          | < 380ug/kg dw  |
| 2 - Chloronaphthalene            | < 380ug/kg dw  |
| 2 - Nitroaniline                 | < 3800ug/kg dw |
| Dimethylphthalate                | < 380ug/kg dw  |
| Acenaphthylene                   | < 380ug/kg dw  |
| 2,6 - Dinitrotoluene             | < 380ug/kg dw  |
| 3 - Nitroaniline                 | < 3800ug/kg dw |
| Acenaphthene                     | < 380ug/kg dw  |
| 2,4 - Dinitrophenol              | < 3800ug/kg dw |
| 4 - Nitrophenol                  | < 3800ug/kg dw |
| Dibenzofuran                     | < 380ug/kg dw  |
| 2,4 - Dinitrotoluene             | < 380ug/kg dw  |
| Diethylphthalate                 | < 380ug/kg dw  |
| 4 - Chlorophenylphenylether      | < 380ug/kg dw  |
| Fluorene                         | < 380ug/kg dw  |
| 4 - Nitroaniline                 | < 3800ug/kg dw |
| 2 - Methyl - 4,6 - dinitrophenol | < 3800ug/kg dw |
| n - Nitrosodiphenylamine         | < 380ug/kg dw  |
| 4 - Bromophenylphenylether       | < 380ug/kg dw  |
| Hexachlorobenzene                | < 380ug/kg dw  |
| Pentachlorophenol                | < 760ug/kg dw  |
| Phenanthrene                     | < 380ug/kg dw  |
| Anthracene                       | < 380ug/kg dw  |
| Carbazole                        | < 380ug/kg dw  |
| di - n - butylphthalate          | < 380ug/kg dw  |
| Fluoranthene                     | < 380ug/kg dw  |
| Pyrene                           | < 380ug/kg dw  |

1st Station Main Entrance

| PARAMETER                  |              |
|----------------------------|--------------|
| Butylbenzylphthalate       | <380ug/kg dw |
| 3,3'-Dichlorobenzidine     | <380ug/kg dw |
| Benzo(a)anthracene         | <380ug/kg dw |
| Chrysene                   | <380ug/kg dw |
| bis(2-Ethylhexyl)phthalate | <380ug/kg dw |
| di-n-octylphthalate        | <380ug/kg dw |
| Benzo(b)fluoranthene       | <380ug/kg dw |
| Benzo(k)fluoranthene       | <380ug/kg dw |
| Benzo(a)pyrene             | <380ug/kg dw |
| Indeno(1,2,3-cd)pyrene     | <380ug/kg dw |
| Dibenzo(a,h)anthracene     | <380ug/kg dw |
| Benzo(ghi)perylene         | <380ug/kg dw |
| BHC (a-isomer)             | <2.0ug/kg dw |
| BHC (b-isomer)             | <2.0ug/kg dw |
| BHC (d-isomer)             | <2.0ug/kg dw |
| BHC (g-isomer)             | <2.0ug/kg dw |
| Heptachlor                 | <2.0ug/kg dw |
| Aldrin                     | <2.0ug/kg dw |
| Heptachlor Epoxide         | <2.0ug/kg dw |
| Endosulfan I               | <2.0ug/kg dw |
| Dieldrin                   | <3.8ug/kg dw |
| 4,4'-DDE                   | <3.8ug/kg dw |
| Endrin                     | <3.8ug/kg dw |
| Endosulfan II              | <3.8ug/kg dw |
| 4,4'-DDD                   | <3.8ug/kg dw |
| Endosulfan Sulfate         | <3.8ug/kg dw |
| 4,4'-DDT                   | <3.8ug/kg dw |
| Methoxychlor               | <20ug/kg dw  |
| Endrin Ketone              | <3.8ug/kg dw |
| Endrin Aldehyde            | <3.8ug/kg dw |
| alpha-Chlordane            | <2.0ug/kg dw |
| gamma-Chlordane            | <2.0ug/kg dw |
| Toxaphene                  | <195ug/kg dw |



# Upstate Laboratories inc.

Shipping: 6034 Corporate Dr. • E. Syracuse, NY 13057-1017 • (315) 437-0255 • Fax (315) 437-1209

Mailing: Box 289 • Syracuse, NY 13206

Albany (518) 459-3134

Binghamton (607) 724-0478

Buffalo (716) 649-2533

Rochester (716) 436-9070

New Jersey (201) 703-1324

June 28, 2000

Ms. Lane Aulick  
Earth Tech  
7870 Villa Park Dr.  
Suite 400  
Richmond, VA 23228

2nd Backfill/  
Topsoil sample

Re: Analysis Report #16400197 - 38340 Lackawanna Foundry

Dear Ms. Aulick:

Please find enclosed the results for your sample which was picked up by ULI personnel on June 9, 2000.

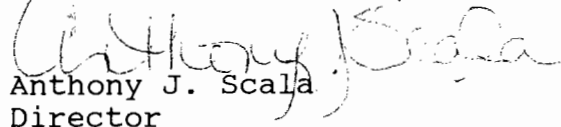
We have included the Chain of Custody Record as part of your report. You may need to reference this form for a more detailed explanation of your sample. Samples will be disposed of approximately one month from final report date.

Should you have any questions, please feel free to give us a call.

Thank you for your patronage.

Sincerely,

UPSTATE LABORATORIES, INC.

  
Anthony J. Scala  
Director

AJS/lw

Enclosures: report, spreadsheets, disk

cc/encs: N. Scala, ULI  
file

✓ K. Matheis, USEPA-Region 2 (copy report, copy spreadsheets)  
M. Kromis, Earth Tech (data deliverables to follow)

Note: Faxed results were given to your office and K. Matheis, USEPA-Region 2 previous to this report. AJS

Disclaimer: The test results and procedures utilized, and laboratory interpretations of data obtained by ULI as contained in this report are believed by ULI to be accurate and reliable for sample(s) tested. In accepting this report, the customer agrees that the full extent of any and all liability for actual and consequential damages of ULI for the services performed shall be equal to the fee charged to the customer for the services as liquidated damages.

DATE: 06/28/00

Upstate Laboratories, Inc.  
Analysis Results  
Report Number: 16400197  
Client I.D.: EARTH TECH  
Sampled by:

APPROVAL: *AS*

QC: *JS*

Lab I.D.: 10170

38340 LACKAWANNA  
FOUNDRY PARISOS TOP SOIL BACK 1430H 05/24/00 C

ULI I.D.: 16400197

Matrix: Soil

| PARAMETERS                       | RESULTS       | DATE ANAL. | KEY | FILE#  |
|----------------------------------|---------------|------------|-----|--------|
| Percent Solids                   | 85%           | 06/13/00   |     | WD0579 |
| Total Aluminum                   | 5000mg/kg dw  | 06/20/00   |     | ME2883 |
| Total Antimony                   | <36mg/kg dw   | 06/20/00   |     | ME2883 |
| Total Arsenic by furnace method  | <1.2mg/kg dw  | 06/20/00   |     | ME2883 |
| Total Barium                     | 60mg/kg dw    | 06/20/00   |     | ME2883 |
| Total Beryllium                  | <0.59mg/kg dw | 06/20/00   |     | ME2883 |
| Total Cadmium                    | 0.74mg/kg dw  | 06/20/00   |     | ME2883 |
| Total Calcium                    | 5100mg/kg dw  | 06/20/00   |     | ME2883 |
| Total Chromium                   | 7.3mg/kg dw   | 06/20/00   |     | ME2883 |
| Total Cobalt                     | <5.9mg/kg dw  | 06/20/00   |     | ME2883 |
| Total Copper                     | 6.9mg/kg dw   | 06/20/00   |     | ME2883 |
| Total Iron                       | 6900mg/kg dw  | 06/20/00   |     | ME2883 |
| Total Lead                       | <12mg/kg dw   | 06/20/00   |     | ME2883 |
| Total Magnesium                  | 1800mg/kg dw  | 06/20/00   |     | ME2883 |
| Total Manganese                  | 120mg/kg dw   | 06/20/00   |     | ME2883 |
| Total Mercury                    | <0.3mg/kg dw  | 06/20/00   |     | ME2494 |
| Total Nickel                     | 7.6mg/kg dw   | 06/20/00   |     | ME2883 |
| Total Potassium                  | 730mg/kg dw   | 06/20/00   |     | ME2883 |
| Total Selenium by furnace method | 5.7mg/kg dw   | 06/20/00   |     | ME2883 |
| Total Silver                     | <6.0mg/kg dw  | 06/20/00   |     | ME2883 |
| Total Sodium                     | 240mg/kg dw   | 06/20/00   |     | ME2883 |
| Total Thallium by furnace method | <0.35mg/kg dw | 06/20/00   |     | ME2883 |
| Total Vanadium                   | <36mg/kg dw   | 06/20/00   |     | ME2883 |
| Total Zinc                       | 38mg/kg dw    | 06/20/00   |     | ME2883 |

PCB (Aroclors) by EPA Method 8080

|              |              |          |  |        |
|--------------|--------------|----------|--|--------|
| Aroclor 1016 | <2.0ug/kg dw | 06/13/00 |  | GA0136 |
| Aroclor 1221 | <2.0ug/kg dw | 06/13/00 |  | GA0136 |
| Aroclor 1232 | <2.0ug/kg dw | 06/13/00 |  | GA0136 |
| Aroclor 1242 | <2.0ug/kg dw | 06/13/00 |  | GA0136 |
| Aroclor 1248 | <2.0ug/kg dw | 06/13/00 |  | GA0136 |
| Aroclor 1254 | <2.0ug/kg dw | 06/13/00 |  | GA0136 |
| Aroclor 1260 | <2.0ug/kg dw | 06/13/00 |  | GA0136 |
| Total PCB    | <2.0ug/kg dw | 06/13/00 |  | GA0136 |

TCL Pesticides by EPA Method 8080

|                |              |          |  |        |
|----------------|--------------|----------|--|--------|
| BHC (a-isomer) | <2.0ug/kg dw | 06/13/00 |  | GA0136 |
| BHC (b-isomer) | <2.0ug/kg dw | 06/13/00 |  | GA0136 |
| BHC (d-isomer) | <2.0ug/kg dw | 06/13/00 |  | GA0136 |
| BHC (g-isomer) | <2.0ug/kg dw | 06/13/00 |  | GA0136 |

dw = Dry weight

DATE: 06/28/00

Upstate Laboratories, Inc.  
Analysis Results  
Report Number: 16400197  
Client I.D.: EARTH TECH  
Sampled by:

APPROVAL: AS  
QC: SD  
Lab I.D.: 10170

38340 LACKAWANNA  
FOUNDRY PARISOS TOP SOIL BACK 1430H 05/24/00 C

ULI I.D.: 16400197

Matrix: Soil

| PARAMETERS         | RESULTS      | DATE ANAL. | KEY | FILE#  |
|--------------------|--------------|------------|-----|--------|
| Heptachlor         | <2.0ug/kg dw | 06/13/00   |     | GA0136 |
| Aldrin             | <2.0ug/kg dw | 06/13/00   |     | GA0136 |
| Heptachlor Epoxide | <2.0ug/kg dw | 06/13/00   |     | GA0136 |
| Endosulfan I       | <2.0ug/kg dw | 06/13/00   |     | GA0136 |
| Dieldrin           | <3.9ug/kg dw | 06/13/00   |     | GA0136 |
| 4,4'-DDE           | 8.9ug/kg dw  | 06/13/00   |     | GA0136 |
| Endrin             | <3.9ug/kg dw | 06/13/00   |     | GA0136 |
| Endosulfan II      | <3.9ug/kg dw | 06/13/00   |     | GA0136 |
| 4,4'-DDD           | <3.9ug/kg dw | 06/13/00   |     | GA0136 |
| Endosulfan Sulfate | <3.9ug/kg dw | 06/13/00   |     | GA0136 |
| 4,4'-DDT           | 6.4ug/kg dw  | 06/13/00   |     | GA0136 |
| Methoxychlor       | <20ug/kg dw  | 06/13/00   |     | GA0136 |
| Endrin Ketone      | <3.9ug/kg dw | 06/13/00   |     | GA0136 |
| Endrin Aldehyde    | <3.9ug/kg dw | 06/13/00   |     | GA0136 |
| alpha-Chlordane    | <2.0ug/kg dw | 06/13/00   |     | GA0136 |
| gamma-Chlordane    | <2.0ug/kg dw | 06/13/00   |     | GA0136 |
| Toxaphene          | <200ug/kg dw | 06/13/00   |     | GA0136 |

dw = Dry weight

| PARAMETER                        | TS - 001 Parisos Top Soil Backup<br>85% |
|----------------------------------|---|
| Percent Solids                   |   |
| Total Aluminum                   | 5000mg/kg dw                            |
| Total Antimony                   | <36mg/kg dw                             |
| Total Arsenic by furnace method  | <1.2mg/kg dw                            |
| Total Barium                     | 60mg/kg dw                              |
| Total Beryllium                  | <0.59mg/kg dw                           |
| Total Cadmium                    | 0.74mg/kg dw                            |
| Total Calcium                    | 5100mg/kg dw                            |
| Total Chromium                   | 7.3mg/kg dw                             |
| Total Cobalt                     | <5.9mg/kg dw                            |
| Total Copper                     | 6.9mg/kg dw                             |
| Total Iron                       | 6900mg/kg dw                            |
| Total Lead                       | <12mg/kg dw                             |
| Total Magnesium                  | 1800mg/kg dw                            |
| Total Manganese                  | 120mg/kg dw                             |
| Total Mercury                    | <0.3mg/kg dw                            |
| Total Nickel                     | 7.6mg/kg dw                             |
| Total Potassium                  | 730mg/kg dw                             |
| Total Selenium by furnace method | 5.7mg/kg dw                             |
| Total Silver                     | <6.0mg/kg dw                            |
| Total Sodium                     | 240mg/kg dw                             |
| Total Thallium by furnace method | <0.35mg/kg dw                           |
| Total Vanadium                   | <36mg/kg dw                             |
| Total Zinc                       | 38mg/kg dw                              |
| Aroclor 1016                     | <2.0ug/kg dw                            |
| Aroclor 1221                     | <2.0ug/kg dw                            |
| Aroclor 1232                     | <2.0ug/kg dw                            |
| Aroclor 1242                     | <2.0ug/kg dw                            |
| Aroclor 1248                     | <2.0ug/kg dw                            |
| Aroclor 1254                     | <2.0ug/kg dw                            |
| Aroclor 1260                     | <2.0ug/kg dw                            |
| Total PCB                        | <2.0ug/kg dw                            |
| BHC (a-isomer)                   | <2.0ug/kg dw                            |
| BHC (b-isomer)                   | <2.0ug/kg dw                            |
| BHC (d-isomer)                   | <2.0ug/kg dw                            |
| BHC (g-isomer)                   | <2.0ug/kg dw                            |
| Heptachlor                       | <2.0ug/kg dw                            |



| PARAMETER          | TS-001 Parisos Top Soil Backup |
|--------------------|--------------------------------|
| Aldrin             | < 2.0ug/kg dw                  |
| Heptachlor Epoxide | < 2.0ug/kg dw                  |
| Endosulfan I       | < 2.0ug/kg dw                  |
| Dieldrin           | < 3.9ug/kg dw                  |
| 4,4'-DDE           | 8.9ug/kg dw                    |
| Endrin             | < 3.9ug/kg dw                  |
| Endosulfan II      | < 3.9ug/kg dw                  |
| 4,4'-DDD           | < 3.9ug/kg dw                  |
| Endosulfan Sulfate | < 3.9ug/kg dw                  |
| 4,4'-DDT           | 6.4ug/kg dw                    |
| Methoxychlor       | < 20ug/kg dw                   |
| Endrin Ketone      | < 3.9ug/kg dw                  |
| Endrin Aldehyde    | < 3.9ug/kg dw                  |
| alpha-Chlordane    | < 2.0ug/kg dw                  |
| gamma-Chlordane    | < 2.0ug/kg dw                  |
| Toxaphene          | < 200ug/kg dw                  |



**ATTACHMENT J**

# The Haseley Companies

10315 LOCKPORT ROAD  
 P.O. BOX 212 L.P.O.  
 NIAGARA FALLS N.Y. 14304

## LETTER OF TRANSMITTAL

|                              |                           |          |      |
|------------------------------|---------------------------|----------|------|
| DATE:                        | 13-May-00                 | JOB NO.: | 9901 |
| ATTENTION: Mr. Vernon Wilson |                           |          |      |
| RE:                          | Gratwick - Riverside Park |          |      |
|                              | Hydric Soil               |          |      |
|                              |                           |          |      |
|                              |                           |          |      |

TO Earth Tech, Inc.  
 3 Elm Street  
 Lackawana, N.Y. 14218

- HASELEY TRUCKING CO.
- HASELEY CONSTRUCTION CO.
- HASELEY CONSULTANTS/CONSTRUCTORS

GENTLEMEN:

WE ARE SENDING YOU  Attached  Under separate cover via \_\_\_\_\_ the following items

- Shop drawings
- Copy of letter
- Prints
- Change order
- Plans
- Reports
- Samples
- Permits
- Specifications
- Test Results

| COPIES | DATE | NO. | DESCRIPTION                                 |
|--------|------|-----|---|
| 1      |      |     | ANALYTICAL TEST DATA (BUCKHORN HYDRIC SOIL) |
|        |      |     |   |
|        |      |     |   |
|        |      |     |   |

THESE ARE TRANSMITTED as checked below

- For Approval
- For your use
- As requested
- For review
- Approved as submitted
- Approved as noted
- Returned for corrections
- Resubmit \_\_\_\_\_ copies for approval
- Submit \_\_\_\_\_ copies for distribution
- Return \_\_\_\_\_ corrected prints

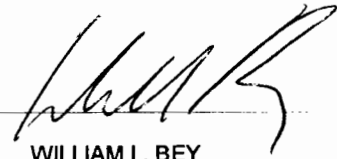
FOR BIDS DUE \_\_\_\_\_ 19\_\_\_\_

REMARKS: \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

PLEASE CONTACT ME @ 694-4755 IF YOU HAVE ANY QUESTIONS

COPY TO: JOB BOOK / TRANSLTR.

SIGNED:



WILLIAM L. BEY  
 PROJECT ENGINEER

Chemical Analysis  
HYDRIC SOIL  
(I PA TEST)



**Severn Trent Laboratories**  
10 Hazelwood Drive  
Suite 106  
Amherst, New York 14228-2298

July 31, 1999

Tel: (716) 691-2600  
Fax: (716) 691-7991

Mr. David Birch  
GZA GeoEnvironmental  
364 Nagel Drive  
Buffalo, New York 14225

RE: Analytical Results

Dear Mr. Birch:

Enclosed are analytical results concerning the samples recently submitted. The pertinent information regarding these analyses is listed below:

Quote #: NY99-237  
Project: HYDRIC SOIL Samples  
Matrix: Soil; Water  
Samples Received: 07/02/99  
Sample Date: 06/02/99

If you have any questions concerning this data, please contact me at (716) 691-2600 and refer to the I.D. number listed below. It has been our pleasure to provide GZA GeoEnvironmental with environmental testing services. We look forward to serving you in the future.

Sincerely,

Severn Trent Laboratories, Inc

Kenneth P. Kinecki  
Program Manager

| DOCUMENT APPROVAL / REVIEW          |                         |
|-------------------------------------|-------------------------|
| <input checked="" type="checkbox"/> | APPROVED                |
| <input type="checkbox"/>            | APPROVED AS NOTED       |
| <input type="checkbox"/>            | NOT APPROVED            |
| <input type="checkbox"/>            | REVISE AND RESUBMIT     |
| <input type="checkbox"/>            | REVIEWED FOR COMPLETION |
| BY                                  |                         |
| DATE                                | 10/22/99                |

KPK/rtv  
Enclosure

I.D. #A99-4287  
#NY9A8465

This report contains 99 pages which are individually numbered.

**Laboratory Locations:**

- Monroe, CT
- Pensacola, FL
- University Park, IL
- Billerica, MA
- Westfield, MA
- Edison, NJ
- Whippany, NJ
- Newburgh, NY
- Houston, TX

**Service Center Locations:**

- Mt. Laurel, NJ
- Glen Cove, NY

**Sales Office Locations:**

- Cantonment, FL
- New Orleans, LA
- Waterford, MI
- Blairstown, NJ
- Schenectady, NY
- Cleveland, OH

a part of

Severn Trent Services Inc



000001

## ANALYTICAL RESULTS

Prepared For:

GZA GeoEnvironmental  
364 Nagel Drive  
Buffalo, New York 14225

Prepared By:

Severn Trent Laboratories, Inc.  
10 Hazelwood Drive, Suite 106  
Amherst, New York 14228-2298

## METHODOLOGY

The specific methodology employed in obtaining the enclosed analytical results is indicated on the specific data table. The method number presented refers to the following U.S. Environmental Protection Agency reference:

- "Test Methods For Evaluating Solid Waste, Physical/Chemical Methods" (SW-846), Third Edition, September 1994, U.S. Environmental Protection Agency Office of Solid Waste.

## COMMENTS

Comments pertain to data on one or all pages of this report.

The enclosed data has been reported utilizing data qualifiers (Q) as defined on the Organic and Inorganic Data Comment Pages.

## METHOD 8260 DATA

The VBLK19 exhibited positive results for Methylene chloride. Affected samples are flagged with "B" qualifiers

No other deviations from protocol were encountered during the analytical procedures.

## METHOD 8270 DATA

The Method Blank exhibited positive results for Bis(2-ethylhexyl) phthalate. Affected samples are flagged with "B" qualifiers

Sample 07029-1A was analyzed at a dilution factor of 5 due to matrix interference.

No other deviations from protocol were encountered during the analytical procedures.



000002

METHOD 8081 DATA

Sample 07029-1A was analyzed at a dilution factor of 4 due to high levels of target compounds.

No other deviations from protocol were encountered during the analytical procedures.

METHOD 8082 DATA

No deviations from protocol were encountered during the analytical procedures.

METALS DATA

The Method Blank (A9B0622401) exhibited results for Mercury. However, all sample results were greater than ten (10) times that of the Method Blank.

The Method Blank (A9B0631702) exhibited results for Lead, Calcium, Manganese, and Zinc. However, all sample results were greater than ten (10) times that of the Method Blank.

Sample 07029-1A exhibited spike recovery results outside quality control limits for Aluminum, Antimony, Beryllium, Cadmium, Chromium, Cobalt, Copper, Lead, Manganese, Nickel, Selenium, Silver, Thallium, Vanadium, and Zinc. The samples were redigested and reanalyzed, and exhibited similar spike recoveries. This suggests matrix interference.

No other deviations from protocol were encountered during the analytical procedures.

WET CHEMISTRY DATA

No deviations from protocol were encountered during the analytical procedures.

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**ORGANIC DATA COMMENT PAGE**

Laboratory Name: SEVERN TRENT LABORATORIES INC.

USEPA Defined Organic Data Qualifiers:

- U - Indicates compound was analyzed for but not detected.
- J - Indicates an estimate value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed, or when the mass spectral data indicate the presence of a compound that meets the identification criteria but the result is less than the sample quantitation limit but greater than zero.
- C - This flag applies to pesticide results where the identification has been confirmed by GC/MS.
- B - This flag is used when the analyte is found in the associated blank as well as in the sample.
- E - This flag identifies compounds whose concentrations exceed the calibration range of the GC/MS instrument for that specific analysis.
- D - This flag identifies all compounds identified in an analysis at a secondary dilution factor.
- T - This flag is used when the analyte is found in the associated TCLP extraction blank as well as in the sample.
- N - Indicates presumptive evidence of a compound. This flag is only used for tentatively identified compounds, where the identification is based on a mass spectral library search. It is applied to all TIC results.
- P - This flag is used for a pesticide/Aroclor target analyte when there is greater than 25% difference for detected concentrations between the two GC columns. The lower of the two values is reported on Form I and flagged with a "P".
- A - This flag indicates that a TIC is a suspected aldol-condensation product.



**INORGANIC DATA COMMENT PAGE**

Laboratory Name: SEVERN TRENT LABORATORIES, INC.

**USEPA Defined Inorganic Data Qualifiers:**

- B - Indicates a value greater than or equal to the instrument detection limit, but less than the contract required detection limit.
- U - Indicates compound was analyzed for but not detected. Report with the detection limit value (e.g., 100).
- N - Indicates spike sample recovery is not within the control limits.
- K - Indicates the post digestion spike recovery is not within the control limits.
- \* - Indicates duplicate analysis is not within the control limits.
- S - Indicates value determined by the Method of Standard Addition.
- + - Indicates the correlation coefficient for the Method of Standard Addition is less than 0.995.
- M - Indicates duplicate injection results exceeded control limits.
- W - Post digestion spike for Furnace AA analysis is out of control limits (85-115%), while sample absorbance is less than 50% of spike absorbance.
- E - Indicates a value estimated or not reported due to the presence of interference.

000005

## Sample Data Package

Date: 07/31/1999

GZA Geoenvironmental

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Time: 10:40:23

Rept: AN1178

HYDRIX SOIL Samples

00000

Sample ID: 07029-1A

Date Received: 07/02/1999

Lab ID: A9428701

Project No: NY9A8465

Date Collected: 07/02/1999

Client No: L11085

Time Collected: 16:00

Site No:

| Parameter                  | Result | Flag | Detection |       | Method | Date/Time |        |     |
|----------------------------|--------|------|-----------|-------|--------|-----------|--------|-----|
|                            |        |      | Limit     | Units |        | Analized  | Analys |     |
| Naphthalene                | ND     |      | 390       | UG/KG | 8270   | 07/21/99  | 13:49  | RCS |
| 2-Nitroaniline             | ND     |      | 1600      | UG/KG | 8270   | 07/21/99  | 13:49  | RCS |
| 3-Nitroaniline             | ND     |      | 1600      | UG/KG | 8270   | 07/21/99  | 13:49  | RCS |
| 4-Nitroaniline             | ND     |      | 1600      | UG/KG | 8270   | 07/21/99  | 13:49  | RCS |
| Nitrobenzene               | ND     |      | 390       | UG/KG | 8270   | 07/21/99  | 13:49  | RCS |
| 2-Nitrophenol              | ND     |      | 780       | UG/KG | 8270   | 07/21/99  | 13:49  | RCS |
| 4-Nitrophenol              | ND     |      | 1600      | UG/KG | 8270   | 07/21/99  | 13:49  | RCS |
| N-nitrosodiphenylamine     | ND     |      | 390       | UG/KG | 8270   | 07/21/99  | 13:49  | RCS |
| N-Nitroso-Di-n-propylamine | ND     |      | 390       | UG/KG | 8270   | 07/21/99  | 13:49  | RCS |
| Pentachlorophenol          | ND     |      | 1600      | UG/KG | 8270   | 07/21/99  | 13:49  | RCS |
| Phenanthrene               | ND     |      | 390       | UG/KG | 8270   | 07/21/99  | 13:49  | RCS |
| Phenol                     | ND     |      | 390       | UG/KG | 8270   | 07/21/99  | 13:49  | RCS |
| Pyrene                     | ND     |      | 390       | UG/KG | 8270   | 07/21/99  | 13:49  | RCS |
| 1,2,4-Trichlorobenzene     | ND     |      | 390       | UG/KG | 8270   | 07/21/99  | 13:49  | RCS |
| 2,4,5-Trichlorophenol      | ND     |      | 800       | UG/KG | 8270   | 07/21/99  | 13:49  | RCS |
| 2,4,6-Trichlorophenol      | ND     |      | 780       | UG/KG | 8270   | 07/21/99  | 13:49  | RCS |

## SOIL-SW8463 8260 - TCL VOLATILES

|                            |    |   |    |       |      |          |       |  |
|----------------------------|----|---|----|-------|------|----------|-------|--|
| Acetone                    | ND |   | 10 | UG/KG | 8260 | 07/08/99 | 17:24 |  |
| Benzene                    | ND |   | 5  | UG/KG | 8260 | 07/08/99 | 17:24 |  |
| Bromodichloromethane       | ND |   | 5  | UG/KG | 8260 | 07/08/99 | 17:24 |  |
| Bromoform                  | ND |   | 5  | UG/KG | 8260 | 07/08/99 | 17:24 |  |
| Bromomethane               | ND |   | 10 | UG/KG | 8260 | 07/08/99 | 17:24 |  |
| 2-Butanone                 | ND |   | 10 | UG/KG | 8260 | 07/08/99 | 17:24 |  |
| Carbon Disulfide           | ND |   | 5  | UG/KG | 8260 | 07/08/99 | 17:24 |  |
| Carbon Tetrachloride       | ND |   | 5  | UG/KG | 8260 | 07/08/99 | 17:24 |  |
| Chlorobenzene              | ND |   | 5  | UG/KG | 8260 | 07/08/99 | 17:24 |  |
| Chloroethane               | ND |   | 10 | UG/KG | 8260 | 07/08/99 | 17:24 |  |
| Chloroform                 | ND |   | 5  | UG/KG | 8260 | 07/08/99 | 17:24 |  |
| Chloromethane              | ND |   | 10 | UG/KG | 8260 | 07/08/99 | 17:24 |  |
| Dibromochloromethane       | ND |   | 5  | UG/KG | 8260 | 07/08/99 | 17:24 |  |
| 1,1-Dichloroethane         | ND |   | 5  | UG/KG | 8260 | 07/08/99 | 17:24 |  |
| 1,2-Dichloroethane         | ND |   | 5  | UG/KG | 8260 | 07/08/99 | 17:24 |  |
| 1,1-Dichloroethene         | ND |   | 5  | UG/KG | 8260 | 07/08/99 | 17:24 |  |
| 1,2-Dichloroethene (Total) | ND |   | 10 | UG/KG | 8260 | 07/08/99 | 17:24 |  |
| 1,2-Dichloropropane        | ND |   | 5  | UG/KG | 8260 | 07/08/99 | 17:24 |  |
| cis-1,3-Dichloropropene    | ND |   | 5  | UG/KG | 8260 | 07/08/99 | 17:24 |  |
| trans-1,3-Dichloropropene  | ND |   | 5  | UG/KG | 8260 | 07/08/99 | 17:24 |  |
| Ethylbenzene               | ND |   | 5  | UG/KG | 8260 | 07/08/99 | 17:24 |  |
| 2-Hexanone                 | ND |   | 10 | UG/KG | 8260 | 07/08/99 | 17:24 |  |
| Methylene chloride         | 17 | B | 5  | UG/KG | 8260 | 07/08/99 | 17:24 |  |
| 4-Methyl-2-pentanone       | ND |   | 10 | UG/KG | 8260 | 07/08/99 | 17:24 |  |
| Styrene                    | ND |   | 5  | UG/KG | 8260 | 07/08/99 | 17:24 |  |
| 1,1,2,2-Tetrachloroethane  | ND |   | 5  | UG/KG | 8260 | 07/08/99 | 17:24 |  |
| Tetrachloroethene          | ND |   | 5  | UG/KG | 8260 | 07/08/99 | 17:24 |  |
| Toluene                    | ND |   | 5  | UG/KG | 8260 | 07/08/99 | 17:24 |  |
| 1,1,1-Trichloroethane      | ND |   | 5  | UG/KG | 8260 | 07/08/99 | 17:24 |  |
| 1,1,2-Trichloroethane      | ND |   | 5  | UG/KG | 8260 | 07/08/99 | 17:24 |  |
| Trichloroethene            | ND |   | 5  | UG/KG | 8260 | 07/08/99 | 17:24 |  |
| Vinyl acetate              | ND |   | 10 | UG/KG | 8260 | 07/08/99 | 17:24 |  |

Date: 07/31/1999  
 Time: 10:40:23

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HYDRIX SOIL Samples

000008

Sample ID: 07029-1A  
 Lab ID: A9428701

Date Collected: 07/02/1999  
 Time Collected: 16:00

Date Received: 07/02/1999

Project No: NY9A8465

Client No: L11085

Site No:

| Parameter                         | Result | Flag | Detection |       |        | Date/Time |       | Analy: |
|-----------------------------------|--------|------|-----------|-------|--------|-----------|-------|--------|
|                                   |        |      | Limit     | Units | Method | Analyzed  |       |        |
| Vinyl chloride                    | ND     |      | 10        | UG/KG | 8260   | 07/08/99  | 17:24 |        |
| Total Xylenes                     | ND     |      | 15        | UG/KG | 8260   | 07/08/99  | 17:24 |        |
| SOIL-SW8463 8082 - PCBS           |        |      |           |       |        |           |       |        |
| Aroclor 1016                      | ND     |      | 40        | UG/KG | 8082   | 07/10/99  | 12:01 | BJ     |
| Aroclor 1221                      | ND     |      | 80        | UG/KG | 8082   | 07/10/99  | 12:01 | BJ     |
| Aroclor 1232                      | ND     |      | 40        | UG/KG | 8082   | 07/10/99  | 12:01 | BJ     |
| Aroclor 1242                      | ND     |      | 40        | UG/KG | 8082   | 07/10/99  | 12:01 | BJ     |
| Aroclor 1248                      | ND     |      | 40        | UG/KG | 8082   | 07/10/99  | 12:01 | BJ     |
| Aroclor 1254                      | 4.0    | J    | 40        | UG/KG | 8082   | 07/10/99  | 12:01 | BJ     |
| Aroclor 1260                      | ND     |      | 40        | UG/KG | 8082   | 07/10/99  | 12:01 | BJ     |
| SOIL-SW8463 8081 - TCL PESTICIDES |        |      |           |       |        |           |       |        |
| Aldrin                            | ND     |      | 8.0       | UG/KG | 8081   | 07/23/99  | 06:43 | KEA    |
| alpha-BHC                         | ND     |      | 8.0       | UG/KG | 8081   | 07/23/99  | 06:43 | KEA    |
| beta-BHC                          | ND     |      | 8.0       | UG/KG | 8081   | 07/23/99  | 06:43 | KEA    |
| gamma-BHC (Lindane)               | ND     |      | 8.0       | UG/KG | 8081   | 07/23/99  | 06:43 | KEA    |
| delta-BHC                         | ND     |      | 8.0       | UG/KG | 8081   | 07/23/99  | 06:43 | KEA    |
| Chlordane                         | ND     |      | 80        | UG/KG | 8081   | 07/23/99  | 06:43 | KEA    |
| 4,4'-DDD                          | ND     |      | 16        | UG/KG | 8081   | 07/23/99  | 06:43 | KEA    |
| 4,4'-DDE                          | ND     |      | 16        | UG/KG | 8081   | 07/23/99  | 06:43 | KEA    |
| 4,4'-DDT                          | ND     |      | 16        | UG/KG | 8081   | 07/23/99  | 06:43 | KEA    |
| Dieldrin                          | ND     |      | 16        | UG/KG | 8081   | 07/23/99  | 06:43 | KEA    |
| Endosulfan I                      | ND     |      | 16        | UG/KG | 8081   | 07/23/99  | 06:43 | KEA    |
| Endosulfan II                     | ND     |      | 16        | UG/KG | 8081   | 07/23/99  | 06:43 | KEA    |
| Endosulfan Sulfate                | ND     |      | 16        | UG/KG | 8081   | 07/23/99  | 06:43 | KEA    |
| Endrin                            | ND     |      | 16        | UG/KG | 8081   | 07/23/99  | 06:43 | KEA    |
| Endrin aldehyde                   | ND     |      | 32        | UG/KG | 8081   | 07/23/99  | 06:43 | KEA    |
| Heptachlor                        | ND     |      | 8.0       | UG/KG | 8081   | 07/23/99  | 06:43 | KEA    |
| Heptachlor epoxide                | ND     |      | 8.0       | UG/KG | 8081   | 07/23/99  | 06:43 | KEA    |
| Methoxychlor                      | ND     |      | 80        | UG/KG | 8081   | 07/23/99  | 06:43 | KEA    |
| Toxaphene                         | ND     |      | 160       | UG/KG | 8081   | 07/23/99  | 06:43 | KEA    |
| Metals Analysis                   |        |      |           |       |        |           |       |        |
| Aluminum - Total                  | 4620   |      | 2.3       | MG/KG | 6010   | 07/16/99  | 18:53 | JL     |
| Antimony - Total                  | ND     |      | 0.47      | MG/KG | 6010   | 07/16/99  | 18:53 | JL     |
| Arsenic - Total                   | 3.1    |      | 0.29      | MG/KG | 6010   | 07/16/99  | 18:53 | JL     |
| Barium - Total                    | 23.0   |      | 0.12      | MG/KG | 6010   | 07/16/99  | 18:53 | JL     |
| Beryllium - Total                 | 0.42   |      | 0.12      | MG/KG | 6010   | 07/16/99  | 18:53 | JL     |
| Cadmium - Total                   | 0.093  |      | 0.058     | MG/KG | 6010   | 07/16/99  | 18:53 | JL     |
| Calcium - Total                   | 10400  |      | 2.3       | MG/KG | 6010   | 07/16/99  | 18:53 | JL     |
| Chromium - Total                  | 6.6    |      | 0.14      | MG/KG | 6010   | 07/16/99  | 18:53 | JL     |
| Cobalt - Total                    | 4.9    |      | 0.12      | MG/KG | 6010   | 07/16/99  | 18:53 | JL     |
| Copper - Total                    | 11.8   |      | 0.17      | MG/KG | 6010   | 07/16/99  | 18:53 | JL     |
| Iron - Total                      | 9880   |      | 2.3       | MG/KG | 6010   | 07/16/99  | 18:53 | JL     |
| Lead - Total                      | 7.4    |      | 0.19      | MG/KG | 6010   | 07/16/99  | 18:53 | JL     |
| Magnesium - Total                 | 5770   |      | 2.3       | MG/KG | 6010   | 07/16/99  | 18:53 | JL     |
| Manganese - Total                 | 129    |      | 0.12      | MG/KG | 6010   | 07/16/99  | 18:53 | JL     |
| Mercury - Total                   | ND     |      | 0.040     | MG/KG | 7471   | 07/13/99  | 13:00 | M      |
| Nickel - Total                    | 12.3   |      | 0.15      | MG/KG | 6010   | 07/16/99  | 18:53 | JL     |

Date: 07/31/1999  
Time: 10:40:23

GZA Geoenvironmental

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Rep: AN1178

00000

HYDRIX SOIL Samples

Sample ID: 07029-1A  
Lab ID: A9428701  
Date Collected: 07/02/1999  
Time Collected: 16:00

Date Received: 07/02/1999  
Project No: NY9A8465  
Client No: L11085  
Site No:

| Parameter                                  | Result | Flag | Detection |         |           | Date/Time      |     | Analys |
|--|--------|------|-----------|---------|-----------|----------------|-----|--------|
|  |        |      | Limit     | Units   | Method    | Analyzed       |     |        |
| Potassium - Total                          | 691    |      | 17.5      | MG/KG   | 6010      | 07/16/99 18:53 | JWT |        |
| Selenium - Total                           | ND     |      | 0.47      | MG/KG   | 6010      | 07/16/99 18:53 | JWT |        |
| Silver - Total                             | ND     |      | 0.17      | MG/KG   | 6010      | 07/16/99 18:53 | JWT |        |
| Sodium - Total                             | 82.8   |      | 29.1      | MG/KG   | 6010      | 07/16/99 18:53 | JWT |        |
| Thallium - Total                           | 1.4    |      | 0.62      | MG/KG   | 6010      | 07/16/99 18:53 | JWT |        |
| Vanadium - Total                           | 9.0    |      | 0.14      | MG/KG   | 6010      | 07/16/99 18:53 | JWT |        |
| Zinc - Total                               | 41.6   |      | 0.12      | MG/KG   | 6010      | 07/16/99 18:53 | JWT |        |
| Wet Chemistry Analysis                     |        |      |           |         |           |                |     |        |
| Leachable Phosphorous                      | ND     |      | 0.20      | UG/G    | 365.2     | 07/13/99       | MA  |        |
| Leachable pH                               | 7.66   |      | 0.00000   | S.U.    | 9045      | 07/09/99       | TB  |        |
| Leachable Total Organic Carbon             | 718    |      | 200       | UG/G    | 9060      | 07/20/99       | PGE |        |
| SHAKE EXTRACTION OF SOLID WASTE WITH WATER | ND     |      | 0         | INVALID | D 3987-85 | 07/13/99       | KMS |        |

Date: 07/31/1999  
Time: 10:40:23

GZA GeoEnvironmental

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Rept: AN1170

HYDRIX SOIL Samples

000010

Sample ID: 07029-1A (L)  
Lab ID: A9428702  
Date Collected: 07/02/1999  
Time Collected: 16:00

Date Received: 07/02/1999  
Project No: NY9A8465  
Client No: L11085  
Site No:

| Parameter              | Result | Flag | Detection Limit | Units   | Method    | Date/Time Analyzed | Analysis |
|------------------------|--------|------|-----------------|---------|-----------|--------------------|----------|
| Wet Chemistry Analysis |        |      |                 |         |           |                    |          |
| Nitrate                | 2.6    |      | 0.050           | MG N/L  | 9200      | 07/14/99           | DPA      |
| Redox Potential        | 205    |      | 0               | M.VOLTS | D-1498-76 | 07/16/99           | PGE      |

000011

# **Chronology and QC Summary Package**

Date: 07/31/1999  
Time: 10:40:50

Rept: AN0326

HYDRIX SOIL Samples  
METHOD 8260 - TCL VOLATILE ORGANICS

| Client ID<br>Job No<br>Sample Date | Lab ID | VBLK19<br>A99-4287 | A9428703           | Reporting<br>Limit | Sample<br>Value    | Reporting<br>Limit | Sample<br>Value    | Reporting<br>Limit | Sample<br>Value    |
|------------------------------------|--------|--------------------|--------------------|--------------------|--------------------|--------------------|--------------------|--------------------|--------------------|
| Analyte                            | Units  | Sample<br>Value    | Reporting<br>Limit | Sample<br>Value    | Reporting<br>Limit | Sample<br>Value    | Reporting<br>Limit | Sample<br>Value    | Reporting<br>Limit |
| Acetone                            | UG/KG  | ND                 | 10                 | NA                 |                    | NA                 |                    | NA                 |                    |
| Benzene                            | UG/KG  | ND                 | 5                  | NA                 |                    | NA                 |                    | NA                 |                    |
| Bromodichloromethane               | UG/KG  | ND                 | 5                  | NA                 |                    | NA                 |                    | NA                 |                    |
| Bromoform                          | UG/KG  | ND                 | 5                  | NA                 |                    | NA                 |                    | NA                 |                    |
| Bromomethane                       | UG/KG  | ND                 | 10                 | NA                 |                    | NA                 |                    | NA                 |                    |
| 2-Butanone                         | UG/KG  | ND                 | 10                 | NA                 |                    | NA                 |                    | NA                 |                    |
| Carbon Disulfide                   | UG/KG  | ND                 | 5                  | NA                 |                    | NA                 |                    | NA                 |                    |
| Carbon Tetrachloride               | UG/KG  | ND                 | 5                  | NA                 |                    | NA                 |                    | NA                 |                    |
| Chlorobenzene                      | UG/KG  | ND                 | 5                  | NA                 |                    | NA                 |                    | NA                 |                    |
| Chloroethane                       | UG/KG  | ND                 | 10                 | NA                 |                    | NA                 |                    | NA                 |                    |
| Chloroform                         | UG/KG  | ND                 | 5                  | NA                 |                    | NA                 |                    | NA                 |                    |
| Chloromethane                      | UG/KG  | ND                 | 10                 | NA                 |                    | NA                 |                    | NA                 |                    |
| Dibromochloromethane               | UG/KG  | ND                 | 5                  | NA                 |                    | NA                 |                    | NA                 |                    |
| 1,1-Dichloroethane                 | UG/KG  | ND                 | 5                  | NA                 |                    | NA                 |                    | NA                 |                    |
| 1,2-Dichloroethane                 | UG/KG  | ND                 | 5                  | NA                 |                    | NA                 |                    | NA                 |                    |
| 1,1-Dichloroethene                 | UG/KG  | ND                 | 5                  | NA                 |                    | NA                 |                    | NA                 |                    |
| 1,2-Dichloroethene (Total)         | UG/KG  | ND                 | 10                 | NA                 |                    | NA                 |                    | NA                 |                    |
| 1,2-Dichloropropane                | UG/KG  | ND                 | 5                  | NA                 |                    | NA                 |                    | NA                 |                    |
| cis-1,3-Dichloropropene            | UG/KG  | ND                 | 5                  | NA                 |                    | NA                 |                    | NA                 |                    |
| trans-1,3-Dichloropropene          | UG/KG  | ND                 | 5                  | NA                 |                    | NA                 |                    | NA                 |                    |
| Ethylbenzene                       | UG/KG  | ND                 | 5                  | NA                 |                    | NA                 |                    | NA                 |                    |
| 2-Hexanone                         | UG/KG  | ND                 | 10                 | NA                 |                    | NA                 |                    | NA                 |                    |
| Methylene chloride                 | UG/KG  | ND                 | 5                  | NA                 |                    | NA                 |                    | NA                 |                    |
| 4-Methyl-2-pentanone               | UG/KG  | ND                 | 10                 | NA                 |                    | NA                 |                    | NA                 |                    |
| Styrene                            | UG/KG  | ND                 | 5                  | NA                 |                    | NA                 |                    | NA                 |                    |
| 1,1,2,2-Tetrachloroethane          | UG/KG  | ND                 | 5                  | NA                 |                    | NA                 |                    | NA                 |                    |
| Toluene                            | UG/KG  | ND                 | 5                  | NA                 |                    | NA                 |                    | NA                 |                    |
| 1,1,1-Trichloroethane              | UG/KG  | ND                 | 5                  | NA                 |                    | NA                 |                    | NA                 |                    |
| 1,1,2-Trichloroethane              | UG/KG  | ND                 | 5                  | NA                 |                    | NA                 |                    | NA                 |                    |
| Trichloroethene                    | UG/KG  | ND                 | 5                  | NA                 |                    | NA                 |                    | NA                 |                    |
| Vinyl acetate                      | UG/KG  | ND                 | 10                 | NA                 |                    | NA                 |                    | NA                 |                    |
| Vinyl chloride                     | UG/KG  | ND                 | 10                 | NA                 |                    | NA                 |                    | NA                 |                    |
| Total Xylenes                      | UG/KG  | ND                 | 15                 | NA                 |                    | NA                 |                    | NA                 |                    |
| IS/SURRGATE(S)                     |        |                    |                    |                    |                    |                    |                    |                    |                    |
| Chlorobenzene-D5                   | %      | 94                 | 50-200             | NA                 |                    | NA                 |                    | NA                 |                    |
| 1,4-Difluorobenzene                | %      | 91                 | 50-200             | NA                 |                    | NA                 |                    | NA                 |                    |
| 1,4-Dichlorobenzene-D4             | %      | 75                 | 50-200             | NA                 |                    | NA                 |                    | NA                 |                    |
| Toluene-D8                         | %      | 98                 | 72-116             | NA                 |                    | NA                 |                    | NA                 |                    |
| p-Bromofluorobenzene               | %      | 89                 | 72-121             | NA                 |                    | NA                 |                    | NA                 |                    |
| 1,2-Dichloroethane-D4              | %      | 113                | 52-157             | NA                 |                    | NA                 |                    | NA                 |                    |

000012





Date: 07/31/1999  
Time: 10:41:25

Rept: AN0326

HYDRIX SOIL Samples  
METHOD 8270 - TCL SEMI-VOLATILE ORGANICS

| Client ID<br>Job No<br>Sample Date                 | Lab ID | Method Blank<br>A99-4287 |                    | A980610902      |                    | Reporting<br>Limit | Sample<br>Value | Reporting<br>Limit | Sample<br>Value |
|--|--------|--------------------------|--------------------|-----------------|--------------------|--------------------|-----------------|--------------------|-----------------|
|  |        | Sample<br>Value          | Reporting<br>Limit | Sample<br>Value | Reporting<br>Limit |                    |                 |                    |                 |
| Hexachloroethane                                   |        | ND                       | 330                | NA              | NA                 |                    | NA              |                    | NA              |
| Indeno(1,2,3-cd)pyrene                             |        | NC                       | 330                | NA              | NA                 |                    | NA              |                    | NA              |
| Isophorone   |        | ND                       | 330                | NA              | NA                 |                    | NA              |                    | NA              |
| 2-Methylnaphthalene                                |        | ND                       | 330                | NA              | NA                 |                    | NA              |                    | NA              |
| 2-Methylphenol                                     |        | ND                       | 330                | NA              | NA                 |                    | NA              |                    | NA              |
| 4-Methylphenol                                     |        | ND                       | 330                | NA              | NA                 |                    | NA              |                    | NA              |
| Naphthalene  |        | ND                       | 330                | NA              | NA                 |                    | NA              |                    | NA              |
| 2-Vinyltoluene                                     |        | ND                       | 1600               | NA              | NA                 |                    | NA              |                    | NA              |
| 3-Vinyltoluene                                     |        | ND                       | 1600               | NA              | NA                 |                    | NA              |                    | NA              |
| 1,2-Dibenzofuran                                   |        | ND                       | 330                | NA              | NA                 |                    | NA              |                    | NA              |
| Nitrobenzene                                       |        | ND                       | 330                | NA              | NA                 |                    | NA              |                    | NA              |
| 2-Nitrophenol                                      |        | ND                       | 330                | NA              | NA                 |                    | NA              |                    | NA              |
| 4-Nitrophenol                                      |        | ND                       | 1600               | NA              | NA                 |                    | NA              |                    | NA              |
| N-nitrosodiphenylamine                             |        | ND                       | 330                | NA              | NA                 |                    | NA              |                    | NA              |
| N-nitroso-Di-n-propylamine                         |        | ND                       | 330                | NA              | NA                 |                    | NA              |                    | NA              |
| Pentachlorophenol                                  |        | ND                       | 1600               | NA              | NA                 |                    | NA              |                    | NA              |
| Phenanthrene                                       |        | ND                       | 330                | NA              | NA                 |                    | NA              |                    | NA              |
| Phenol   |        | ND                       | 330                | NA              | NA                 |                    | NA              |                    | NA              |
| Pyrene   |        | ND                       | 330                | NA              | NA                 |                    | NA              |                    | NA              |
| 1,2,4-Trichlorobenzene                             |        | ND                       | 330                | NA              | NA                 |                    | NA              |                    | NA              |
| 2,4,5-Trichlorophenol                              |        | ND                       | 800                | NA              | NA                 |                    | NA              |                    | NA              |
| 2,4,6-Trichlorophenol                              |        | ND                       | 330                | NA              | NA                 |                    | NA              |                    | NA              |
| 1,2,3,4,6-Pentachlorocyclohexane (IS/SURROGATE(S)) |        |                          |                    |                 |                    |                    |                 |                    |                 |
| 1,4-Dichlorobenzene-D4                             |        | 115                      | 50-200             | NA              | NA                 |                    | NA              |                    | NA              |
| Naphthalene-D8                                     |        | 122                      | 50-200             | NA              | NA                 |                    | NA              |                    | NA              |
| Arenanthrene-D10                                   |        | 112                      | 50-200             | NA              | NA                 |                    | NA              |                    | NA              |
| Phenanthrene-D10                                   |        | 107                      | 50-200             | NA              | NA                 |                    | NA              |                    | NA              |
| Chrysene-D12                                       |        | 114                      | 50-200             | NA              | NA                 |                    | NA              |                    | NA              |
| Perylene-D12                                       |        | 92                       | 50-200             | NA              | NA                 |                    | NA              |                    | NA              |
| Nitrobenzene-D5                                    |        | 69                       | 20-114             | NA              | NA                 |                    | NA              |                    | NA              |
| Terphenyl-D14                                      |        | 78                       | 31-125             | NA              | NA                 |                    | NA              |                    | NA              |
| 2-Fluorobiphenyl                                   |        | 70                       | 38-116             | NA              | NA                 |                    | NA              |                    | NA              |
| 2-Fluorophenol                                     |        | 34                       | 23-105             | NA              | NA                 |                    | NA              |                    | NA              |
| Phenol-D5  |        | 26                       | 18-105             | NA              | NA                 |                    | NA              |                    | NA              |
| 2,4,6-Tribromophenol                               |        | 98                       | 25-122             | NA              | NA                 |                    | NA              |                    | NA              |

000014

Date: 07/31/1999  
 Time: 10:41:46

Rept: AN0326

HYDRIX SOIL Samples  
 METHOD 8082 - POLYCHLORINATED BIPHENYLS

| Client ID                     | Lab ID | Method Blank | Method Blank    | Reporting Limit | Sample Value    | Reporting Limit | Sample Value    | Reporting Limit | Sample Value    | Reporting Limit | Sample Value    |
|-------------------------------|--------|--------------|-----------------|-----------------|-----------------|-----------------|-----------------|-----------------|-----------------|-----------------|-----------------|
| Job No                        |        | A99-4287     | A980610302      |                 |                 |                 |                 |                 |                 |                 |                 |
| Sample Date                   |        |              |                 |                 |                 |                 |                 |                 |                 |                 |                 |
| Analyte                       | Units  | Sample Value | Reporting Limit | Sample Value    | Reporting Limit | Sample Value    | Reporting Limit | Sample Value    | Reporting Limit | Sample Value    | Reporting Limit |
| Aroclor 1216                  | UG/KG  | ND           | 40              | NA              |                 | NA              |                 | NA              |                 | NA              |                 |
| Aroclor 1221                  | UG/KG  | ND           | 80              | NA              |                 | NA              |                 | NA              |                 | NA              |                 |
| Aroclor 1232                  | UG/KG  | ND           | 40              | NA              |                 | NA              |                 | NA              |                 | NA              |                 |
| Aroclor 1242                  | UG/KG  | ND           | 40              | NA              |                 | NA              |                 | NA              |                 | NA              |                 |
| Aroclor 1248                  | UG/KG  | ND           | 40              | NA              |                 | NA              |                 | NA              |                 | NA              |                 |
| Aroclor 1254                  | UG/KG  | ND           | 40              | NA              |                 | NA              |                 | NA              |                 | NA              |                 |
| Aroclor 1260                  | UG/KG  | ND           | 40              | NA              |                 | NA              |                 | NA              |                 | NA              |                 |
| SURDATE(S)                    |        |              |                 |                 |                 |                 |                 |                 |                 |                 |                 |
| Tetrachloro- <i>n</i> -xylene | %      | 80           | 55-137          | NA              |                 | NA              |                 | NA              |                 | NA              |                 |
| Dibachlorodiphenyl            | %      | 76           | 50-150          | NA              |                 | NA              |                 | NA              |                 | NA              |                 |

000015

Date: 07/31/1999  
Time: 10:41:46

Rept: AN0326

HYDRIX SOIL Samples  
METHOD 8081 - TCL PESTICIDES

| Client ID<br>Job No<br>Sample Date | Lab ID | Method Blank<br>A99-4287 | A980626603         | Reporting<br>Limit | Sample<br>Value    | Reporting<br>Limit | Sample<br>Value    | Reporting<br>Limit | Sample<br>Value    |
|------------------------------------|--------|--------------------------|--------------------|--------------------|--------------------|--------------------|--------------------|--------------------|--------------------|
| Analyte                            | Units  | Sample<br>Value          | Reporting<br>Limit | Sample<br>Value    | Reporting<br>Limit | Sample<br>Value    | Reporting<br>Limit | Sample<br>Value    | Reporting<br>Limit |
| Aldrin                             | UG/KG  | ND                       | 8.0                | NA                 |                    | NA                 |                    | NA                 |                    |
| alpha-BHC                          | UG/KG  | ND                       | 8.0                | NA                 |                    | NA                 |                    | NA                 |                    |
| beta-BHC                           | UG/KG  | ND                       | 8.0                | NA                 |                    | NA                 |                    | NA                 |                    |
| gamma-BHC (Lindane)                | UG/KG  | ND                       | 8.0                | NA                 |                    | NA                 |                    | NA                 |                    |
| delta-BHC                          | UG/KG  | ND                       | 8.0                | NA                 |                    | NA                 |                    | NA                 |                    |
| Chlordane                          | UG/KG  | ND                       | 80                 | NA                 |                    | NA                 |                    | NA                 |                    |
| 4,4'-DDD                           | UG/KG  | ND                       | 16                 | NA                 |                    | NA                 |                    | NA                 |                    |
| 4,4'-DDE                           | UG/KG  | ND                       | 16                 | NA                 |                    | NA                 |                    | NA                 |                    |
| 4,4'-DDT                           | UG/KG  | ND                       | 16                 | NA                 |                    | NA                 |                    | NA                 |                    |
| Heptachlor                         | UG/KG  | ND                       | 16                 | NA                 |                    | NA                 |                    | NA                 |                    |
| Endosulfan I                       | UG/KG  | ND                       | 16                 | NA                 |                    | NA                 |                    | NA                 |                    |
| Endosulfan II                      | UG/KG  | ND                       | 16                 | NA                 |                    | NA                 |                    | NA                 |                    |
| Endosulfan Sulfate                 | UG/KG  | ND                       | 16                 | NA                 |                    | NA                 |                    | NA                 |                    |
| Endrin                             | UG/KG  | ND                       | 16                 | NA                 |                    | NA                 |                    | NA                 |                    |
| Endrin aldehyde                    | UG/KG  | ND                       | 32                 | NA                 |                    | NA                 |                    | NA                 |                    |
| Heptachlor                         | UG/KG  | ND                       | 8.0                | NA                 |                    | NA                 |                    | NA                 |                    |
| Heptachlor epoxide                 | UG/KG  | ND                       | 8.0                | NA                 |                    | NA                 |                    | NA                 |                    |
| Methoxychlor                       | UG/KG  | ND                       | 80                 | NA                 |                    | NA                 |                    | NA                 |                    |
| Toxaphene                          | UG/KG  | ND                       | 160                | NA                 |                    | NA                 |                    | NA                 |                    |
| SURrogate(S)                       |        |                          |                    |                    |                    |                    |                    |                    |                    |
| Tetrachloro-mylene                 | %      | 78                       | 30-130             | NA                 |                    | NA                 |                    | NA                 |                    |
| Decachlorobiphenyl                 | %      | 91                       | 21-169             | NA                 |                    | NA                 |                    | NA                 |                    |

000016

Date: 07/31/1999  
Time: 10:43:27

Rept: AN0326

HYDRIX SOIL Samples  
GZA-SW8463 - TAL METALS

| Client ID<br>Job No<br>Sample Date | Lab ID | Method Blank<br>A99-4287 | Reporting<br>Limit | Sample<br>Value | Method Blank<br>A99-4287<br>07/15/1999 | Reporting<br>Limit | Sample<br>Value | Reporting<br>Limit | Sample<br>Value |
|------------------------------------|--------|--------------------------|--------------------|-----------------|--|--------------------|-----------------|--------------------|-----------------|
| Aluminum - Total                   |        | NA                       |                    | NA              | 2.0                                    |                    | NA              |                    | NA              |
| Antimony - Total                   |        | NA                       |                    | ND              | 0.40                                   |                    | NA              |                    | NA              |
| Iron - Total                       |        | NA                       |                    | ND              | 2.0                                    |                    | NA              |                    | NA              |
| Magnesium - Total                  |        | NA                       |                    | ND              | 2.0                                    |                    | NA              |                    | NA              |
| Potassium - Total                  |        | NA                       |                    | ND              | 15.0                                   |                    | NA              |                    | NA              |
| Sodium - Total                     |        | NA                       |                    | ND              | 25.0                                   |                    | NA              |                    | NA              |
| Vanadium - Total                   |        | NA                       |                    | ND              | 0.12                                   |                    | NA              |                    | NA              |
| Thallium - Total                   |        | NA                       |                    | ND              | 0.53                                   |                    | NA              |                    | NA              |
| Lead - Total                       |        | NA                       |                    | 0.24            | 0.16                                   |                    | NA              |                    | NA              |
| Cadmium - Total                    |        | NA                       |                    | 0.5             | 2.0                                    |                    | NA              |                    | NA              |
| Barium - Total                     |        | NA                       |                    | ND              | 0.10                                   |                    | NA              |                    | NA              |
| Arsenic - Total                    |        | NA                       |                    | ND              | 0.25                                   |                    | NA              |                    | NA              |
| Beryllium - Total                  |        | NA                       |                    | ND              | 0.10                                   |                    | NA              |                    | NA              |
| Cobalt - Total                     |        | NA                       |                    | ND              | 0.10                                   |                    | NA              |                    | NA              |
| Manganese - Total                  |        | NA                       |                    | 5.4             | 0.10                                   |                    | NA              |                    | NA              |
| Mercury - Total                    |        | 0.033                    |                    | NA              |  |                    | NA              |                    | NA              |
| Nickel - Total                     |        | NA                       |                    | ND              | 0.13                                   |                    | NA              |                    | NA              |
| Selenium - Total                   |        | NA                       |                    | ND              | 0.40                                   |                    | NA              |                    | NA              |
| Silver - Total                     |        | NA                       |                    | ND              | 0.15                                   |                    | NA              |                    | NA              |
| Zinc - Total                       |        | NA                       |                    | 1.6             | 0.10                                   |                    | NA              |                    | NA              |

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Date: 07/31/1999  
Time: 10:43:27

HYDRIX SOIL Samples  
GZA-SW8463 - TAL METALS

Rept: AN0326

| Client ID<br>Job No<br>Sample Date | Lab ID | 07029-1A<br>A99-4287<br>07/02/1999 | A9428701MD         | 07029-1A<br>A99-4287<br>07/02/1999 | A9428701MS         | LCS<br>A99-4287 | A980622402         | LCS CLP Soils<br>A99-4287<br>07/15/1999 | A980631701         |
|------------------------------------|--------|------------------------------------|--------------------|------------------------------------|--------------------|-----------------|--------------------|---|--------------------|
| Analyte                            | Units  | Sample<br>Value                    | Reporting<br>Limit | Sample<br>Value                    | Reporting<br>Limit | Sample<br>Value | Reporting<br>Limit | Sample<br>Value                         | Reporting<br>Limit |
| Aluminum - Total                   | MG/KG  | 4970                               | 2.4                | 5200                               | 2.3                | NA              |                    | 3000                                    | 2.0                |
| Antimony - Total                   | MG/KG  | ND                                 | 0.49               | 21.5                               | 0.46               | NA              |                    | 21.4                                    | 0.40               |
| Arsenic - Total                    | MG/KG  | 3.3                                | 0.30               | 42.6                               | 0.29               | NA              |                    | 110                                     | 0.25               |
| Barium - Total                     | MG/KG  | 23.9                               | 0.12               | 69.1                               | 0.11               | NA              |                    | 136                                     | 0.099              |
| Beryllium - Total                  | MG/KG  | 0.44                               | 0.12               | 39.8                               | 0.11               | NA              |                    | 52.5                                    | 0.099              |
| Cadmium - Total                    | MG/KG  | 0.11                               | 0.061              | 40.5                               | 0.057              | NA              |                    | NA                                      | 2.0                |
| Calcium - Total                    | MG/KG  | 11400                              | 2.4                | 20700                              | 2.3                | NA              |                    | 768                                     |                    |
| Chromium - Total                   | MG/KG  | 7.4                                | 0.15               | 44.7                               | 0.14               | NA              |                    | NA                                      |                    |
| Cobalt - Total                     | MG/KG  | 5.3                                | 0.12               | 43.6                               | 0.11               | NA              |                    | 47.8                                    | 0.099              |
| Copper - Total                     | MG/KG  | 12.8                               | 0.16               | 53.6                               | 0.17               | NA              |                    | NA                                      |                    |
| Iron - Total                       | MG/KG  | 10600                              | 2.4                | 10300                              | 2.3                | NA              |                    | 4510                                    | 2.0                |
| Lead - Total                       | MG/KG  | 7.4                                | 0.19               | 45.9                               | 0.18               | NA              |                    | 45.4                                    | 0.16               |
| Magnesium - Total                  | MG/KG  | 6180                               | 2.4                | 13400                              | 2.3                | NA              |                    | 597                                     | 2.0                |
| Manganese - Total                  | MG/KG  | 136                                | 0.12               | 199                                | 0.11               | NA              |                    | 169                                     | 0.099              |
| Mercury - Total                    | MG/KG  | ND                                 | 0.040              | 0.77                               | 0.040              | 2.3             | 0.32               | NA                                      |                    |
| Nickel - Total                     | MG/KG  | 12.6                               | 0.16               | 49.6                               | 0.15               | NA              |                    | 42.9                                    | 0.13               |
| Potassium - Total                  | MG/KG  | 687                                | 18.2               | 3080                               | 17.2               | NA              |                    | 910                                     | 14.8               |
| Selenium - Total                   | MG/KG  | ND                                 | 0.49               | 35.7                               | 0.46               | NA              |                    | 77.2                                    | 0.40               |
| Silver - Total                     | MG/KG  | ND                                 | 0.18               | 42.0                               | 0.17               | NA              |                    | 36.8                                    | 0.15               |
| Sodium - Total                     | MG/KG  | 98.2                               | 30.4               | 2500                               | 28.7               | NA              |                    | 840                                     | 24.7               |
| Thallium - Total                   | MG/KG  | 1.6                                | 0.64               | 42.2                               | 0.61               | NA              |                    | 60.7                                    | 0.52               |
| Vanadium - Total                   | MG/KG  | 9.8                                | 0.15               | 49.4                               | 0.14               | NA              |                    | 61.0                                    | 0.12               |
| Zinc - Total                       | MG/KG  | 40.9                               | 0.12               | 73.9                               | 0.11               | NA              |                    | 125                                     | 0.15               |

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Date: 07/31/1999  
 Time: 10:43:37

HYDRIX SOIL Samples  
 WET CHEMISTRY ANALYSIS

Rept: AN0326

| Client ID<br>Job No<br>Sample Date                                 | Lab ID | Method Blank<br>A99-4287 |                 | Method Blank<br>A99-4287 |                 | Method Blank<br>A99-4287 |                 | Method Blank<br>A980652504 |                 |
|--|--------|--------------------------|-----------------|--------------------------|-----------------|--------------------------|-----------------|----------------------------|-----------------|
|  |        | Sample Value             | Reporting Limit | Sample Value             | Reporting Limit | Sample Value             | Reporting Limit | Sample Value               | Reporting Limit |
| Leachable Phosphorous<br>Nitrate<br>Leachable Total Organic Carbon | Units  |                          |                 |                          |                 |                          |                 |                            |                 |
|  | UG/G   | ND                       | 0.20            | NA                       | 0.050           | NA                       | NA              | NA                         | NA              |
|  | MG N/L | NA                       |                 | ND                       |                 | NA                       |                 | NA                         |                 |
|  |        | NA                       |                 | NA                       |                 | ND                       |                 | NA                         |                 |

00001

Date: 07/31/1999  
 Time: 10:43:37  
 Rept: AN0326  
 HYDRIX SOIL Samples  
 WET CHEMISTRY ANALYSIS

| Client ID<br>Job No<br>Sample Date | Lab ID   | LCS<br>A99-4287 | A980614501      | LCS<br>A99-4287 | A980624801      | LCS<br>A99-4287 | A980630501      | LCS<br>A99-4287 | A980635201      |
|------------------------------------|----------|-----------------|-----------------|-----------------|-----------------|-----------------|-----------------|-----------------|-----------------|
| Analyte                            | Units    | Sample Value    | Reporting Limit | Sample Value    | Reporting Limit | Sample Value    | Reporting Limit | Sample Value    | Reporting Limit |
| Leachable PH                       | S.U.     | 7.03            | 0.00000         | NA              | 0.20            | NA              |                 | NA              |                 |
| Leachable Phosphorous              | UG/G     | NA              |                 | 16.2            |                 | NA              |                 | NA              |                 |
| Nitrate                            | MG N/L   | NA              |                 | NA              |                 | 2.4             | 0.050           | NA              |                 |
| Redox Potential                    | M. VOLTS | NA              |                 | NA              |                 | NA              |                 | 228             | 0               |

| Client ID<br>Job No<br>Sample Date | Lab ID | LCS<br>A99-4287 | A980652501      | Matrix Spike B<br>A99-4287 | A980624802      | Matrix Spike B<br>A99-4287 | A980630502      | Matrix Spike B<br>A99-4287 | A980652502      |
|------------------------------------|--------|-----------------|-----------------|----------------------------|-----------------|----------------------------|-----------------|----------------------------|-----------------|
| Analyte                            | Units  | Sample Value    | Reporting Limit | Sample Value               | Reporting Limit | Sample Value               | Reporting Limit | Sample Value               | Reporting Limit |
| Leachable Total Organic Carbon     | UG/G   | 11400           | 200             | NA                         | 0.20            | NA                         |                 | 1940                       | 200             |
| Leachable Phosphorous              | UG/G   | NA              |                 | 9.6                        |                 | NA                         |                 | NA                         |                 |
| Nitrate                            | MG N/L | NA              |                 | NA                         |                 | 1.0                        | 0.050           | NA                         |                 |

| Client ID<br>Job No<br>Sample Date | Lab ID | Matrix Spike B<br>A99-4287 | A980624803      | Matrix Spike B<br>A99-4287 | A980630503      | Matrix Spike B<br>A99-4287 | A980652503      |
|------------------------------------|--------|----------------------------|-----------------|----------------------------|-----------------|----------------------------|-----------------|
| Analyte                            | Units  | Sample Value               | Reporting Limit | Sample Value               | Reporting Limit | Sample Value               | Reporting Limit |
| Leachable Phosphorous              | UG/G   | 10.6                       | 0.20            | NA                         | 0.050           | NA                         |                 |
| Nitrate                            | MG N/L | NA                         |                 | 1.0                        |                 | NA                         |                 |
| Leachable Total Organic Carbon     | UG/G   | NA                         |                 | NA                         |                 | 1920                       | 200             |



Date : 07/31/99 10:42  
 Job No: A99-4287

GZA GEOENVIRONMENTAL, INC. SOIL ANALYSES

Rept: AN0364

SDG: HYDRIX  
 Client Sample ID: VBLK19  
 Lab Sample ID: A9428703

MSB19  
 A9428704

| Analyte                             | Units of Measure | Concentration |              | % Recovery | QC LIMITS |
|-------------------------------------|------------------|---------------|--------------|------------|-----------|
|                                     |                  | Blank Spike   | Spike Amount |            |           |
| METHOD 8260 - TCL VOLATILE ORGANICS |                  |               |              |            |           |
| 1,1-Dichloroethene                  | UG/KG            | 60.9          | 50.0         | 122        | 64-147    |
| Trichloroethene                     | UG/KG            | 47.1          | 50.0         | 94         | 66-126    |
| Benzene                             | UG/KG            | 54.0          | 50.0         | 108        | 66-129    |
| Toluene                             | UG/KG            | 48.9          | 50.0         | 98         | 61-124    |
| Chlorobenzene                       | UG/KG            | 49.5          | 50.0         | 99         | 65-117    |

Date : 07/31/90 10:42  
 Job No: A99-428

GZA GEGENVIRONMENTAL, INC. SOIL ANALYSES

Rept: AN0364

SDG HYDRIX  
 Client Sample ID: Method Blank  
 Lab Sample ID: A980610902

Matrix Spike Blank  
 A980610901

| Analyte                                 | Units of Measure | Blank Spike | Concentration Spike Amount | % Recovery Blank Spike | QC LIMITS |
|---|------------------|-------------|----------------------------|------------------------|-----------|
| METHOD 8270 - TL SEMI-VOLATILE ORGANICS |                  |             |                            |                        |           |
| Phenol                                  | UG/KG            | 2031        | 3311                       | 61                     | 40-114    |
| 2-Chloropheno                           | UG/KG            | 2205        | 3311                       | 66                     | 40-115    |
| 1,4-Dichlorobenzene                     | UG/KG            | 1923        | 3311                       | 58                     | 28-104    |
| N-Nitroso-Di-n-propylamine              | UG/KG            | 2541        | 3311                       | 77                     | 47-104    |
| 1,2,4-Trichlorobenzene                  | UG/KG            | 2156        | 3311                       | 65                     | 49-107    |
| 4-Chloro-3-methylphenol                 | UG/KG            | 2417        | 3311                       | 73                     | 51-129    |
| Acenaphthene                            | UG/KG            | 2385        | 3311                       | 72                     | 52-104    |
| 4-Nitrophenol                           | UG/KG            | 2789        | 3311                       | 84                     | 47-114    |
| 2,4-Dinitrochlorobenzene                | UG/KG            | 2672        | 3311                       | 81                     | 60-114    |
| Pentachloropheno.                       | UG/KG            | 2872        | 3311                       | 87                     | 57-141    |
| Pyrene                                  | UG/KG            | 2767        | 3311                       | 84                     | 59-128    |

000022  
 STL Buffalo

\* Indicates Result is outside QC Limits

Date : 07/31/99 10:42  
Job No: A99-4287

GZA GEONVIRONMENTAL, INC. SOIL ANALYSES

Rept: AN0364

SDG: HYDRIX

Client Sample ID: Method Blank  
Lab Sample ID: A9B0610302

Matrix Spike Blank  
A9B0610301

| Analyte   | Units of Measure | Blank Spike | Concentration Spike Amount | % Recovery Blank Spike | QC LIMITS |
|---|------------------|-------------|----------------------------|------------------------|-----------|
| METHOD 8082 - POLYCHLORINATED BIPHENYLS<br>Aroclor 1254 | UG/KG            | 162         | 164                        | 99                     | 57-134    |

Rept: AN0364

GZA GEONVIRONMENTAL, INC. SOIL ANALYSES

Date : 07/31/99 10:42  
Job No: A99-4287

SDG: HYDRIX  
Client Sample ID: Method Blank Matrix Spike Blk Dup  
Lab Sample ID: A980626603 A980626601 A980626602

| Analyte                      | Units of Measure | Concentration |                 | Spike Amount |      | % Recovery |     | % RPD |     | QC LIMITS |        |
|------------------------------|------------------|---------------|-----------------|--------------|------|------------|-----|-------|-----|-----------|--------|
|                              |                  | Spike Blank   | Spike Blank Dup | SB           | SBD  | SB         | SBD | Avg   | RPD | RPD       | REC.   |
| METHOD 8081 - TCL PESTICIDES |                  |               |                 |              |      |            |     |       |     |           |        |
| Aldrin                       | UG/KG            | 32.9          | 26.3            | 32.8         | 32.5 | 100        | 81  | 91    | 21  | 30.0      | 48-128 |
| alpha-BHC                    | UG/KG            | 31.9          | 25.4            | 32.8         | 32.5 | 97         | 78  | 88    | 22  | 30.0      | 47-123 |
| beta-BHC                     | UG/KG            | 33.3          | 27.0            | 32.8         | 32.5 | 101        | 83  | 92    | 20  | 30.0      | 56-129 |
| delta-BHC                    | UG/KG            | 31.9          | 25.3            | 32.8         | 32.5 | 97         | 78  | 88    | 22  | 30.0      | 52-127 |
| gamma-BHC (Lindane)          | UG/KG            | 32.6          | 26.1            | 32.8         | 32.5 | 99         | 80  | 90    | 21  | 30.0      | 47-124 |
| 4,4'-DDD                     | UG/KG            | 34.8          | 27.7            | 32.8         | 32.5 | 106        | 85  | 96    | 22  | 30.0      | 52-133 |
| 4,4'-DDE                     | UG/KG            | 34.4          | 27.3            | 32.8         | 32.5 | 105        | 84  | 95    | 22  | 30.0      | 54-136 |
| 4,4'-DDT                     | UG/KG            | 34.6          | 27.4            | 32.8         | 32.5 | 105        | 84  | 95    | 22  | 30.0      | 59-148 |
| Dieldrin                     | UG/KG            | 33.7          | 28.4            | 32.8         | 32.5 | 109        | 88  | 99    | 21  | 30.0      | 51-132 |
| Endosulfan I                 | UG/KG            | 32.0          | 25.8            | 32.8         | 32.5 | 98         | 80  | 89    | 20  | 30.0      | 52-132 |
| Endosulfan II                | UG/KG            | 34.9          | 27.9            | 32.8         | 32.5 | 106        | 86  | 96    | 21  | 30.0      | 54-135 |
| Endosulfan Sulfate           | UG/KG            | 33.7          | 27.0            | 32.8         | 32.5 | 103        | 83  | 93    | 22  | 30.0      | 52-136 |
| Endrin aldehyde              | UG/KG            | 30.9          | 25.3            | 32.8         | 32.5 | 94         | 78  | 86    | 19  | 30.0      | 37-123 |
| Endrin                       | UG/KG            | 35.3          | 28.1            | 32.8         | 32.5 | 108        | 86  | 97    | 23  | 30.0      | 61-132 |
| Heptachlor                   | UG/KG            | 33.2          | 26.9            | 32.8         | 32.5 | 101        | 83  | 92    | 20  | 30.0      | 53-127 |
| Heptachlor epoxide           | UG/KG            | 34.4          | 27.8            | 32.8         | 32.5 | 105        | 86  | 96    | 20  | 30.0      | 55-128 |
| Methoxychlor                 | UG/KG            | 36.0          | 29.2            | 32.8         | 32.5 | 110        | 90  | 100   | 20  | 30.0      | 72-180 |

\* Indicates Result is outside QC Limits

Date : 07/31/99 10:43  
 Job No: A99-4287

GZA GEOENVIRONMENTAL, INC. SOIL ANALYSES  
 SAMPLE DATE 07/02/99

Rept: AN0364

SDC: HYDRIX  
 Client Sample ID: 07029-1A  
 Lab Sample ID: A9428701MS

| Analyte                | Units of Measure | Concentration |              | Spike Amount | % Recovery MS | QC LIMITS |
|------------------------|------------------|---------------|--------------|--------------|---------------|-----------|
|                        |                  | Sample        | Matrix Spike |              |               |           |
| GZA-S48463 - TA METALS |                  |               |              |              |               |           |
| TOTAL ANTIMONY         | MG/KG            | 0             | 21.47        | 57.5         | 37 *          | 80-120    |
| TOTAL ARSENIC          | MG/KG            | 3.08          | 42.62        | 57.5         | 68 *          | 80-120    |
| TOTAL BARIUM           | MG/KG            | 22.99         | 69.06        | 57.5         | 80            | 80-120    |
| TOTAL BERYLLIUM        | MG/KG            | 0.419         | 39.81        | 57.5         | 68 *          | 80-120    |
| TOTAL CADMIUM          | MG/KG            | 0.0932        | 40.54        | 57.5         | 70 *          | 80-120    |
| TOTAL CHROMIUM         | MG/KG            | 6.59          | 44.69        | 57.5         | 66 *          | 80-120    |
| TOTAL COBALT           | MG/KG            | 4.87          | 43.63        | 57.5         | 67 *          | 80-120    |
| TOTAL COPPER           | MG/KG            | 11.80         | 53.65        | 57.5         | 72 *          | 80-120    |
| TOTAL LEAD             | MG/KG            | 7.35          | 45.88        | 57.5         | 67 *          | 80-120    |
| TOTAL MANGANESE        | MG/KG            | 129.2         | 198.6        | 57.5         | 120 *         | 80-120    |
| TOTAL MERCURY          | MG/KG            | 0             | 0.773        | 0.80         | 96            | 80-120    |
| TOTAL NICKEL           | MG/KG            | 2.31          | 49.56        | 57.5         | 64 *          | 80-120    |
| TOTAL SELENIUM         | MG/KG            | 0             | 35.66        | 57.5         | 62 *          | 80-120    |
| TOTAL SILVER           | MG/KG            | 0.0349        | 41.97        | 57.5         | 72 *          | 80-120    |
| TOTAL THALLIUM         | MG/KG            | 1.44          | 42.19        | 57.5         | 70 *          | 80-120    |
| TOTAL VANADIUM         | MG/KG            | 9.01          | 49.40        | 57.5         | 70 *          | 80-120    |
| TOTAL ZINC             | MG/KG            | 41.56         | 73.94        | 57.5         | 56 *          | 80-120    |

\* Indicates value is outside QC Limits  
 NC = Not Calculated ND = Not Calculated

Date : 07/31/99 10:44  
 Job No: A99-4287

GZA GEOENVIRONMENTAL, INC. SOIL ANALYSES

Rept: AN0364

SDB: HYDRIX  
 Client Sample ID: Method Blank  
 Lab Sample ID: A980624804

Matrix Spike Blank  
 A980624802

Matrix Spike Blk Dup  
 A980624803

| Analyte  | Units of Measure | Concentration |                 | Spike Amount |     | % Recovery |     | QC LIMITS |      |        |
|--|------------------|---------------|-----------------|--------------|-----|------------|-----|-----------|------|--------|
|  |                  | Spike Blank   | Spike Blank Dup | SB           | SBD | SB         | SBD | RPD       | REC. |        |
| WET CHEMISTRY ANALYSIS<br>METHOD 365.2 - LEACHABLE PHOSPHOROUS | UG/G             | 9.60          | 10.60           | 10           | 10  | 96         | 106 | 101       | 20.0 | 89-124 |

Date : 07/31/99 10:44  
 Job No: A99-4287

GZA GEOENVIRONMENTAL, INC. SOIL ANALYSES

Rept: AN0364

Client Sample ID: Method Blank  
 Lab Sample ID: A980630504

SDI: HYDRIX  
 Matrix Spike Blank  
 A980630502

Matrix Spike Blk Dup  
 A980630503

| Analyte                                       | Units of Measure | Concentration |                 | % Recovery |     |     | QC LIMITS |        |
|---|------------------|---------------|-----------------|------------|-----|-----|-----------|--------|
|   |                  | Spike Blank   | Spike Blank Dup | SB         | SBD | Avg | RPD       | REC.   |
| NET CHEMISTRY ANALYSIS<br>METHOD 9200 NITRATE | MG N/L           | 1.00          | 1.00            | 100        | 100 | 100 | 20.0      | 74-122 |

\* Indicates result is outside QC Limits

000027

STL Buffalo

Date : 07/31/99 10:44  
 Job No: A99-428

GZA GEOENVIRONMENTAL, INC. SOIL ANALYSES

Rept: AN0364

SDI: HYDRIX  
 Client Sample ID: Method Blank Matrix Spike Blk Dup  
 Lab Sample ID: A980652504 A980652502 A980652503

| Analyte   | Units of Measure | Concentration |                 | % Recovery |      |      | QC LIMITS |      |    |   |      |        |
|---|------------------|---------------|-----------------|------------|------|------|-----------|------|----|---|------|--------|
|   |                  | Spike Blank   | Spike Blank Dup | SB         | SBD  | Avg  | RPD       | REC. |    |   |      |        |
| WET CHEMISTRY ANALYSIS<br>METHOD 9060 LEACHABLE TOTAL ORGANIC | UG/G             | 1940          | 1920            | 2000       | 2000 | 2000 | 97        | 96   | 97 | 1 | 20.0 | 81-117 |

\* Indicates Result is outside QC Limits



Date: 07/31/99  
Time: 10:42:41

GZA GEOENVIRONMENTAL  
SAMPLE CHRONOLOGY

Rept: AN0374  
Page: 1

METHOD 8260 - ICL VOLATILE ORGANICS

|  |                               |  |  |  |
|--|-------------------------------|--|--|--|
| Client Sample ID<br>Job No & Lab Sample ID | 07029-1A<br>A99-4287 A9428701 |  |  |  |
| Sample Date                                | 07/02/99 16:00                |  |  |  |
| Received Date                              | 07/02/99 16:40                |  |  |  |
| Extraction Date                            |                               |  |  |  |
| Analysis Date                              | 07/08/99 17:24                |  |  |  |
| Extraction HI Mt?                          |                               |  |  |  |
| Analytical HI Mt?                          | YES                           |  |  |  |
| Sample Matrix                              | SOIL                          |  |  |  |
| Dilution Factor                            | 1.0                           |  |  |  |
| Sample wt/vol                              | 5.15 GRAMS                    |  |  |  |
| % DRY                                      | 81.69                         |  |  |  |

000029

METHOD 8260 - TCL VOLATILE ORGANICS

| Sample Date | Client Sample ID | Job No & Lab Sample ID | Received Date | Extraction Date | Analysis Date  | Extraction HT Met? | Analytical HT Met? | Sample Matrix | Dilution Factor | Sample wt/vol | % DRY |
|-------------|------------------|------------------------|---------------|-----------------|----------------|--------------------|--------------------|---------------|-----------------|---------------|-------|
|             | VBLK19           | A99-4287 A9428703      |               |                 | 07/08/99 12:22 |                    |                    | SOIL          | LOW             |               |       |
|             |                  |                        |               |                 |                |                    |                    | 1.0           |                 |               |       |
|             |                  |                        |               |                 |                |                    |                    | 5.0           |                 | GRAMS         |       |
|             |                  |                        |               |                 |                |                    |                    | 100.00        |                 |               |       |

000030

Date: 07/31/99  
Time: 10:42:41

GZA GEOENVIRONMENTAL  
QC SAMPLE CHRONOLOGY

Rept: AN0374  
Page: 2

METHOD 8260 - TCL VOLATILE ORGANICS

|  |                            |       |  |  |
|--|----------------------------|-------|--|--|
| Client Sample ID<br>Job No & Lab Sample ID | MSB19<br>A99-4287 A9428704 |       |  |  |
| Sample Date                                |                            |       |  |  |
| Received Date                              |                            |       |  |  |
| Extraction Date                            | 07/08/99 11:15             |       |  |  |
| Analysis Date                              |                            |       |  |  |
| Extraction HF Mt?                          |                            |       |  |  |
| Analytical HF Mt?                          | SOIL                       | LOW   |  |  |
| Sample Matrix                              | 1.0                        |       |  |  |
| Dilution Factor                            | 5.0                        | GRAMS |  |  |
| Sample wt/vol                              | 100.00                     |       |  |  |
| % Dry                                      |                            |       |  |  |

METHOD 8270 - ICL SEMI-VOLATILE ORGANICS

| Client Sample ID<br>Job No & Lab Sample ID | 07029-1A<br>A99-4287 A9428701 |       |  |  |
|--|-------------------------------|-------|--|--|
| Sample Date                                | 07/02/99                      | 16:00 |  |  |
| Received Date                              | 07/02/99                      | 16:40 |  |  |
| Extraction Date                            | 07/12/99                      | 07:00 |  |  |
| Analysis Date                              | 07/21/99                      | 13:49 |  |  |
| Extraction #? Mt?                          | YES                           |       |  |  |
| Analytical HT Mt?                          | YES                           |       |  |  |
| Sample Matrix                              | SOIL                          | LOW   |  |  |
| Dilution Factor                            | 5.0                           |       |  |  |
| Sample wt/vol                              | 30.84                         | GRAMS |  |  |
| % Dry                                      | 83.15                         |       |  |  |

000032

METHOD 8270 - ICL SEMI-VOLATILE ORGANICS

| Client Sample ID<br>Job No & Lab Sample ID | Method Blank<br>A99-4287 A980610902 |       |  |  |
|--|-------------------------------------|-------|--|--|
| Sample Date                                |                                     |       |  |  |
| Received Date                              | 07/12/99 07:00                      |       |  |  |
| Extraction Date                            | 07/21/99 10:44                      |       |  |  |
| Analysis Date                              |                                     |       |  |  |
| Extraction HT Met?                         |                                     |       |  |  |
| Analytical HT Met?                         |                                     |       |  |  |
| Sample Matrix                              | SOIL                                | LOW   |  |  |
| Dilution Factor                            | 1.0                                 |       |  |  |
| Sample wt/vol                              | 30.39                               | GRAMS |  |  |
| % Dry                                      | 100.00                              |       |  |  |

000033

METHOD 8270 - TCL SEMI-VOLATILE ORGANICS

| Client: Sample ID<br>Job No & Lab Sample ID | Matrix Spike Blank<br>A99-4287 A980610901 |  |  |  |
|---|---|--|--|--|
| Sample Date                                 |   |  |  |  |
| Received Date                               | 07/12/99 07:00                            |  |  |  |
| Extraction Date                             | 07/21/99 14:56                            |  |  |  |
| Analysis Date                               |   |  |  |  |
| Extraction HI Met?                          |   |  |  |  |
| Analytical HI Met?                          |   |  |  |  |
| Sample Matrix                               | SOIL                                      |  |  |  |
| Dilution Factor                             | 1.0                                       |  |  |  |
| Sample wt/vol                               | 30.2                                      |  |  |  |
| % Dry                                       | 100.00                                    |  |  |  |

000034

METHOD 8081 - ICL PESTICIDES

| Job No & Lab Sample ID | Client Sample ID | 07029-1A<br>A99-4287 A9428701 |
|------------------------|------------------|-------------------------------|
| Sample Date            | 07/02/99         | 16:00                         |
| Received Date          | 07/02/99         | 16:40                         |
| Extraction Date        | 07/14/99         | 07:00                         |
| Analysis Date          | 07/23/99         | 06:43                         |
| Extraction HT Met?     | YES              |                               |
| Analytical HT Met?     | YES              |                               |
| Sample Matrix          | SOIL             | LOW                           |
| Dilution Factor        | 4.0              |                               |
| Sample Wt/vol          | 30.37            | GRAMS                         |
| % Dry                  | 83.15            |                               |

METHOD 8082 - POLYCHLORINATED BIPHENYLS

| Job No & Lab Sample ID | Client Sample ID | 07029-1A<br>A99-4287 A9428701 |
|------------------------|------------------|-------------------------------|
| Sample Date            | 07/02/99         | 16:00                         |
| Received Date          | 07/02/99         | 16:40                         |
| Extraction Date        | 07/09/99         | 07:00                         |
| Analysis Date          | 07/10/99         | 12:01                         |
| Extraction HT Met?     | YES              |                               |
| Analytical HT Met?     | YES              |                               |
| Sample Matrix          | SOIL             | LOW                           |
| Dilution Factor        | 1.0              |                               |
| Sample Wt/vol          | 30.81            | GRAMS                         |
| % Dry                  | 65.05            |                               |

Date: 07/31/99 10:44  
 Jobno: A99-4237

GZA GEONVIRONMENTAL  
 QC CHRONOLOGY

Rept: AN0369

| Lab ID     | Sample ID      | Units | Analyte           | Method | Dilution Factor | Sample Date    | Receive Date | TCLP Date | THI | Analysis Date | AHT | Matrix |
|------------|----------------|-------|-------------------|--------|-----------------|----------------|--------------|-----------|-----|---------------|-----|--------|
| A980631702 | Method Blank   | MG/KG | Antimony - Total  | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|            |                | MG/KG | Arsenic - Total   | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|            |                | MG/KG | Barium - Total    | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|            |                | MG/KG | Beryllium - Total | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|            |                | MG/KG | Calcium - Total   | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|            |                | MG/KG | Cobalt - Total    | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|            |                | MG/KG | Iron - Total      | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|            |                | MG/KG | Lead - Total      | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|            |                | MG/KG | Magnesium - Total | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|            |                | MG/KG | Manganese - Total | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|            |                | MG/KG | Nickel - Total    | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|            |                | MG/KG | Potassium - Total | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|            |                | MG/KG | Selenium - Total  | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|            |                | MG/KG | Silver - Total    | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|            |                | MG/KG | Sodium - Total    | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|            |                | MG/KG | Thallium - Total  | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|            |                | MG/KG | Vanadium - Total  | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|            |                | MG/KG | Zinc - Total      | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
| A980622402 | LCS            | MG/KG | Mercury - Total   | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
| A980631701 | LCS TCLP Soils | MG/KG | Aluminum - Total  | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|            |                | MG/KG | Antimony - Total  | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|            |                | MG/KG | Arsenic - Total   | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|            |                | MG/KG | Barium - Total    | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|            |                | MG/KG | Beryllium - Total | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|            |                | MG/KG | Calcium - Total   | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|            |                | MG/KG | Cobalt - Total    | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|            |                | MG/KG | Iron - Total      | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|            |                | MG/KG | Lead - Total      | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|            |                | MG/KG | Magnesium - Total | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|            |                | MG/KG | Manganese - Total | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|            |                | MG/KG | Nickel - Total    | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|            |                | MG/KG | Potassium - Total | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|            |                | MG/KG | Selenium - Total  | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|            |                | MG/KG | Silver - Total    | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|            |                | MG/KG | Sodium - Total    | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|            |                | MG/KG | Thallium - Total  | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|            |                | MG/KG | Vanadium - Total  | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|            |                | MG/KG | Zinc - Total      | 6010   | 1.00            | 07/15/99 08:00 | 07/15 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |

AHT = Analyte Holding Time Met  
 THI = TCLP Holding Time Met



Date: 07/31/99 10:44  
 Jobno: A99-4287

GZA GEONVIRONMENTAL  
 SAMPLE CHRONOLOGY

Rept: AN0369

| Lab ID   | Sample ID | Units | Analyte           | Method | Dilution Factor | Sample Date    | Receive Date | TCLP Date | THT | Analysis Date | AHT | Matrix |
|----------|-----------|-------|-------------------|--------|-----------------|----------------|--------------|-----------|-----|---------------|-----|--------|
| A9428701 | 07089-1A  | MG/KG | Aluminum - Total  | 6010   | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|          |           | MG/KG | Antimony - Total  | 6010   | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|          |           | MG/KG | Arsenic - Total   | 6010   | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|          |           | MG/KG | Barium - Total    | 6010   | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|          |           | MG/KG | Beryllium - Total | 6010   | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|          |           | MG/KG | Cadmium - Total   | 6010   | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|          |           | MG/KG | Calcium - Total   | 6010   | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|          |           | MG/KG | Chromium - Total  | 6010   | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|          |           | MG/KG | Cobalt - Total    | 6010   | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|          |           | MG/KG | Copper - Total    | 6010   | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|          |           | MG/KG | Iron - Total      | 6010   | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|          |           | MG/KG | Lead - Total      | 6010   | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|          |           | MG/KG | Magnesium - Total | 6010   | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|          |           | MG/KG | Manganese - Total | 6010   | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|          |           | MG/KG | Mercury - Total   | 7471   | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|          |           | MG/KG | Nickel - Total    | 6010   | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|          |           | MG/KG | Potassium - Total | 6010   | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|          |           | MG/KG | Selenium - Total  | 6010   | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|          |           | MG/KG | Silver - Total    | 6010   | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|          |           | MG/KG | Sodium - Total    | 6010   | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|          |           | MG/KG | Thallium - Total  | 6010   | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|          |           | MG/KG | Vanadium - Total  | 6010   | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |
|          |           | MG/KG | Zinc - Total      | 6010   | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA        | NA  | 07/16         | Yes | SOIL   |

000038

Date: 07/31/99  
Time: 10:43:06

GZA GEOENVIRONMENTAL  
QC SAMPLE CHRONOLOGY

Rept.: AN0374  
Page: 3

METHOD 8081 - TCL PESTICIDES

| Client Sample ID<br>Job No & Lab Sample ID | Method Blank<br>A99-4287 A980610302 | Method Blank<br>A99-4287 A980626603 |
|--|-------------------------------------|-------------------------------------|
| Sample Date                                |                                     |                                     |
| Received Date                              |                                     |                                     |
| Extraction Date                            |                                     | 07/14/99 07:00                      |
| Analysis Date                              |                                     | 07/23/99 04:38                      |
| Extraction HT Met?                         | NA                                  |                                     |
| Analytical HT Met?                         |                                     |                                     |
| Sample Matrix                              |                                     | SOIL LOW                            |
| Dilution Factor                            |                                     | 1.0                                 |
| Sample wt/vol                              |                                     | 30.3 GRAMS                          |
| % Dry                                      |                                     | 100.00                              |

METHOD 8082 - POLYCHLORINATED BIPHENYLS

| Client Sample ID<br>Job No & Lab Sample ID | Method Blank<br>A99-4287 A980610302 | Method Blank<br>A99-4287 A980626603 |
|--|-------------------------------------|-------------------------------------|
| Sample Date                                |                                     |                                     |
| Received Date                              |                                     |                                     |
| Extraction Date                            | 07/09/99 07:00                      |                                     |
| Analysis Date                              | 07/10/99 11:14                      |                                     |
| Extraction HT Met?                         |                                     | NA                                  |
| Analytical HT Met?                         |                                     |                                     |
| Sample Matrix                              | SOIL LOW                            |                                     |
| Dilution Factor                            | 1.0                                 |                                     |
| Sample wt/vol                              | 30.8 GRAMS                          |                                     |
| % Dry                                      | 100.00                              |                                     |

Date: 07/31/99  
Time: 10:43:06

GZA GEOENVIRONMENTAL  
QC SAMPLE CHRONOLOGY

Rept: AN0374  
Page: 2

METHOD 8081 - ICL PESTICIDES

| Client Sample ID<br>Job No & Lab Sample ID   | Matrix Spike Blank<br>A99-4287 A980626601                                    | Matrix Spike Blk Dup<br>A99-4287 A980626602                                  |
|--|--|--|
| Sample Date<br>Received Date<br>Extraction Date<br>Analysis Date<br>Extraction HI Met?<br>Analytical HI Met?<br>Sample Matrix<br>Dilution Factor<br>Sample Wt/vol<br>% Dry | 07/14/99 07:00<br>07/23/99 03:15<br>SOIL LOW<br>1.0 GRAMS<br>30.43<br>100.00 | 07/14/99 07:00<br>07/23/99 03:57<br>SOIL LOW<br>1.0 GRAMS<br>30.74<br>100.00 |

METHOD 8082 - POLYCHLORINATED BIPHENYLS

| Client Sample ID<br>Job No & Lab Sample ID   | Matrix Spike Blank<br>A99-4287 A980626601                                    | Matrix Spike Blk Dup<br>A99-4287 A980626602 |
|--|--|---|
| Sample Date<br>Received Date<br>Extraction Date<br>Analysis Date<br>Extraction HI Met?<br>Analytical HI Met?<br>Sample Matrix<br>Dilution Factor<br>Sample Wt/vol<br>% Dry | 07/09/99 07:00<br>07/10/99 11:38<br>SOIL LOW<br>1.0 GRAMS<br>30.41<br>100.00 | NA  |

000037

Date: 07/31/99 10:44  
 Jobno: A99-427

GZA GEOENVIRONMENTAL  
 QC CHRONOLOGY

Rept: AN0369

| Lab ID     | Sample ID         | Units      | Analyte           | Method         | Dilution Factor | Sample Date    | Receive Date | TCLP Date      | THT         | Analysis Date | AHT | Matrix |
|------------|-------------------|------------|-------------------|----------------|-----------------|----------------|--------------|----------------|-------------|---------------|-----|--------|
| A9428701MD | 07029-1A          | MG/KG      | Aluminum - Total  | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
|            |                   | MG/KG      | Antimony - Total  | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
|            |                   | MG/KG      | Arsenic - Total   | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
|            |                   | MG/KG      | Barium - Total    | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
|            |                   | MG/KG      | Beryllium - Total | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
|            |                   | MG/KG      | Cadmium - Total   | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
|            |                   | MG/KG      | Calcium - Total   | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
|            |                   | MG/KG      | Chromium - Total  | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
|            |                   | MG/KG      | Cobalt - Total    | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
|            |                   | MG/KG      | Copper - Total    | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
|            |                   | MG/KG      | Iron - Total      | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
|            |                   | MG/KG      | Lead - Total      | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
|            |                   | MG/KG      | Magnesium - Total | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
|            |                   | MG/KG      | Manganese - Total | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
|            |                   | MG/KG      | Mercury - Total   | 7471           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
|            |                   | MG/KG      | Nickel - Total    | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
|            |                   | MG/KG      | Potassium - Total | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
|            |                   | MG/KG      | Selenium - Total  | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
|            |                   | MG/KG      | Silver - Total    | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
|            |                   | A9428701MS | 07029-1A          | MG/KG          | Sodium - Total  | 6010           | 1.00         | 07/02/99 16:00 | 07/02 16:40 | NA            | NA  | 07/16  |
| MG/KG      | Thallium - Total  |            |                   | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
| MG/KG      | Vanadium - Total  |            |                   | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
| MG/KG      | Zinc - Total      |            |                   | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
| MG/KG      | Aluminum - Total  |            |                   | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
| MG/KG      | Antimony - Total  |            |                   | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
| MG/KG      | Arsenic - Total   |            |                   | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
| MG/KG      | Barium - Total    |            |                   | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
| MG/KG      | Beryllium - Total |            |                   | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
| MG/KG      | Cadmium - Total   |            |                   | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
| MG/KG      | Calcium - Total   |            |                   | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
| MG/KG      | Chromium - Total  |            |                   | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
| MG/KG      | Cobalt - Total    |            |                   | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
| MG/KG      | Copper - Total    |            |                   | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
| MG/KG      | Iron - Total      |            |                   | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
| MG/KG      | Lead - Total      |            |                   | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
| MG/KG      | Magnesium - Total |            |                   | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
| MG/KG      | Manganese - Total |            |                   | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
| MG/KG      | Mercury - Total   |            |                   | 7471           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
| MG/KG      | Nickel - Total    |            |                   | 6010           | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA             | NA          | 07/16         | Yes | SOIL   |
| MG/KG      | Potassium - Total | 6010       | 1.00              | 07/02/99 16:00 | 07/02 16:40     | NA             | NA           | 07/16          | Yes         | SOIL          |     |        |
| MG/KG      | Selenium - Total  | 6010       | 1.00              | 07/02/99 16:00 | 07/02 16:40     | NA             | NA           | 07/16          | Yes         | SOIL          |     |        |
| MG/KG      | Silver - Total    | 6010       | 1.00              | 07/02/99 16:00 | 07/02 16:40     | NA             | NA           | 07/16          | Yes         | SOIL          |     |        |
| MG/KG      | Sodium - Total    | 6010       | 1.00              | 07/02/99 16:00 | 07/02 16:40     | NA             | NA           | 07/16          | Yes         | SOIL          |     |        |
| MG/KG      | Thallium - Total  | 6010       | 1.00              | 07/02/99 16:00 | 07/02 16:40     | NA             | NA           | 07/16          | Yes         | SOIL          |     |        |
| MG/KG      | Vanadium - Total  | 6010       | 1.00              | 07/02/99 16:00 | 07/02 16:40     | NA             | NA           | 07/16          | Yes         | SOIL          |     |        |
| MG/KG      | Zinc - Total      | 6010       | 1.00              | 07/02/99 16:00 | 07/02 16:40     | NA             | NA           | 07/16          | Yes         | SOIL          |     |        |
| MG/KG      | Mercury - Total   | 7471       | 1.00              | 07/02/99 16:00 | 07/02 16:40     | NA             | NA           | 07/16          | Yes         | SOIL          |     |        |
| MG/KG      | Aluminum - Total  | 6010       | 1.00              | 07/15/99 08:00 | 07/15 16:40     | NA             | NA           | 07/16          | NA          | 07/16         | Yes | SOIL   |

AHT = Analysis Holding Time Met  
 THT = TCLP Holding Time Met  
 NA = Not Applicable

Date: 07/31/99 10:44  
 Jobno: A99-4287

GZA GEONVIRONMENTAL  
 SAMPLE CHRONOLOGY

Rept: AN0369

| Lab ID   | Sample ID    | Units             | Analyte                                    | Method            | Dilution Factor | Sample Date    | Receive Date | TCLP Date | THT | Analysis Date | AHT Matrix |
|----------|--------------|-------------------|--|-------------------|-----------------|----------------|--------------|-----------|-----|---------------|------------|
| A9428701 | 07029-1A     | UG/G<br>S.U.      | Leachable Phosphorous<br>Leachable pH      | 365.2<br>9045     | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA        | NA  | 07/13         | Yes SOIL   |
|          |              | UG/G              | Leachable Total Organic Carbon             | 9060              | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA        | NA  | 07/09         | Yes SOIL   |
|          |              | INVALID           | SHAKE EXTRACTION OF SOLID WASTE WITH WATER | D 3987-85         | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA        | NA  | 07/13         | Yes SOIL   |
| A9428702 | 07029-1A (L) | MG N/L<br>M.VOLTS | Nitrate<br>Redox Potential                 | 9200<br>D-1498-76 | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA        | NA  | 07/14         | No WATER   |
|          |              |                   |  |                   | 1.00            | 07/02/99 16:00 | 07/02 16:40  | NA        | NA  | 07/16         | Yes WATER  |

Date: 07/31/99 10:44  
 Jobno: A99-4287

GZA GEOENVIRONMENTAL  
 QC CHRONOLOGY

Rept: AN0369

| Lab ID     | Sample ID            | Units   | Analyte                        | Method    | Dilution Factor | Sample Date | Receive Date | TCLP Date | THT | Analysis Date | AHT Matrix |
|------------|----------------------|---------|--------------------------------|-----------|-----------------|-------------|--------------|-----------|-----|---------------|------------|
| A980624802 | Matrix Spike Blank   | UG/G    | Leachable Phosphorous          | 365.2     | 1.00            |             | - 16:40      | NA        | NA  | 07/13         | Yes SOIL   |
| A980630502 | Matrix Spike Blank   | MG N/L  | Nitrate                        | 9200      | 1.00            |             | - 16:40      | NA        | NA  | 07/14         | Yes WATER  |
| A980652502 | Matrix Spike Blank   | UG/G    | Leachable Total Organic Carbon | 9060      | 1.00            |             | - 16:40      | NA        | NA  | 07/20         | Yes SOIL   |
| A980624803 | Matrix Spike Blk Dup | UG/G    | Leachable Phosphorous          | 365.2     | 1.00            |             | - 16:40      | NA        | NA  | 07/13         | Yes SOIL   |
| A980630503 | Matrix Spike Blk Dup | MG N/L  | Nitrate                        | 9200      | 1.00            |             | - 16:40      | NA        | NA  | 07/14         | Yes WATER  |
| A980652503 | Matrix Spike Blk Dup | UG/G    | Leachable Total Organic Carbon | 9060      | 1.00            |             | - 16:40      | NA        | NA  | 07/20         | Yes SOIL   |
| A980630504 | Method Blank         | MG N/L  | Nitrate                        | 9200      | 1.00            |             | - 16:40      | NA        | NA  | 07/14         | Yes WATER  |
| A980652504 | Method Blank         | UG/G    | Leachable Total Organic Carbon | 9060      | 1.00            |             | - 16:40      | NA        | NA  | 07/20         | Yes SOIL   |
| A980614501 | LCS                  | S.U.    | Leachable pH                   | 9045      | 1.00            |             | - 16:40      | NA        | NA  | 07/09         | Yes SOIL   |
| A980624801 | LCS                  | UG/G    | Leachable Phosphorous          | 365.2     | 1.00            |             | - 16:40      | NA        | NA  | 07/13         | Yes SOIL   |
| A980630501 | LCS                  | MG N/L  | Nitrate                        | 9200      | 1.00            |             | - 16:40      | NA        | NA  | 07/14         | Yes WATER  |
| A980635201 | LCS                  | M.VOLTS | Redox Potential                | D-1498-76 | 1.00            |             | - 16:40      | NA        | NA  | 07/16         | Yes WATER  |
| A980652501 | LCS                  | UG/G    | Leachable Total Organic Carbon | 9060      | 1.00            |             | - 16:40      | NA        | NA  | 07/20         | Yes SOIL   |
| A980624804 | Method Blank         | UG/G    | Leachable Phosphorous          | 365.2     | 1.00            |             | - 16:40      | NA        | NA  | 07/13         | Yes SOIL   |

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AHT = Analysis Holding Time Met  
 THT = TCLP Holding Time Met  
 NA = Not Applicable

STL Buffalo

# Chain of Custody





DATE: / /

Upstate Laboratories, Inc.  
Analysis Results  
Report Number: 28199058  
Client I.D.: GZA

APPROVAL: \_\_\_\_\_  
QC: *JS* \_\_\_\_\_  
Lab I.D.: 10170  
Sampled by: Client

ID:28199058 Mat:Soil 55350/GRATWICK PARK S-1 1015H 10/07/99 G

| PARAMETERS        | RESULTS      | DATE ANAL. | KEY | FILE#  |
|-------------------|--------------|------------|-----|--------|
| Percent Solids    | 79%          | 10/08/99   |     | WC7698 |
| EPA Method 8150   |              |            |     |        |
| 2,4-D             | <420ug/kg dw | 10/21/99   |     | PA5287 |
| 2,4,5-T           | <42ug/kg dw  | 10/21/99   |     | PA5287 |
| 2,4,5-TP (Silvex) | <42ug/kg dw  | 10/21/99   |     | PA5287 |
| Dinoseb           | <42ug/kg dw  | 10/21/99   |     | PA5287 |

dw = Dry weight

REFERENCE NO. 10

21 November 2000

Mr. Dennis Munhall  
U.S. Environmental Protection Agency  
290 Broadway - 18th Floor  
New York, NY 10007-1866

**DOCUMENT CONTROL NO: SAT-1002.014**  
**SUBJECT: SAMPLING TRIP REPORT -**  
**LACKAWANNA FOUNDRY**

Dear Mr. Munhall:

Enclosed please find the Sampling Trip Report for samples collected during the Brownfields investigation conducted at the Lackawanna Foundry site on 7 November 2000. The samples were collected under Case No. 28706.

If you have any questions, do not hesitate to call me at (732) 417-5806.

Very truly yours,

ROY F. WESTON, INC.



Donna Janda  
Project Manager

enclosure

cc: J. Bulich, Lockheed ESAT  
W.S. Butterfield, SAT Program Manager, w/o enclosure  
SAT TDD file

## SAMPLING TRIP REPORT

**SITE NAME:** Lackawanna Foundry  
DCN #: SAT-1002.014  
CASE NO.: 28706

**EPA I.D. NO.:** NYSFN0204209

**SAMPLING DATE:** November 7, 2000

1. Site Location: Refer to Figure 1
2. Sample Locations: Refer to Figure 2
3. Sample Descriptions: Refer to Table 1
4. Laboratories Receiving Samples:

Sample Type

Name and Address of Laboratory

Target Compound List  
(TCL) Organics

American Analytical & Technical Services, Inc.  
11950 Industriplex Boulevard  
Baton Rouge, LA 70809

Target Analyte List (TAL)  
Inorganics

Chemtech Consulting Group  
Raritan Center  
205 Campus Plaza 1  
Edison, NJ 08837

5. Sample Dispatch Data:

Fourteen soil samples and one aqueous samples for TCL volatiles, BNA and Pesticides/PCB analysis, were shipped to American Analytical & Technical Services, Inc., on 11/07/00 at 1900 hours via Federal Express (Airbill Nos. 4955287293 and 4955287304).

Fourteen soil samples and one aqueous samples for TAL inorganics analysis were shipped to Chemtech Consulting Group on 11/07/00 at 1900 hours via Federal Express (Airbill Nos. 4955287234 and 4955287245).

6. On-Site Personnel:

| <u>Name</u>   | <u>Company</u> | <u>Duties on Site</u>                   |
|---------------|----------------|---|
| Donna Janda   | Region II SAT  | Project Manager, Sampler                |
| David Lewitt  | Region II SAT  | Site Health and Safety Officer, Sampler |
| Jess Anderson | Region II SAT  | Sample Management Officer, Sampler      |

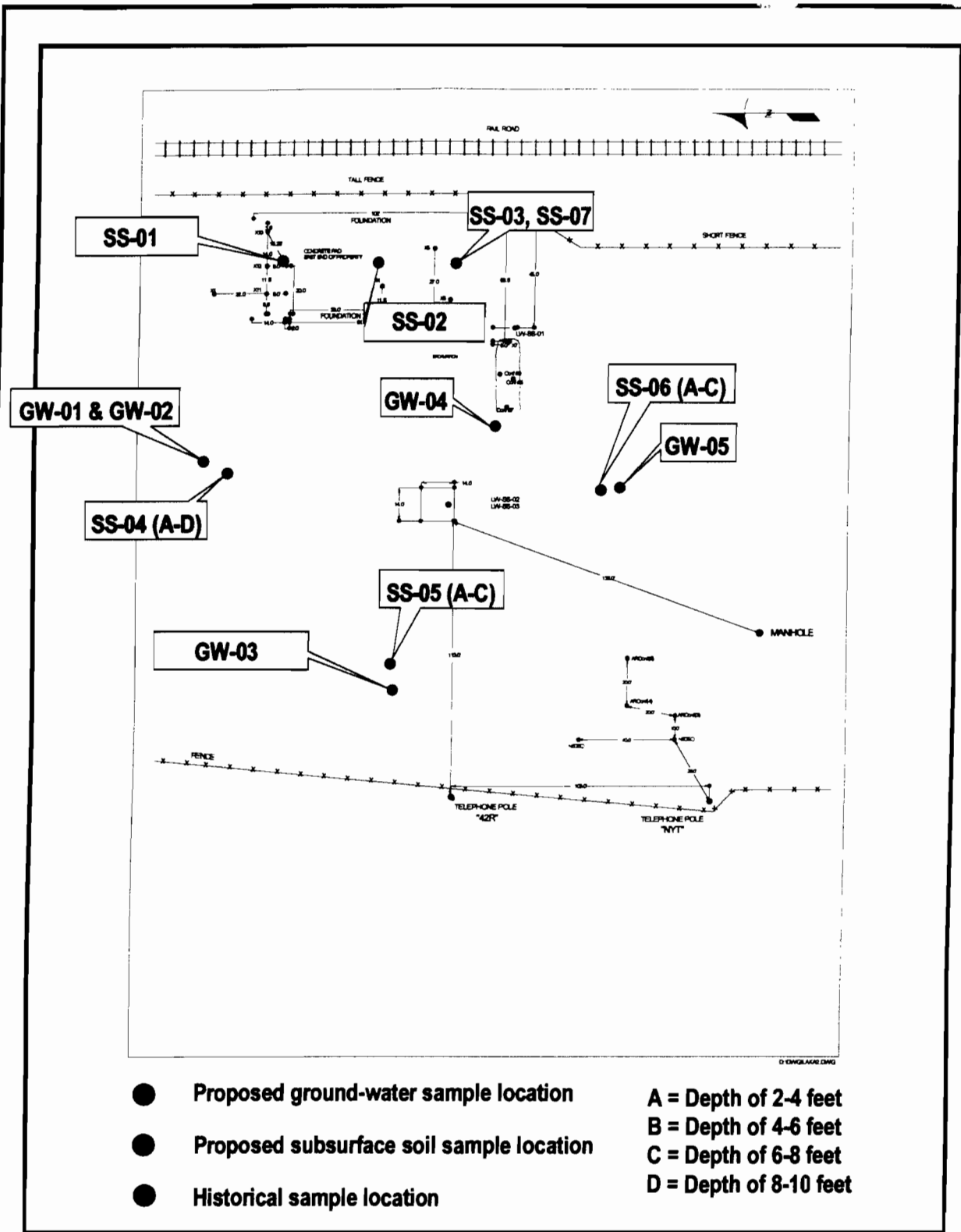
7. Additional Comments:

On November 7, 2000, the Region II Site Assessment Team (SAT) collected subsurface soil samples from the Lackawanna Foundry site. These samples were collected as part of a Brownfields investigation conducted on site. All samples collected by Region II SAT were analyzed for Target Compound List (TCL) and Target Analyte List (TAL), parameters through the U.S. EPA Contract Laboratory Program. Organic and Inorganic Traffic Reports are presented in Attachment 1.

8. Report Prepared by: Donna Janda Date: 11/21/00  
Donna Janda
9. Report Approved by: W. S. Butterfield Date: 11/21/00  
W. S. Butterfield, CHMM

**SITE LOCATION AND SAMPLE  
LOCATION MAPS**







**TABLE 1**  
**SAMPLE DESCRIPTIONS**  
**LACKAWANNA FOUNDRY**  
**LACKAWANNA, ERIE COUNTY, NEW YORK**

| <b>SAMPLE NUMBER<br/>1002-</b> | <b>ORGANIC<br/>CLP NO.</b> | <b>INORGANIC<br/>CLP NO.</b> | <b>DATE</b> | <b>TIME</b> | <b>COMMENTS</b>   |
|--------------------------------|----------------------------|------------------------------|-------------|-------------|---|
| SS-01                          | BNP05                      | MBNJ51                       | 11/07/00    | 1005        | Subsurface soil sample collected beneath concrete pad at 7-9 foot depth.        |
| SS-02                          | BNP06                      | MBNJ52                       | 11/07/00    | 1055        | Subsurface soil sample collected beneath concrete pad at 6-8 foot depth.        |
| SS-03                          | BNP07                      | MBNJ53                       | 11/07/00    | 1125        | Subsurface soil sample collected beneath concrete pad at 6-8 foot depth.        |
| SS-04A                         | BNP08                      | MBNJ54                       | 11/07/00    | 1440        | Subsurface soil sample collected from north section of site at 2-4 foot depth.  |
| SS-04B                         | BNP09                      | MBNJ55                       | 11/07/00    | 1450        | Subsurface soil sample collected from north section of site at 4-6 foot depth.  |
| SS-04C                         | BNP10                      | MBNJ56                       | 11/07/00    | 1455        | Subsurface soil sample collected from north section of site at 6-8 foot depth.  |
| SS-04D                         | BNP11                      | MBNJ57                       | 11/07/00    | 1515        | Subsurface soil sample collected from north section of site at 8-10 foot depth. |
| SS-05A                         | BNP12                      | MBNJ58                       | 11/07/00    | 1335        | Subsurface soil sample collected from west section of site at 2-4 foot depth.   |
| SS-05B                         | BNP13                      | MBNJ59                       | 11/07/00    | 1340        | Subsurface soil sample collected from west section of site at 4-6 foot depth.   |
| SS-05C                         | BNQ85                      | MBNJ60                       | 11/07/00    | 1345        | Subsurface soil sample collected from west section of site at 6-8 foot depth.   |
| SS-06A                         | BNQ87                      | MBNJ62                       | 11/07/00    | 1300        | Subsurface soil sample collected from south section of site at 2-4 foot depth.  |
| SS-06B                         | BNQ88                      | MBNJ63                       | 11/07/00    | 1305        | Subsurface soil sample collected from south section of site at 4-6 foot depth.  |
| SS-06C                         | BNQ89                      | MBNJ64                       | 11/07/00    | 1310        | Subsurface soil sample collected from south section of site at 6-8 foot depth.  |
| SS-07                          | BNQ91                      | MBNJ66                       | 11/07/00    | 1135        | Duplicate sample of 1002-SS-03.   |
| RB-01                          | BNQ92                      | MBNJ67                       | 11/07/00    | 1225        | Rinsate blank collected from split spoon and stainless steel bowl and trowel.   |

**ATTACHMENT 1**  
**TRAFFIC REPORTS/CHAIN OF CUSTODY RECORDS**



United States Environmental Protection Agency  
Contract Laboratory Program

**Organic Traffic Report  
& Chain of Custody Record**  
(For Organic CLP Analysis)

Case No. **28706**

|   |  |   |  |   |  |   |  |  |  |  |  |   |  |
|---|--|---|--|---|--|---|--|--|--|--|--|---|--|
| 1. Project Code<br><b>1003</b>                            |  | 2. Region No.<br><b>2</b>               |  | 3. Sampling Co.<br><b>Key-E-Lab</b>   |  | 4. Date Shipped<br><b>11/7/00</b>   |  | 5. Carrier<br><b>FedEx</b>   |  | 6. Matrix<br>(Enter in Column A)<br>1. Surface Water<br>2. Ground Water<br>3. Leachate<br>4. Field QC<br>5. Soil/Sediment<br>6. PE-water<br>7. PE-soil<br>8. Other (specify in Column A) |  | 7. Preservative<br>(Enter in Column D)<br>1. HCl<br>2. HNO3<br>3. NaHSO4<br>4. H2SO4<br>5. Ice only<br>6. CH3OH<br>7. Other (specify in Column D)<br>N. Not Preserved |  |
| Account Code  |  | Sampler (Name)<br><b>DM McJannet</b>    |  | Airbill Number<br><b>495 5287304</b>  |  | 5. Ship To:<br><b>American Analytical + Technical Serv. Inc.<br/>11950 Industriplex Blvd.<br/>Baton Rouge, LA 70809</b>   |  |  |  |  |  |   |  |
| Site Name<br><b>Lackawanna Foundry<br/>Lackawanna, NY</b> |  | Sampler Signature<br><i>[Signature]</i> |  | Long-Term Action<br>IA <input type="checkbox"/> PA <input type="checkbox"/> REM <input type="checkbox"/> RFS <input type="checkbox"/><br>RO <input type="checkbox"/> RI <input type="checkbox"/> RA <input type="checkbox"/> O&M <input type="checkbox"/> |  | 3. Purpose**<br>SF <input type="checkbox"/> PRP <input type="checkbox"/> ST <input type="checkbox"/> FED <input type="checkbox"/> BZ <input checked="" type="checkbox"/>  |  | Regional Specific Tracking Number or Tag Numbers<br><b>Tag Nos. 87-77-79</b> |  | H Mo/Day/Year/Time Sample Collection<br><b>11/7/00 1005</b>  |  | J Sampler Initials<br><b>MSJ</b>  |  |
| City, State<br><b>Lackawanna, NY</b>                      |  | Site Spill ID<br><b>400</b>             |  | Op Unit<br><b>400</b>   |  | E RAS Analysis<br>TA (circle one) PR* 7, 14, 21, 28, 35, 42, 49, 56, 63, 70, 77, 84, 91, 98, 105, 112, 119, 126, 133, 140, 147, 154, 161, 168, 175, 182, 189, 196, 203, 210, 217, 224, 231, 238, 245, 252, 259, 266, 273, 280, 287, 294, 301, 308, 315, 322, 329, 336, 343, 350, 357, 364, 371, 378, 385, 392, 399, 406, 413, 420, 427, 434, 441, 448, 455, 462, 469, 476, 483, 490, 497, 504, 511, 518, 525, 532, 539, 546, 553, 560, 567, 574, 581, 588, 595, 602, 609, 616, 623, 630, 637, 644, 651, 658, 665, 672, 679, 686, 693, 700, 707, 714, 721, 728, 735, 742, 749, 756, 763, 770, 777, 784, 791, 798, 805, 812, 819, 826, 833, 840, 847, 854, 861, 868, 875, 882, 889, 896, 903, 910, 917, 924, 931, 938, 945, 952, 959, 966, 973, 980, 987, 994, 1001, 1008, 1015, 1022, 1029, 1036, 1043, 1050, 1057, 1064, 1071, 1078, 1085, 1092, 1099, 1106, 1113, 1120, 1127, 1134, 1141, 1148, 1155, 1162, 1169, 1176, 1183, 1190, 1197, 1204, 1211, 1218, 1225, 1232, 1239, 1246, 1253, 1260, 1267, 1274, 1281, 1288, 1295, 1302, 1309, 1316, 1323, 1330, 1337, 1344, 1351, 1358, 1365, 1372, 1379, 1386, 1393, 1400, 1407, 1414, 1421, 1428, 1435, 1442, 1449, 1456, 1463, 1470, 1477, 1484, 1491, 1498, 1505, 1512, 1519, 1526, 1533, 1540, 1547, 1554, 1561, 1568, 1575, 1582, 1589, 1596, 1603, 1610, 1617, 1624, 1631, 1638, 1645, 1652, 1659, 1666, 1673, 1680, 1687, 1694, 1701, 1708, 1715, 1722, 1729, 1736, 1743, 1750, 1757, 1764, 1771, 1778, 1785, 1792, 1799, 1806, 1813, 1820, 1827, 1834, 1841, 1848, 1855, 1862, 1869, 1876, 1883, 1890, 1897, 1904, 1911, 1918, 1925, 1932, 1939, 1946, 1953, 1960, 1967, 1974, 1981, 1988, 1995, 2002, 2009, 2016, 2023, 2030, 2037, 2044, 2051, 2058, 2065, 2072, 2079, 2086, 2093, 2100, 2107, 2114, 2121, 2128, 2135, 2142, 2149, 2156, 2163, 2170, 2177, 2184, 2191, 2198, 2205, 2212, 2219, 2226, 2233, 2240, 2247, 2254, 2261, 2268, 2275, 2282, 2289, 2296, 2303, 2310, 2317, 2324, 2331, 2338, 2345, 2352, 2359, 2366, 2373, 2380, 2387, 2394, 2401, 2408, 2415, 2422, 2429, 2436, 2443, 2450, 2457, 2464, 2471, 2478, 2485, 2492, 2499, 2506, 2513, 2520, 2527, 2534, 2541, 2548, 2555, 2562, 2569, 2576, 2583, 2590, 2597, 2604, 2611, 2618, 2625, 2632, 2639, 2646, 2653, 2660, 2667, 2674, 2681, 2688, 2695, 2702, 2709, 2716, 2723, 2730, 2737, 2744, 2751, 2758, 2765, 2772, 2779, 2786, 2793, 2800, 2807, 2814, 2821, 2828, 2835, 2842, 2849, 2856, 2863, 2870, 2877, 2884, 2891, 2898, 2905, 2912, 2919, 2926, 2933, 2940, 2947, 2954, 2961, 2968, 2975, 2982, 2989, 2996, 3003, 3010, 3017, 3024, 3031, 3038, 3045, 3052, 3059, 3066, 3073, 3080, 3087, 3094, 3101, 3108, 3115, 3122, 3129, 3136, 3143, 3150, 3157, 3164, 3171, 3178, 3185, 3192, 3199, 3206, 3213, 3220, 3227, 3234, 3241, 3248, 3255, 3262, 3269, 3276, 3283, 3290, 3297, 3304, 3311, 3318, 3325, 3332, 3339, 3346, 3353, 3360, 3367, 3374, 3381, 3388, 3395, 3402, 3409, 3416, 3423, 3430, 3437, 3444, 3451, 3458, 3465, 3472, 3479, 3486, 3493, 3500, 3507, 3514, 3521, 3528, 3535, 3542, 3549, 3556, 3563, 3570, 3577, 3584, 3591, 3598, 3605, 3612, 3619, 3626, 3633, 3640, 3647, 3654, 3661, 3668, 3675, 3682, 3689, 3696, 3703, 3710, 3717, 3724, 3731, 3738, 3745, 3752, 3759, 3766, 3773, 3780, 3787, 3794, 3801, 3808, 3815, 3822, 3829, 3836, 3843, 3850, 3857, 3864, 3871, 3878, 3885, 3892, 3899, 3906, 3913, 3920, 3927, 3934, 3941, 3948, 3955, 3962, 3969, 3976, 3983, 3990, 3997, 4004, 4011, 4018, 4025, 4032, 4039, 4046, 4053, 4060, 4067, 4074, 4081, 4088, 4095, 4102, 4109, 4116, 4123, 4130, 4137, 4144, 4151, 4158, 4165, 4172, 4179, 4186, 4193, 4200, 4207, 4214, 4221, 4228, 4235, 4242, 4249, 4256, 4263, 4270, 4277, 4284, 4291, 4298, 4305, 4312, 4319, 4326, 4333, 4340, 4347, 4354, 4361, 4368, 4375, 4382, 4389, 4396, 4403, 4410, 4417, 4424, 4431, 4438, 4445, 4452, 4459, 4466, 4473, 4480, 4487, 4494, 4501, 4508, 4515, 4522, 4529, 4536, 4543, 4550, 4557, 4564, 4571, 4578, 4585, 4592, 4599, 4606, 4613, 4620, 4627, 4634, 4641, 4648, 4655, 4662, 4669, 4676, 4683, 4690, 4697, 4704, 4711, 4718, 4725, 4732, 4739, 4746, 4753, 4760, 4767, 4774, 4781, 4788, 4795, 4802, 4809, 4816, 4823, 4830, 4837, 4844, 4851, 4858, 4865, 4872, 4879, 4886, 4893, 4900, 4907, 4914, 4921, 4928, 4935, 4942, 4949, 4956, 4963, 4970, 4977, 4984, 4991, 4998, 5005, 5012, 5019, 5026, 5033, 5040, 5047, 5054, 5061, 5068, 5075, 5082, 5089, 5096, 5103, 5110, 5117, 5124, 5131, 5138, 5145, 5152, 5159, 5166, 5173, 5180, 5187, 5194, 5201, 5208, 5215, 5222, 5229, 5236, 5243, 5250, 5257, 5264, 5271, 5278, 5285, 5292, 5299, 5306, 5313, 5320, 5327, 5334, 5341, 5348, 5355, 5362, 5369, 5376, 5383, 5390, 5397, 5404, 5411, 5418, 5425, 5432, 5439, 5446, 5453, 5460, 5467, 5474, 5481, 5488, 5495, 5502, 5509, 5516, 5523, 5530, 5537, 5544, 5551, 5558, 5565, 5572, 5579, 5586, 5593, 5600, 5607, 5614, 5621, 5628, 5635, 5642, 5649, 5656, 5663, 5670, 5677, 5684, 5691, 5698, 5705, 5712, 5719, 5726, 5733, 5740, 5747, 5754, 5761, 5768, 5775, 5782, 5789, 5796, 5803, 5810, 5817, 5824, 5831, 5838, 5845, 5852, 5859, 5866, 5873, 5880, 5887, 5894, 5901, 5908, 5915, 5922, 5929, 5936, 5943, 5950, 5957, 5964, 5971, 5978, 5985, 5992, 5999, 6006, 6013, 6020, 6027, 6034, 6041, 6048, 6055, 6062, 6069, 6076, 6083, 6090, 6097, 6104, 6111, 6118, 6125, 6132, 6139, 6146, 6153, 6160, 6167, 6174, 6181, 6188, 6195, 6202, 6209, 6216, 6223, 6230, 6237, 6244, 6251, 6258, 6265, 6272, 6279, 6286, 6293, 6300, 6307, 6314, 6321, 6328, 6335, 6342, 6349, 6356, 6363, 6370, 6377, 6384, 6391, 6398, 6405, 6412, 6419, 6426, 6433, 6440, 6447, 6454, 6461, 6468, 6475, 6482, 6489, 6496, 6503, 6510, 6517, 6524, 6531, 6538, 6545, 6552, 6559, 6566, 6573, 6580, 6587, 6594, 6601, 6608, 6615, 6622, 6629, 6636, 6643, 6650, 6657, 6664, 6671, 6678, 6685, 6692, 6699, 6706, 6713, 6720, 6727, 6734, 6741, 6748, 6755, 6762, 6769, 6776, 6783, 6790, 6797, 6804, 6811, 6818, 6825, 6832, 6839, 6846, 6853, 6860, 6867, 6874, 6881, 6888, 6895, 6902, 6909, 6916, 6923, 6930, 6937, 6944, 6951, 6958, 6965, 6972, 6979, 6986, 6993, 7000, 7007, 7014, 7021, 7028, 7035, 7042, 7049, 7056, 7063, 7070, 7077, 7084, 7091, 7098, 7105, 7112, 7119, 7126, 7133, 7140, 7147, 7154, 7161, 7168, 7175, 7182, 7189, 7196, 7203, 7210, 7217, 7224, 7231, 7238, 7245, 7252, 7259, 7266, 7273, 7280, 7287, 7294, 7301, 7308, 7315, 7322, 7329, 7336, 7343, 7350, 7357, 7364, 7371, 7378, 7385, 7392, 7399, 7406, 7413, 7420, 7427, 7434, 7441, 7448, 7455, 7462, 7469, 7476, 7483, 7490, 7497, 7504, 7511, 7518, 7525, 7532, 7539, 7546, 7553, 7560, 7567, 7574, 7581, 7588, 7595, 7602, 7609, 7616, 7623, 7630, 7637, 7644, 7651, 7658, 7665, 7672, 7679, 7686, 7693, 7700, 7707, 7714, 7721, 7728, 7735, 7742, 7749, 7756, 7763, 7770, 7777, 7784, 7791, 7798, 7805, 7812, 7819, 7826, 7833, 7840, 7847, 7854, 7861, 7868, 7875, 7882, 7889, 7896, 7903, 7910, 7917, 7924, 7931, 7938, 7945, 7952, 7959, 7966, 7973, 7980, 7987, 7994, 8001, 8008, 8015, 8022, 8029, 8036, 8043, 8050, 8057, 8064, 8071, 8078, 8085, 8092, 8099, 8106, 8113, 8120, 8127, 8134, 8141, 8148, 8155, 8162, 8169, 8176, 8183, 8190, 8197, 8204, 8211, 8218, 8225, 8232, 8239, 8246, 8253, 8260, 8267, 8274, 8281, 8288, 8295, 8302, 8309, 8316, 8323, 8330, 8337, 8344, 8351, 8358, 8365, 8372, 8379, 8386, 8393, 8400, 8407, 8414, 8421, 8428, 8435, 8442, 8449, 8456, 8463, 8470, 8477, 8484, 8491, 8498, 8505, 8512, 8519, 8526, 8533, 8540, 8547, 8554, 8561, 8568, 8575, 8582, 8589, 8596, 8603, 8610, 8617, 8624, 8631, 8638, 8645, 8652, 8659, 8666, 8673, 8680, 8687, 8694, 8701, 8708, 8715, 8722, 8729, 8736, 8743, 8750, 8757, 8764, 8771, 8778, 8785, 8792, 8799, 8806, 8813, 8820, 8827, 8834, 8841, 8848, 8855, 8862, 8869, 8876, 8883, 8890, 8897, 8904, 8911, 8918, 8925, 8932, 8939, 8946, 8953, 8960, 8967, 8974, 8981, 8988, 8995, 9002, 9009, 9016, 9023, 9030, 9037, 9044, 9051, 9058, 9065, 9072, 9079, 9086, 9093, 9100, 9107, 9114, 9121, 9128, 9135, 9142, 9149, 9156, 9163, 9170, 9177, 9184, 9191, 9198, 9205, 9212, 9219, 9226, 9233, 9240, 9247, 9254, 9261, 9268, 9275, 9282, 9289, 9296, 9303, 9310, 9317, 9324, 9331, 9338, 9345, 9352, 9359, 9366, 9373, 9380, 9387, 9394, 9401, 9408, 9415, 9422, 9429, 9436, 9443, 9450, 9457, 9464, 9471, 9478, 9485, 9492, 9499, 9506, 9513, 9520, 9527, 9534, 9541, 9548, 9555, 9562, 9569, 9576, 9583, 9590, 9597, 9604, 9611, 9618, 9625, 9632, 9639, 9646, 9653, 9660, 9667, 9674, 9681, 9688, 9695, 9702, 9709, 9716, 9723, 9730, 9737, 9744, 9751, 9758, 9765, 9772, 9779, 9786, 9793, 9800, 9807, 9814, 9821, 9828, 9835, 9842, 9849, 9856, 9863, 9870, 9877, 9884, 9891, 9898, 9905, 9912, 9919, 9926, 9933, 9940, 9947, 9954, 9961, 9968, 9975, 9982, 9989, 9996, 10003, 10010, 10017, 10024, 10031, 10038, 10045, 10052, 10059, 10066, 10073, 10080, 10087, 10094, 10101, 10108, 10115, 10122, 10129, 10136, 10143, 10150, 10157, 10164, 10171, 10178, 10185, 10192, 10199, 10206, 10213, 10220, 10227, 10234, 10241, 10248, 10255, 10262, 10269, 10276, 10283, 10290, 10297, 10304, 10311, 10318, 10325, 10332, 10339, 10346, 10353, 10360, 10367, 10374, 10381, 10388, 10395, 10402, 10409, 10416, 10423, 10430, 10437, 10444, 10451, 10458, 10465, 10472, 10479, 10486, 10493, 10500, 10507, 10514, 10521, 10528, 10535, 10542, 10549, 10556, 10563, 10570, 10577, 10584, 10591, 10598, 10605, 10612, 10619, 10626, 10633, 10640, 10647, 10654, 10661, 10668, 10675, 10682, 10689, 10696, 10703, 10710, 10717, 10724, 10731, 10738, 10745, 10752, 10759, 10766, 10773, 10780, 10787, 10794, 10801, 10808, 10815, 10822, 10829, 10836, 10843, 10850, 10857, 10864, 10871, 10878, 10885, 10892, 10899, 10906, 10913, 10920, 10927, 10934, 10941, 10948, 10955, 10962, 10969, 10976, 10983, 10990, 10997, 11004, 11011, 11018, 11025, 11032, 11039, 11046, 11053, 11060, 11067, 11074, 11081, 11088, 11095, 11102, 11109, 11116, 11123, 11130, 11137, 11144, 11151, 11158, 11165, 11172, 11179, 11186, 11193, 11200, 11207, 11214, 11221, 11228, 11235, 11242, 11249, 11256, 11263, 11270, 11277, 11284, 11291, 11298, 11305, 11312, 11319, 11326, 11333, 11340, 11347, 11354, 11361, 11368, 11375, 11382, 11389, 11396, 11403, 11410, 11417, 11424, 11431, 11438, 11445, 11452, 11459, 11466, 11473, 11480, 11487, 11494, 11501, 11508, 11515, 11522, 11529, 11536, 11543, 11550, 11557, 11564, 11571, 11578, 11585, 11592, 11599, 11606, 11613, 11620, 11627, 11634, 11641, 11648, 11655, 11662, 11669, 11676, 11683, 11690, 11697, 11704, 11711, 11718, 11725, 11732, 11739, 11746, 11753, 11760, 11767, 11774, 11781, 11788, 11795, 11802, 11809, 11816, 11823, 11830, 11837, 11844, 11851, 11858, 11865, 11872, 11879, 11886, 11893, 11900, 11907, 11914, 11921, 11928, 11935, 11942, 11949, 11956, 11963, 11970, 11977, 11984, 11991, 11998, 12005, 12012, 12019, 12026, 12033, 12040, 12047, 12054, 12061, 12068, 12075, 12082, 12089, 12096, 12103, 12110, 12117, 12124, 12131, 12138, 12145, 12152, 12159, 12166, 12173, 12180, 12187, 12194, 12201, 12208, 12215, 12222, 12229, 12236, 12243, 12250, 12257, 12264, 12271, 12278, 12285, 12292, 12299, 12306, 12313, 12320, 12327, 12334, 12341, 12348, 12355, 12362, 12369, 12376, 12383, 12390, 12397, 12404, 12411, 12418, 12425, 12432, 12439, 12446, 12453, 12460, 12467, 12474, 12481, 12488, 12495, 12502, 12509, 12516, 12523, 12530, 12537, 12544, 12551, 12558, 12565, 12572, 12579, 12586, 12593, 12600, 12607, 12614, 12621, 12628, 12635, 12642, 12649, 12656, 12663, 12670, 12677, 12684, 12691, 12698, 12705, 12712, 12719, 12726, 12733, 12740, 12747, 12754, 12761, 12768, 12775, 12782, 12789, 12796, 12803, 12810, 12817, 12824, 12831, 12838, 12845, 12852, 12859, 12866, 12873, 12880, 12887, 12894, 12901, 12908, 12915, 12922, 12929, 12936, 12943, 129 |  |  |  |  |  |   |  |



United States Environmental Protection Agency  
Contract Laboratory Program

**Organic Traffic Report  
& Chain of Custody Record**  
(For Organic CLP Analysis)

Case No. **28706**

|  |  |  |  |                          |   |
|--|--|--|--|--------------------------|---|
| 1. Project Code<br><b>100-2</b>        | 2. Region No.<br><b>2</b>  | Sampling Co.<br><b>Ray F. Weston</b>   | 4. Date Shipped<br><b>11/7/00</b>  | Carrier<br><b>Fed Ex</b> | 7. Preservative<br>(Enter in Column D)<br>1. HCl<br>2. HNO3<br>3. NaHSO4<br>4. H2SO4<br>5. Ice only<br>6. CH3OH<br>7. Other (specify in Column D)<br>N. Not Preserved |
| Account Code                           | Sampler (Name)<br><b>Diana Janda</b>   | Airbill Number<br><b>4455287301</b>  | 6. Matrix<br>(Enter in Column A)<br>1. Surface Water<br>2. Ground Water<br>3. Leachate<br>4. Field QC<br>5. Soil/Sediment<br>6. PE-water<br>7. PE-soil<br>8. Other (specify in Column A) |                          |   |
| Site Name<br><b>Lackawanna Foundry</b> | Sampler Signature<br><i>[Signature]</i>  | 5. Ship To:<br><b>American Analytical Technical Services, Inc.<br/>11950 Industrial Blvd<br/>Baton Rouge, LA 70809</b>                 |  |                          |   |
| City, State<br><b>Lackawanna, NY</b>   | 3. Purpose**<br>SF <input type="checkbox"/> PA <input type="checkbox"/> REIM <input type="checkbox"/> RI <input type="checkbox"/> ST <input type="checkbox"/> FED <input type="checkbox"/> EBZ <input checked="" type="checkbox"/> | Long-Term Action<br>RIFS <input type="checkbox"/> RD <input type="checkbox"/> RA <input type="checkbox"/> O&M <input type="checkbox"/> |  |                          |   |

| CLP Sample Numbers (from labels) | Matrix (from Box 6) Other: | B Conc.: Low Med | C Sample Type: Comp./Grab | D Preservative (from Box 7) Other: | E RAS Analysis                              |   | F Regional Specific Tracking Number or Tag Numbers | G Station Location Identifier | H Mo/Day/Year/Time Sample Collection | I Corresponding CLP Inorganic Sample No. | J Sampler Initials | K Field QC Qualifier |
|----------------------------------|----------------------------|------------------|---------------------------|------------------------------------|---|---|--|-------------------------------|--------------------------------------|--|--------------------|----------------------|
|                                  |                            |                  |                           |                                    | TA (circle one) PR* 7 14 (21) PR* 7 14 (21) | TA (circle one) PR* 7 14 (21) PR* 7 14 (21) |  |                               |                                      |  |                    |                      |
| B089                             | 5                          | Low              | Comp. N                   | N                                  | X   | X   | Tag Nos. 109-131                                   | 1002-SS-06C                   | 11/7/00 1310                         | MBAJ04                                   | BA                 |                      |
| B0913                            | 5                          | Low              | Comp. N                   | N                                  | X   | X   | Tag Nos. 105-107                                   | 1002-SS-05A                   | 11/7/00 1335                         | MBAJ58                                   | BA                 |                      |
| B0913                            | 5                          | Low              | Comp. N                   | N                                  | X   | X   | Tag Nos. 109-111                                   | 1002-SS-05B                   | 11/7/00 1340                         | MBAJ59                                   | BA                 |                      |
| B0985                            | 5                          | Low              | Comp. N                   | N                                  | X   | X   | Tag Nos. 113-115                                   | 1002-SS-05C                   | 11/7/00 1345                         | MBAJ60                                   | BA                 |                      |

Additional Sampler Signatures

Chain of Custody Seal Number(s)

Shipment for Case Complete? (Y/N) **Y** Page **2** of **2**

VOA MS/MSD Required? **Y/N** Sample #: \_\_\_\_\_

BNA MS/MSD Required? **Y/N** Sample #: \_\_\_\_\_

Pest/PCB MS/MSD Required? **Y/N** Sample #: \_\_\_\_\_

\*PR provides 7-day data turnaround in addition to preliminary results. Requests for preliminary results will increase analytical costs.

**Chain of Custody Record**

|  |                                    |   |             |
|--|------------------------------------|---|-------------|
| Relinquished by: (Signature)<br><i>[Signature]</i> | Date / Time<br><b>11/7/00 1730</b> | Received by: (Signature)                | Date / Time |
| Relinquished by: (Signature)                       | Date / Time                        | Received by: (Signature)                | Date / Time |
| Relinquished by: (Signature)                       | Date / Time                        | Received for Laboratory by: (Signature) | Date / Time |

Remarks: Is custody seal intact? **Y/N/none**







United States Environmental Protection Agency  
Contract Laboratory Program

### Inorganic Traffic Report & Chain of Custody Record

(For Inorganic CLP Analysis)

Case No. **28706**

1. Project Code: **1003** Account Code: \_\_\_\_\_  
 Regional Information: **Ray, Erikston**  
 Non-Superfund Program: **DDM, Kilda**  
 Site Name: **Lickhousen Ferryway** Site Spill ID: \_\_\_\_\_  
 City, State: **Lickhousen Ferryway**

2. Region No.: **2** Sampling Co.: **Ray, Erikston**  
 Sampler (Name): **DDM, Kilda**  
 Sampler Signature: *[Signature]*  
 3. Purpose:  SF  PRP  ST  FED

4. Date Shipped: **11/10/00** Carrier: **Fed Ex**  
 Airbill Number: **4455287234**  
 5. Ship To: **Chemtech Consulting Group**  
**Route Center**  
**205 Comp's Plaza 1**  
**Edison, NJ 08837**  
 ATTN: \_\_\_\_\_

6. Matrix (Enter in Column A)  
 1. Surface Water  
 2. Ground Water  
 3. Leachate  
 4. Field QC  
 5. Soil/Sediment  
 6. Oil (High only)  
 7. Waste (High only)  
 8. Other (Specify in Column A)

7. Preservative (Enter in Column D)  
 1. HCl  
 2. HNO3  
 3. NaOH  
 4. H2SO4  
 5. K2CR2O7  
 6. Ice only  
 7. Other (specify in Column D)  
 N. Not preserved

| CLP Sample Numbers (from labels) | A Matrix (from Box 6) | B Conc.: Low Med High | C Sample Type: Comp./Grab | D Preservative (from Box 7) | E - RAS Analysis |             |         |                 |          | G Station Location Identifier | H Mo/Day/Year/Time Sample Collection | I Corresponding CLP Organic Sample No. | J Sampler Initials | K Field QC Qualifier |
|----------------------------------|-----------------------|-----------------------|---------------------------|-----------------------------|------------------|-------------|---------|-----------------|----------|-------------------------------|--------------------------------------|--|--------------------|----------------------|
|                                  |                       |                       |                           |                             | Total Metals     | Diss Metals | Cyanide | NO <sub>3</sub> | Fluoride |                               |                                      |  |                    |                      |
| MANTJ64                          | 5                     | Low                   | Comp                      | N                           | XX               |             |         |                 |          | 1002-SS-02C                   | 11/7/00 1310                         | BWQ89                                  | BJ                 |                      |
| MANTJ58                          | 5                     | Low                   | Comp                      | N                           | XX               |             |         |                 |          | 1002-SS-02A                   | 11/7/00 1315                         | BWP12                                  | BJ                 |                      |
| MANTJ59                          | 5                     | Low                   | Comp                      | N                           | XX               |             |         |                 |          | 1002-SS-05B                   | 11/7/00 1340                         | BWP13                                  | BJ                 |                      |
| MANTJ60                          | 5                     | Low                   | Comp                      | N                           | XX               |             |         |                 |          | 1002-SS-05C                   | 11/7/00 1345                         | BWQ85                                  | BJ                 |                      |

Sample(s) to be Used for Laboratory QC: \_\_\_\_\_  
 Additional Sampler Signatures: \_\_\_\_\_  
 Chain of Custody Seal Number(s): \_\_\_\_\_  
 Page **2** of **2**

#### CHAIN OF CUSTODY RECORD

| Relinquished by: (Signature) | Date / Time  | Received by: (Signature) | Date / Time | Relinquished by: (Signature) | Date / Time | Received by: (Signature) | Date / Time |
|------------------------------|--------------|--------------------------|-------------|------------------------------|-------------|--------------------------|-------------|
| <i>[Signature]</i>           | 11/7/00 1730 |                          |             |                              |             |                          |             |
|                              |              |                          |             |                              |             |                          |             |
|                              |              |                          |             |                              |             |                          |             |



United States Environmental Protection Agency  
Contract Laboratory Program

### Inorganic Traffic Report & Chain of Custody Record

(For Inorganic CLP Analysis)

Case No. **28706**

|   |  |   |  |  |  |   |  |
|---|--|---|--|--|--|---|--|
| 1. Project Code<br><b>1002</b>                |  | 2. Region No./Sampling Co.<br><b>2</b> <i>Confidential</i>  |  | 4. Date Shipped<br><b>Fed Ex</b>   |  | 7. Preservative<br>(Enter in Column D)<br>1. HCl<br>2. HNO3<br>3. NaOH<br>4. H2SO4<br>5. K2CR2O7<br>6. Ice only<br>7. Other (specify in Column D)<br>N. Not preserved                                     |  |
| Regional Information                          |  | Sampler (Name)<br><b>Donna Jemel</b>  |  | Airbill Number<br><b>4955287285</b>  |  | 6. Matrix<br>(Enter in Column A)<br>1. Surface Water<br>2. Ground Water<br>3. Leachate<br>4. Field QC<br>5. Soil/Sediment<br>6. Oil (High only)<br>7. Waste (High only)<br>8. Other (specify in Column A) |  |
| Non-Superfund Program                         |  | Sampler Signature<br><i>[Signature]</i>   |  | 5. Ship To<br><b>Clientech Consulting Group<br/>Riverton Center<br/>205 Camps Plaza 1<br/>Edison, NJ 08837<br/>ATTN:</b> |  | H Mo/Day/Year/Time Sample Collection  |  |
| Site Name<br><b>Lackawanna Fencing</b>        |  | 3. Purpose<br><input type="checkbox"/> SF <input type="checkbox"/> PRP <input type="checkbox"/> ST <input type="checkbox"/> FED |  | F Regional Specific Tracking Number or Tag Numbers   |  | I Corresponding CLP Organic Sample No.  |  |
| City, State<br><b>Lackawanna, NY</b>          |  | Site Spill ID   |  | G Station Location Identifier  |  | J Sampler Initials  |  |
| CLP Sample Numbers (from labels)              |  | D Sample Preservative (from Box 7) Other:   |  | F Tracking Number or Tag Numbers   |  | K Field QC Qualifier<br>B = Blank S = Spike<br>D = Duplicate<br>R = Rinstate<br>PE = Perform Eval<br>N = Not a QC Sample  |  |
| A Matrix (from Box 6) Other:                  |  | E - RAS Analysis<br>Low only High only<br>Cyanide NO <sub>2</sub> -NO <sub>3</sub> Fluoride pH Conduct                          |  | F Tag No. 146<br>Tag No. 147   |  | H Mo/Day/Year/Time Sample Collection  |  |
| B Conc.: Low Med High                         |  | Diss. Metals Total Metals   |  | G 1002-RS-01<br>1002-RS-01   |  | I BAW92<br>BAW92  |  |
| C Sample Type: Comp./Grab (from Box 7) Other: |  | Lead SF PRP ST FED  |  | F Tag No. 146<br>Tag No. 147   |  | J Initials<br>AGG<br>AGG  |  |
| M.B.N.J.67 4 Low Grab 2                       |  |   |  |  |  |   |  |
| M.B.N.J.67 4 Low Grab 3                       |  |   |  |  |  |   |  |
|   |  |   |  |  |  |   |  |
|   |  |   |  |  |  |   |  |
|   |  |   |  |  |  |   |  |
|   |  |   |  |  |  |   |  |
| Shipment for Case Complete? (Y/N)             |  | Page <b>1</b> of <b>1</b>   |  | Additional Sampler Signatures  |  | Chain of Custody Seal Number(s)   |  |
|   |  |   |  |  |  |   |  |

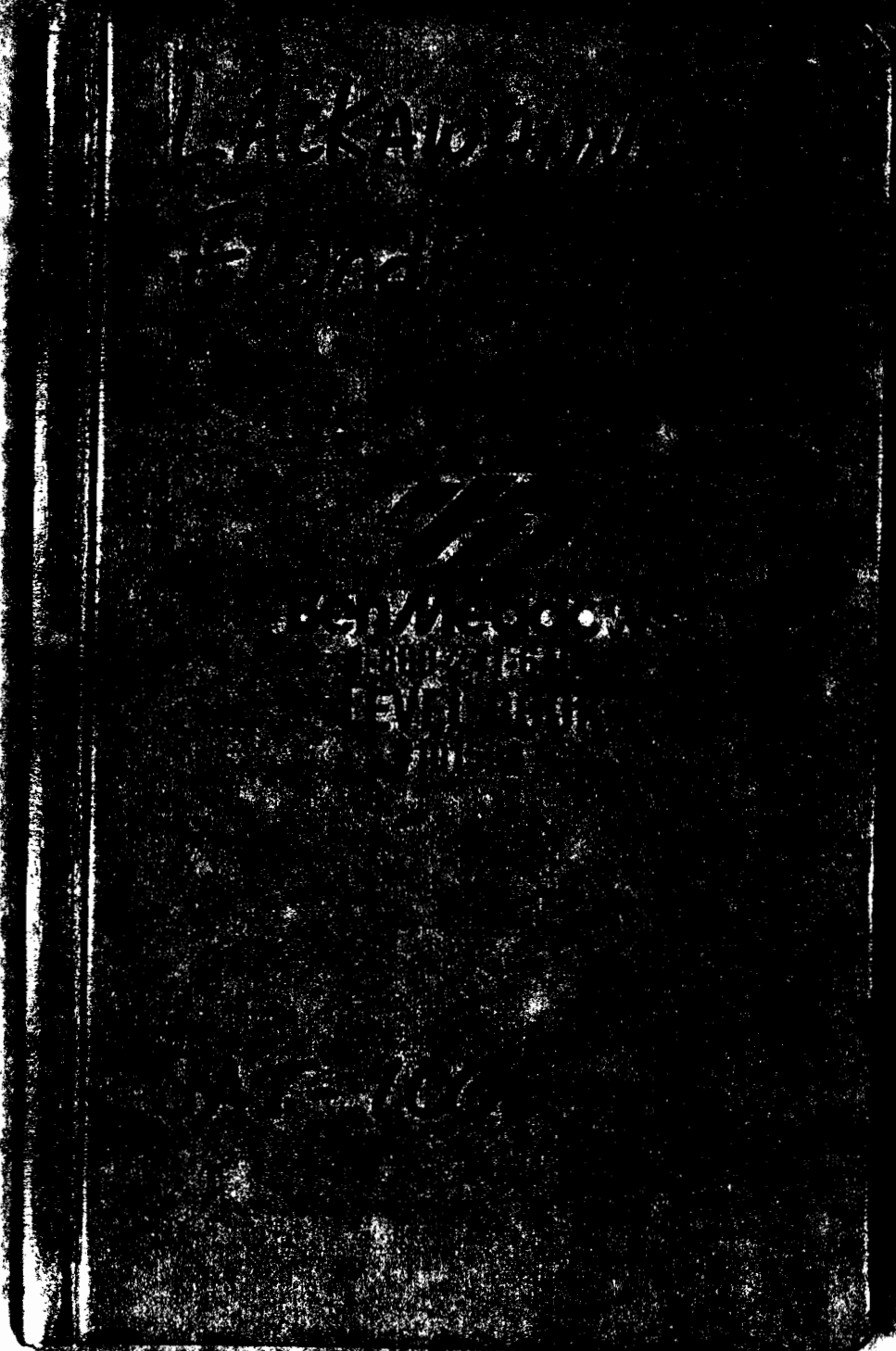
### CHAIN OF CUSTODY RECORD

|  |                                     |                          |             |
|--|-------------------------------------|--------------------------|-------------|
| Relinquished by: (Signature)<br><i>Donna Jemel</i> | Date / Time<br><b>11/16/00 1730</b> | Received by: (Signature) | Date / Time |
| Relinquished by: (Signature)                       | Date / Time                         | Received by: (Signature) | Date / Time |
| Relinquished by: (Signature)                       | Date / Time                         | Received by: (Signature) | Date / Time |

P-13



REFERENCE NO. 11





27 11/7/00

- 1105 Core taken at 6-8'. Blow count = w, 1, 2, 1  
Recovery = 1.5 of 2 feet. Over reading 0.0.
- 1115 Silty clay with a little brown soils. 1/4" samples.  
Dippers set up on third string location metals  
(SS0.3) ——— or
- 1115 0.0M reading 0.0 at base hole. — 5
- 1125 Core taken at 6.5 to 8.5'. Blow count =  
w, 4, 5, 7. Recovery 2 of 2 feet. 0.0M  
reading 0.0 Duplicate core taken at 1125.  
6.5 to 8.5'. Recovery 1.5 of 2 feet. 0.0M  
reading 0.0 Firm, grayish brown clay take  
VDA's, semis, metals (1) samples - SS0.3 + SS0.7  
All samples managed with CAP EPW protocol.  
Pipelines → Liberty, VT. Two Juncos →  
Baton Rouge, LA. ——— 2
- 1130 Lunch. Dippers prepare to steam clean  
auger & gears will collect inside blank after  
cleaning. ——— 2
- 1135 Binstate blank collected. 2 VDA vials,  
4 ambis for semis, 1 poly for metals, 1 bcd.  
VDA's preserved in HCL. Metals preserved  
in HNO<sub>3</sub> to pH < 2. ——— 2N  
Dippers setup on string for continuous  
SS0.6 sample ——— 2

D. J. [Signature]

3

11/7/00

- 1300 Core taken 2-4'. Blow count = 5, 4, 3. Recovery  
is 1 of 2 feet. 0.0M reading 0.0. Silt, clay  
brownish ——— 2
- 1305 Core taken 4-6'. Blow count = 1, 3, 3, 3.  
Recovery is 0 of 2 feet. ——— 2
- 1310 Core taken 6-8'. Blow count = 6, 6, 5, 3.  
Recovery 2 of 2 feet. Will use top 2 as 4-6 and  
bottom 2 as 6-8'. Brownish, gray clay  
1315 Core taken 8-10'. Very little recovery (~ 2")  
and bit water. No sample taken. (No. 2) clay  
silt. ——— 2
- 1320 Dippers set up on continuous string for SS0.6.
- 1325 Core taken 2-4'. Blow count = 2, 3, 4, 6  
Recovery = 2 of 2'. 0.0M readings 0.0  
Greenish, firm clay. ——— 2
- 1330 Core taken 4-6'. Blow count = 7, 1, 5, 3.  
Recovery = 1 of 2'. 0.0M reading 0.0.  
Green clay. ——— 2
- 1335 Core taken 6-8'. Blow count = 3, 7, 5, 5  
Recovery = 1.5 of 2'. 0.0M readings 0.0.  
Gray, silty clay ——— 2
- 1345 Core taken 8-10'. 0 of 2' recovery. 0.0  
sample taken ——— 2
- 1350 Dippers set up on bucket for continuous  
sample SS0.4. To be deep bedrock well.  
——— 2

D. J. [Signature]

4 11/7/00

- 1440 Core taken at 2-4'. Blow count = 1, 3, 5.
- Recovery = 1 of 2. Over reading 0.5
- Silty sand. Black at top in color.
- 1450 Core taken at 4-6'. Blow count = 2, 1, 5, 9
- Recovery = .5 of 2. Over reading 0.0.
- Brown, black silty sand.
- 1455 Core taken at 6-8'. Blow count = 7, 5, 1, 9.
- Recovery = 2 of 2. Over reading 0.0
- Gray, silty clay.
- 1515 Core taken at 8-10'. Blow count = 1, 5, 7, 10.
- Recovery = 1.5 of 2. Over reading 0.0.
- Gray, silty clay.
- 1515 Continue angers 0-8' angers to reach bedrock.
- 1105 30' of angers drilled - no bedrock reached.
- Drillers have no more 5' angers on truck.
- Will continue toward bedrock tomorrow.
- 1615 Drillers leave site.
- 1705 P. Lawitt, J. Anderson, J. Janke leave site.
- Bring samples to local FedEx to ship out to labs.
- Took six soil borings (3 continuous samplings) for UDA, SEM, metals. Samples shipped to EPA CUP labs. Angers sunk to 3' in prep for bedrock well installation.

9/20/00

11/17/00 weather: overcast ~45°

- 0750 Weston onsite. D. Lawitt / J. Anderson
- 0800 SJB (Drillers) onsite. Ken & Scott Teller
- 0815 Resume drilling for top of bedrock. Already at 30' ± 19" angers. Location is the sample 550%. Cuttings are 100% clay after 10 feet of angle.
- 0850 Rock hit at 38 feet. Soft, black shale.
- Angers advanced ~2' into bedrock. Are hole at 40' total depth. Will start from Ann
- 0915 hole to test rock competency.
- Spore from 40-42' feet containing weathered bedrock. Unconsolidated gravel, sand, + black shale. Will continue angering to reach more competent rock. Try to set steel casing at least 3' into compacted bedrock.
- 0945 Drillers have to pick up additional angers from office.
- 1010 Drillers return w/ 2 angers (8" x 5') Resume drilling from 40'.
- 1040 Anger hitting firmer resistance at ~42 feet. Paul slow advancement.
- 1056 45 Anger stopped at 43'. Attached roller bit to core into bedrock.
- 1115 EPA DSC K. Mathis + Region I EPA coord - Inscr. D. Pimball onsite.

9/20/00

6/1/81 1100

1130. Weston & EPA walk site and discuss progress & future plans, i.e. sampling on 117, well installations. Finish bedrock well today and go out. Get set for deep bedrock boring tomorrow. Begin installation of slatons with today. Drillers off-site to pick up compressor for air compressor line.

1145. EPA, OSC, Mattie's, lower site to meet with students. Will discuss document info with students using this site as example of environmental remediation.

1156. Driller's return to site.

1158. Reconn. EPA OSC Mattie's lower site

1200. Reconn. cutting of bedrock borehole in

after bit and air compression

All drill cuttings dumped in 6 by 30 drums.

1300 Problem with air compressor. Air not

completely flowing to bottom of borehole. Drill

bit may be stuck in upper casing.

1350. Reconn. cutting w/ roller bit.

1400. Auger bit stuck in borehole again. Pumping

causing problems. Start pouring into larger casing

and filling borehole.

1450. Reconn. casing from 13'. Using larger

casing for more torque & air flow.

1500

1400

1600. Ralk 411 engaged to 4 1/2". Approx. 6 feet below auger into bedrock. Will sink steel casing tomorrow and gravel. Airtight seal will come to 469' (20 feet further into bedrock) to obtain walk water production from bedrock zone.

1620 Weston & driller's leave site

Completed bedrock seal at 42'. Ralk bit came to 46' for installation of steel casing.

1500

- 8/14/00 weather: overcast, rain ~ 50"
- 0755 Weston D. Lewis & 5 JD's drillers on site  
Begin blasting standing water out of  
backrock with compressed air.  
well is producing water + overnight rain.
- 0840 Begin sinking steel casing into well to  
depth of 21' —
- 0900 2 sections of 21' each to be welded together  
+ additional length —
- 0900 Steel casing sunk to 49' by 5. Begin  
prep of grout —
- 1045 Begin grouting steel casing of backrock  
well —
- 1045 Backrock grouted - begin pulling augers. All  
more grout as augers are pulled up —
- 1046 Backrock grouted to top around steel  
casing. All augers pulled. waiting for  
grout to settle —
- 1105 Preparation for steam cleaning augers
- 1120 Add more grout to backrock well (GW01)
- 1215 Final steam clean of augers + truck
- 1230 Drillers set up on GW03 - over system  
well —
- 1240 Begin augering at GW03 - 7' augers  
Soil cuttings of 10' are gray clay. All cuttings  
being placed in 55g drums. —
- D. Lewis

- 1140
- 1300 Well at 20' depth. Cuttings are gray/slat  
clay. One driller leaves site to pick up 10  
slot screen + riser —
- 1350 10' of screen + riser in place. Begin sand  
filter grouting. PVC 50 schedule, 10 slot  
screen. —
- 1430 2 feet of bentonite seal in place —
- 1500 Bentonite - cement grout in place. GW03 well  
completed. Will set grout set with development  
GW03 20' deep, 18 foot screen - 10 feet below -  
(0.5 inch) sand pack to 2 feet above screen, 2 feet bentonite  
seal, grout to surface —
- 1530 Begin setting on overburden well GW05.  
Augering to 2' auger depth at 20'. Soil  
cuttings are brown clay. —
- 1600 Soil cuttings are gray clay at 20'. Rain  
coming down hard for last hour. —
- 1615 Well at 20'. Drillers install 10' of 10 slot  
screen + 10' feed off over - schedule 50  
PVC. —
- 1630 Sand filter (0 sand) installed to 1 foot  
above screen. Bentonite seal placed (2')  
Will grout tomorrow —
- 1840 Weston & drillers leave site
- D. Lewis

- 11/10/00  
1300 Sand filter (0 sand) up to 1 foot above screen
- 1295 GWOZ complet. 11 feet of sand filter pack, 6" 00 sand choke, 2 feet bentonite seal, 6" 00 sand choke, grout to grade
- 1400 Drills set up on overburden well G604.
- 1405 Begin angering to depth of 20'
- 1425 Cell cuttings are brown clay. After 10' cuttings become gray clay
- 1475 Well at 20' depth
- 1500 10' of 10 slot screen, 10' riser installed in well - 70 schedule PVC.
- 1505 Begin pouring 0 sand filter pack to 1 foot above screen. Pulling augers as sand is poured
- 1545 Well GWOZ completed. 11 feet of sand filter pack, 6" 00 sand choke, 2' bentonite seal, 6" 00 sand choke, grout to grade
- 1600 Drills begin cleanup. Will develop shallow wells, install curb boxes, & drum remaining
- 1605 well cuttings Monday 11/17
- 1630 Weirton & dr:llors leave site
- Cond bed rock well to 67', installed two overburden wells to 20' each

Edgett

- 11/10/00 weather: overcast, very windy ~ 95°
- 1300 Weirton 2 levels + 500 drillers onsite
- 1330 Finish grouting GWOZ. Jet set for double point on Monday
- 1375 Drills set up on bedrock well G601. Will core through bedrock with roller-bit from 49' until water is reached.
- First clear well of standing water is on compressor
- 1395 Begin coring into bedrock - roller-bit engaged
- 1330 Well producing water after 4' of cutting. (total depth = 53')
- 1400 Continue roller-bit coring into bedrock.
- 1435 Boring into bedrock completed. Final well depth = 67'. 20' total feet into bedrock. Well producing water (~ 1/2 in) since 53'
- 1445 Begin pulling up rods & roller-bit from well
- 1450 Set up drill rig on GWOZ - overburden well. Begin angering to depth of 20'.
- 1450 Drill cuttings are brown clay and after 12' go gray
- 1475 Well at 20' depth
- 1480 Level
- 1530 Resume work on GWOZ. Sink 10' of 10 slot screen, 10 feet of PVC - 70 schedule PVC

Edgett



12 11/13/00 water: increased ~ 25"

0730 Wesleyan D. Conitt on site. Awaiting 530 dollars

0820 553 on site. Scott + Ken Fuller

0830 Drillers set up on G403. Only a few inches of water in well. Can not develop

0940 Remaining soil cuttings are drummed

2900 Set up on G405. ~ 6' of water in well

0910 Begin development of G405.

0630 Water clears up after ~ 15 minutes.

Pumping volume decreasing. Shut off pump to allow well to recover. Already pumped ~ 3 well volumes.

0946 Soil cuttings completely drummed

1000 G406 brought to grade in cement. Cover installed 1140. Still developing well & waiting for cement to harden.

1020 G403 brought to grade in cement and cover installed

1030 Site development on G406 (pumped ~ 60g for 1 hour 15 min). Water slightly turbid probably from clay

1040 Set up on G407. Approximately 6 feet of water in well. Begin development and drumming of soil cuttings.

11/12/00

13

1166 Site development on G404. Pumped ~ 20g water into ground. Soil slightly turbid.

G404 brought to grade in cement. Cover installed. All soil cuttings drummed.

143 Set up on G402. Very little water & weak.

1200 Curb box installed on G401 + G402

1430 G401 + G402 brought to grade in cement.

Curb boxes installed. All soil cuttings placed in 55g drums

1500 Drillers clean up rest of their material on site and leave. Wesley leaves site & locks up gate

Installed 5 curb boxes and cemented to grade - 11 well (G401 - G405). All soil

cuttings placed in 55g drums on site for future disposal. G402 + G403 are too dry to develop. G404 + G405 each developed for 1 hour + 15 minutes.

Q20A

10/10/00 weather sunny + windy 10-20  
0730 Weston E. Sa/920 - 2 D. L. 1000

0745 on site  
0755 1.45 m. 0.3 static water level taken  
9:45 ft

0800 level goes for general - begin setting o-  
0805 4.2.3 for groundwater

| Water | Depth | Count | Temp | Depth | Temp |
|-------|-------|-------|------|-------|------|
| 0835  | 11.75 | 6.3   | 72.7 | 12.8  | 0.7  |
| 0840  | 22.4  | 6.5   | 76.1 | 13.15 | 20   |
| 0845  | 33.5  | 6.3   | 76.3 | 13.2  | 2.5  |
| 0850  | 44.0  | 6.5   | 76.5 | 13.2  | 3.5  |

0855 Collected sample for 0855  
0905 set up with CWS  
0915 Static water level = 102 ft

| Water | Depth | Count | Temp | Depth | Temp |
|-------|-------|-------|------|-------|------|
| 0920  | 8.75  | 6.4   | 76.2 | 14    | 10.2 |
| 0925  | 9.72  | 6.7   | 76.7 | 16.2  | 11.5 |
| 0930  | 7.79  | 6.7   | 76.9 | 17.7  | 11.5 |
| 0935  | 9.42  | 6.5   | 76.5 | 18.1  | 12.6 |
| 0940  | 8.79  | 6.5   | 76.9 | 19    | 12.8 |

0945 collect sample & CWS  
0955 set up on ground  
Static water level = 89.8 ft

| Water | Depth | Count | Temp | Depth | Temp |
|-------|-------|-------|------|-------|------|
| 1000  | 11.75 | 6.3   | 76.1 | 13.15 | 20   |
| 1005  | 22.4  | 6.5   | 76.1 | 13.15 | 20   |
| 1010  | 33.5  | 6.3   | 76.3 | 13.2  | 2.5  |
| 1015  | 44.0  | 6.5   | 76.5 | 13.2  | 3.5  |

1020 collect sample & CWS  
1030 set up on ground  
Static water level = 89.8 ft

1035

1040 water level count 1040 1040 1040

1045 10.54 6.4 62.8 60 11.45 17 95.0  
1050 11.04 6.4 62.6 61 11.37 10.3 90.0

1055 11.14 6.1 62.7 5.2 11.31 10.1 88.2  
1100 11.24 6.4 62.4 4.7 11.03 12.6 82.3

1105 11.29 6.4 62.0 5.8 11.03 10.7 86.6  
1110 11.40 6.5 63.2 3.3 10.25 10.8 80.3

1100 Collect sample CWS  
1130 Lunch  
1220 set up 2 CWS

Static water level = 17.81 ft  
0.75  
13.8 6.7 90.9 5.0 10.5 7.5  
1400 14.3 6.7 97.4 3.0 10.5 7.4

Well run dry. did not reach CWS  
to collect sample later

1515 set up 2 CWS  
Static water level = 12.81 ft

1525 13.71 6.2 2.9 2.4 11.1 3  
1530 13.71 6.2 2.9 2.4 11.1 3

1535 13.81 6.7 3.5 2.1 11.1 7.7  
1400 14.3 6.7 3.5 2.1 11.1 7.7

1540 14.3 6.7 3.5 2.1 11.1 7.7

1545 14.3 6.7 3.5 2.1 11.1 7.7

1550 14.3 6.7 3.5 2.1 11.1 7.7

1555 14.3 6.7 3.5 2.1 11.1 7.7

1600 14.3 6.7 3.5 2.1 11.1 7.7

16  
12/1/00

1100 Collect sample C-201, m.s. in SP, + Dye

~~1100~~ \_\_\_\_\_

1120 lower did not retrace. can not correct

oil sample from C-202 \_\_\_\_\_

1130 Collect sample from Skunk C-302 \_\_\_\_\_

1130 Cleanup + back wells. D. Leming + E.

Salvage leave site \_\_\_\_\_

3 shallow wells + 1 bedrock well

sampled for TEL + TAC water samples.

PCB/PCIT + CN in depth. C-202 not

sampled but well dry-up \_\_\_\_\_

D. Leming

REFERENCE NO. 12

**APPENDIX A of TAGM #4046**

**TABLE 1**  
**Recommended soil cleanup objectives (mg/kg or ppm)**  
**Volatile Organic Contaminants**

Shortcut to TAGM 4046 Tables for SVOCs | Pesticides/PCBs | Heavy Metals

| Contaminant                                       | Partition Coefficient, Koc | Groundwater Standards/ Criteria, Cw (ug/l or ppb) | a Allowable soil conc., Cs (ppm) | b ** Soil cleanup objectives to protect GW quality (ppm) | USEPA Health Based (ppm) |                    | CRQL (ppb) | *** Rec. Soil Cleanup Objective (ppm) |
|---|----------------------------|---|----------------------------------|--|--------------------------|--------------------|------------|---------------------------------------|
|   |                            |   |                                  |  | Carcinogens              | Systemic Toxicants |            |                                       |
| Acetone   | 2.2                        | 50  | 0.0011                           | 0.11   | N/A                      | 8,000              | 10         | 0.2                                   |
| Benzene   | 83                         | 0.7   | 0.0006                           | 0.06   | 24                       | N/A                | 5          | 0.06                                  |
| Benzoic Acid                                      | 54 *                       | 50  | 0.027                            | 2.7  | N/A                      | 300,000            | 5          | 2.7                                   |
| 2-Butanone  | 4.5 *                      | 50  | 0.003                            | 0.3  | N/A                      | 4,000              | 10         | 0.3                                   |
| Carbon Disulfide                                  | 54 *                       | 50  | 0.027                            | 2.7  | N/A                      | 8,000              | 5          | 2.7                                   |
| Carbon Tetrachloride                              | 110 *                      | 5   | 0.006                            | 0.6  | 5.4                      | 60                 | 5          | 0.6                                   |
| Chlorobenzene                                     | 330                        | 5   | 0.017                            | 1.7  | N/A                      | 2,000              | 5          | 1.7                                   |
| Chloroethane                                      | 37 *                       | 50  | 0.019                            | 1.9  | N/A                      | N/A                | 10         | 1.9                                   |
| Chloroform  | 31                         | 7   | 0.003                            | 0.30   | 114                      | 800                | 5          | 0.3                                   |
| Dibromochloromethane                              | N/A                        | 50  | N/A                              | N/A  | N/A                      | N/A                | 5          | N/A                                   |
| 1,2-Dichlorobenzene                               | 1,700                      | 4.7   | 0.079                            | 7.9  | N/A                      | N/A                | 330        | 7.9                                   |
| 1,3-Dichlorobenzene                               | 310 *                      | 5   | 0.0155                           | 1.55   | N/A                      | N/A                | 330        | 1.6                                   |
| 1,4-Dichlorobenzene                               | 1,700                      | 5   | 0.085                            | 8.5  | N/A                      | N/A                | 330        | 8.5                                   |
| 1,1-Dichloroethane                                | 30                         | 5   | 0.002                            | 0.2  | N/A                      | N/A                | 5          | 0.2                                   |
| 1,2-Dichloroethane                                | 14                         | 5   | 0.001                            | 0.1  | 7.7                      | N/A                | 5          | 0.1                                   |
| 1,1-Dichloroethene                                | 65                         | 5   | 0.004                            | 0.4  | 12                       | 700                | 5          | 0.4                                   |
| 1,2-Dichloroethene (trans)                        | 59                         | 5   | 0.003                            | 0.3  | N/A                      | 2,000              | 5          | 0.3                                   |
| 1-3 dichloropropane                               | 51                         | 5   | 0.003                            | 0.3  | N/A                      | N/A                | 5          | 0.3                                   |
| Ethylbenzene                                      | 1,100                      | 5   | 0.055                            | 5.5  | N/A                      | 8,000              | 5          | 5.5                                   |
| 113 Freon (1,1,2 Trichloro-1,2,2 Trifluoroethane) | 1,230 *                    | 5   | 0.060                            | 6.0  | N/A                      | 200,000            | 5          | 6.0                                   |
| Methylene chloride                                | 21                         | 5   | 0.001                            | 0.1  | 93                       | 5,000              | 5          | 0.1                                   |
| 4-Methyl-2-Pentanone                              | 19 *                       | 50  | 0.01                             | 1.0  | N/A                      | N/A                | 10         | 1.0                                   |
| Tetrachloroethene                                 | 277                        | 5   | 0.014                            | 1.4  | 14                       | 800                | 5          | 1.4                                   |
| 1,1,1-Trichloroethane                             | 152                        | 5   | 0.0076                           | 0.76   | N/A                      | 7,000              | 5          | 0.8                                   |

0.1

| 1,1,1-trichloroethane     | 118   | 5 | 0.0070 | 0.70 | N/A | 1,000   | 5   | 0.8 |
|---------------------------|-------|---|--------|------|-----|---------|-----|-----|
| 1,1,2,2-Tetrachloroethane | 118   | 5 | 0.006  | 0.6  | 35  | N/A     | 5   | 0.6 |
| 1,2,3-trichloropropane    | 68    | 5 | 0.0034 | 0.34 | N/A | 80      | 5   | 0.4 |
| 1,2,4-trichlorobenzene    | 670 * | 5 | 0.034  | 3.4  | N/A | N/A     | 330 | 3.4 |
| Toluene                   | 300   | 5 | 0.015  | 1.5  | N/A | 20,000  | 5   | 1.5 |
| Trichloroethene           | 126   | 5 | 0.007  | 0.70 | 64  | N/A     | 5   | 0.7 |
| Vinyl chloride            | 57    | 2 | 0.0012 | 0.12 | N/A | N/A     | 10  | 0.2 |
| Xylenes                   | 240   | 5 | 0.012  | 1.2  | N/A | 200,000 | --  | 1.2 |

- a. Allowable Soil Concentration  $C_s = f \times C_w \times K_{oc}$
  - b. Soil cleanup objective =  $C_s \times$  Correction Factor (CF)
- N/A is not available

- \* Partition coefficient is calculated by using the following equation:  
 $\log K_{oc} = -0.55 \log S + 3.64$ , where S is solubility in water in ppm.  
 All other  $K_{oc}$  values are experimental values.
- \*\* Correction Factor (CF) of 100 is used as per TAGM #4046
- \*\*\* As per TAGM #4046, Total VOCs < 10 ppm.

Note: Soil cleanup objectives are developed for soil organic carbon content (f) of 1%, and should be adjusted for the actual soil organic carbon content if it is known.

**APPENDIX A of TAGM #4046**

**TABLE 2**  
**Recommended soil cleanup objectives (mg/kg or ppm)**  
**Semi-Volatile Organic Contaminants**

Shortcut to TAGM 4046 Tables for VOCs | Pesticides/PCBs | Heavy Metals

| Contaminant                 | Partition Coefficient, Koc | Groundwater Standards/ Criteria, Cw (ug/l or ppb) | a Allowable soil conc., Cs (ppm) | b ** Soil cleanup objectives to protect GW quality (ppm) | USEPA Health Based (ppm) |                    | CRQL (ppb) | *** Rec. Soil Cleanup Objective (ppm) |
|-----------------------------|----------------------------|---|----------------------------------|--|--------------------------|--------------------|------------|---------------------------------------|
|                             |                            |   |                                  |  | Carcin-ogens             | Systemic Toxicants |            |                                       |
| Acenaphthene                | 4,600                      | 20  | 0.9                              | 90.0   | N/A                      | 5,000              | 330        | 50.0<br>***                           |
| Acenaphthylene              | 2,056 *                    | 20  | 0.41                             | 41.0   | N/A                      | N/A                | 330        | 41.0                                  |
| Aniline                     | 13.8                       | 5   | 0.001                            | 0.1  | 123                      | N/A                | 330        | 0.1                                   |
| Anthracene                  | 14,000                     | 50  | 7.00                             | 700.0  | N/A                      | 20,000             | 330        | 50.0<br>***                           |
| Benzo(a)anthracene          | 1,380,000                  | 0.002   | 0.03                             | 3.0  | 0.224                    | N/A                | 330        | 0.224<br>or<br>MDL                    |
| Benzo (a) pyrene            | 5,500,000                  | 0.002 (ND)  | 0.110                            | 11.0   | 0.0609                   | N/A                | 330        | 0.061<br>or<br>MDL                    |
| Benzo (b) fluoranthene      | 550,000                    | 0.002   | 0.011                            | 1.1  | N/A                      | N/A                | 330        | 1.1                                   |
| Benzo (g,h,i) perylene      | 1,600,000                  | 5   | 8.0                              | 800  | N/A                      | N/A                | 330        | 50.0<br>***                           |
| Benzo (k) fluoranthene      | 550,000                    | 0.002   | 0.011                            | 1.1  | N/A                      | N/A                | 330        | 1.1                                   |
| bis(2-ethylhexyl) phthalate | 8,706 *                    | 50  | 4.35                             | 435.0  | 50                       | 2,000              | 330        | 50.0<br>***                           |
| Butylbenzylphthlate         | 2,430                      | 50  | 1.215                            | 122.0  | N/A                      | 20,000             | 330        | 50.0<br>***                           |
| Chrysene                    | 200,000                    | 0.002   | 0.004                            | 0.4  | N/A                      | N/A                | 330        | 0.4                                   |
| 4- Chloroaniline            | 43 ****                    | 5   | 0.0022                           | 0.22   | 200                      | 300                | 330        | 0.220<br>or<br>MDL                    |
| 4-Chloro-3-methylphenol     | 47                         | 5   | 0.0024                           | 0.24   | N/A                      | N/A                | 330        | 0.240<br>or<br>MDL                    |
| 2-Chlorophenol              | 15 *                       | 50  | 0.008                            | 0.8  | N/A                      | 400                | 330        | 0.8                                   |

|                          |                        |       |        |         |        |        |       |                    |
|--------------------------|------------------------|-------|--------|---------|--------|--------|-------|--------------------|
| Dibenzofuran             | 1,230 *                | 5     | 0.062  | 6.2     | N/A    | N/A    | 330   | 6.2                |
| Dibenzo(a,h)anthracene   | 33,000,000             | 50    | 1,650  | 165,000 | 0.0143 | N/A    | 330   | 0.014<br>or<br>MDL |
| 3,3'-Dichlorobenzidine   | N/A                    | N/A   | N/A    | N/A     | N/A    | N/A    | N/A   | N/A                |
| 2,4-Dichlorophenol       | 380                    | 1     | 0.004  | 0.4     | N/A    | 200    | 330   | 0.4                |
| 2,4-Dinitrophenol        | 38                     | 5     | 0.002  | 0.2     | N/A    | 200    | 1,600 | 0.200<br>or<br>MDL |
| 2,6 Dinitrotoluene       | 198*                   | 5     | 0.01   | 1.0     | 1.03   | N/A    | 330   | 1.0                |
| Diethylphthlate          | 142                    | 50    | 0.071  | 7.1     | N/A    | 60,000 | 330   | 7.1                |
| Dimethylphthlate         | 40                     | 50    | 0.020  | 2.0     | N/A    | 80,000 | 330   | 2.0                |
| Di-n-butyl phthalate     | 162*                   | 50    | 0.081  | 8.1     | N/A    | 8,000  | 330   | 8.1                |
| Di-n-octyl phthlate      | 2,346 *                | 50    | 1.2    | 120.0   | N/A    | 2,000  | 330   | 50.0<br>***        |
| Fluoranthene             | 38,000                 | 50    | 19     | 1900.0  | N/A    | 3,000  | 330   | 50.0<br>***        |
| Fluorene                 | 7,300                  | 50    | 3.5    | 350.0   | N/A    | 3,000  | 330   | 50.0<br>***        |
| Hexachlorobenzene        | 3,900                  | 0.35  | 0.014  | 1.4     | 0.41   | 60     | 330   | 0.41               |
| Indeno (1,2,3-cd) pyrene | 1,600,000              | 0.002 | 0.032  | 3.2     | N/A    | N/A    | 330   | 3.2                |
| Isophorone               | 88.31 *                | 50    | 0.044  | 4.40    | 1,707  | 20,000 | 330   | 4.40               |
| 2-methylnaphthalene      | 727 *                  | 50    | 0.364  | 36.4    | N/A    | N/A    | 330   | 36.4               |
| 2-Methylphenol           | 15                     | 5     | 0.001  | 0.1     | N/A    | N/A    | 330   | 0.100<br>or<br>MDL |
| 4-Methylphenol           | 17                     | 50    | 0.009  | 0.9     | N/A    | 4,000  | 330   | 0.9                |
| Naphthalene              | 1,300                  | 10    | 0.130  | 13.0    | N/A    | 300    | 330   | 13.0               |
| Nitrobenzene             | 36                     | 5     | 0.002  | 0.2     | N/A    | 40     | 330   | 0.200<br>or<br>MDL |
| 2-Nitroaniline           | 86                     | 5     | 0.0043 | 0.43    | N/A    | N/A    | 1,600 | 0.430<br>or<br>MDL |
| 2-Nitrophenol            | 65                     | 5     | 0.0033 | 0.33    | N/A    | N/A    | 330   | 0.330<br>or<br>MDL |
| 4-Nitrophenol            | 21                     | 5     | 0.001  | 0.1     | N/A    | N/A    | 1,600 | 0.100<br>or<br>MDL |
| 3-Nitroaniline           | 93<br><del>88.31</del> | 5     | 0.005  | 0.5     | N/A    | N/A    | 1,600 | 0.500<br>OR MDL    |



| Contaminant           | Concentration (ppm) | Soil Cleanup Objective (ppm) | Soil Organic Carbon Content (f) | Soil Cleanup Objective (ppm) | Soil Cleanup Objective (ppm) | Soil Cleanup Objective (ppm) | Soil Cleanup Objective (ppm) | Soil Cleanup Objective (ppm) |
|-----------------------|---------------------|------------------------------|---------------------------------|------------------------------|------------------------------|------------------------------|------------------------------|------------------------------|
| Pentachlorophenol     | 1,022               | 1                            | 0.01                            | 1.0                          | N/A                          | 2,000                        | 1,600                        | 1.0 or MDL                   |
| Phenanthrene          | 4,365 *             | 50                           | 2.20                            | 220.0                        | N/A                          | N/A                          | 330                          | 50.0 ***                     |
| Phenol                | 27                  | 1                            | 0.0003                          | 0.03                         | N/A                          | 50,000                       | 330                          | 0.03 or MDL                  |
| Pyrene                | 13,295 *            | 50                           | 6.65                            | 665.0                        | N/A                          | 2,000                        | 330                          | 50.0 ***                     |
| 2,4,5-Trichlorophenol | 89 *                | 1                            | 0.001                           | 0.1                          | N/A                          | 8,000                        | 330                          | 0.1                          |

- a. Allowable Soil Concentration  $C_s = f \times C_w \times K_{oc}$   
b. Soil cleanup objective =  $C_s \times$  Correction Factor (CF)  
N/A is not available  
MDL is Method Detection Limit

- \* Partition coefficient is calculated by using the following equation:  
 $\log K_{oc} = -0.55 \log S + 3.64$ , where S is solubility in water in ppm.  
Other  $K_{oc}$  values are experimental values.
- \*\* Correction Factor (CF) of 100 is used as per TAGM #4046
- \*\*\* As per TAGM #4046, Total VOCs < 10 ppm., Total Semi- VOCs < 500ppm. and Individual Semi-VOCs < 50 ppm.
- \*\*\*\*  $K_{oc}$  is derived from the correlation  $K_{oc} = 0.63 K_{ow}$  (Determining Soil Response Action Levels..... EPA/540/2-89/057).  $K_{ow}$  is obtained from the USEPA computer database 'MAIN'.

Note: Soil cleanup objectives are developed for soil organic carbon content (f) of 1%, and should be adjusted for the actual soil organic carbon content if it is known.

**APPENDIX A of TAGM #4046**

**TABLE 3**  
**Recommended soil cleanup objectives (mg/kg or ppm)**  
**Organic Pesticides / Herbicides and PCBs**

Shortcut to TAGM 4046 Tables for [VOCs](#) | [SVOCs](#) | [Heavy Metals](#)

| Contaminant                           | Partition Coefficient, Koc | Groundwater Standards/ Criteria, Cw (ug/l or ppb) | a Allowable soil conc., Cs (ppm) | b ** Soil cleanup objectives to protect GW quality (ppm) | USEPA Health Based (ppm) |                    | CRQL (ppb) | *** Rec. Soil Cleanup Objective (ppm) |
|---------------------------------------|----------------------------|---|----------------------------------|--|--------------------------|--------------------|------------|---------------------------------------|
|                                       |                            |   |                                  |  | Carcinogens              | Systemic Toxicants |            |                                       |
| Aldrin                                | 96,000                     | ND (<0.01)  | 0.005                            | 0.5  | 0.041                    | 2                  | 8          | 0.041                                 |
| alpha- BHC                            | 3,800                      | ND (<0.05)  | 0.002                            | 0.2  | 0.111                    | N/A                | 8          | 0.11                                  |
| beta - BHC                            | 3,800                      | ND (<0.05)  | 0.002                            | 0.2  | 3.89                     | N/A                | 8          | 0.2                                   |
| delta - BHC                           | 6,600                      | ND (<0.05)  | 0.003                            | 0.3  | N/A                      | N/A                | 8          | 0.3                                   |
| Chlordane                             | 21,305 *                   | 0.1   | 0.02                             | 2.0  | 0.54                     | 50                 | 80         | 0.54                                  |
| 2,4-D                                 | 104 *                      | 4.4   | 0.005                            | 0.5  | N/A                      | 800                | 800        | 0.5                                   |
| 4,4'- DDD                             | 770,000 *                  | ND (<0.01)  | 0.077                            | 7.7  | 2.9                      | N/A                | 16         | 2.9                                   |
| 4,4'-DDE                              | 440,000 *                  | ND (<0.01)  | 0.0440                           | 4.4  | 2.1                      | N/A                | 16         | 2.1                                   |
| 4,4'-DDT                              | 243,000 *                  | ND (<0.01)  | 0.025                            | 2.5  | 2.1                      | 40                 | 16         | 2.1                                   |
| Dibenzo-P-dioxins (PCDD) 2,3,7,8 TCDD | 1709800                    | 0.000035  | 0.0006                           | 0.06   | N/A                      | N/A                | N/A        | N/A                                   |
| Dieldrin                              | 10,700 *                   | ND (<0.01)  | 0.0010                           | 0.1  | 0.044                    | 4                  | 16         | 0.044                                 |
| Endosulfan I                          | 8,168 *                    | 0.1   | 0.009                            | 0.9  | N/A                      | N/A                | 16         | 0.9                                   |
| Endosulfan II                         | 8,031 *                    | 0.1   | 0.009                            | 0.9  | N/A                      | N/A                | 16         | 0.9                                   |
| Endosulfan Sulfate                    | 10,038 *                   | 0.1   | 0.01                             | 1.0  | N/A                      | N/A                | 16         | 1.0                                   |
| Endrin                                | 9,157 *                    | ND (<0.01)  | 0.001                            | 0.1  | N/A                      | 20                 | 8          | 0.10                                  |
| Endrin keytone                        | N/A                        | N/A   | N/A                              | N/A  | N/A                      | N/A                | N/A        | N/A                                   |
| gamma - BHC                           |                            | ND  |                                  |  |                          |                    |            |                                       |

|  |          |                   |        |      |       |     |     |                                       |
|--|----------|-------------------|--------|------|-------|-----|-----|---------------------------------------|
| (Lindane)                                    | 1,080    | ND<br>( $<0.05$ ) | 0.0006 | 0.06 | 5.4   | 20  | 8   | 0.06                                  |
| gamma -<br>chlordane                         | 140,000  | 0.1               | 0.14   | 14.0 | 0.54  | 5   | 80  | 0.54                                  |
| Heptachlor                                   | 12,000   | ND<br>( $<0.01$ ) | 0.0010 | 0.1  | 0.16  | 40  | 8   | 0.10                                  |
| Heptachlor<br>epoxide                        | 220      | ND<br>( $<0.01$ ) | 0.0002 | 0.02 | 0.077 | 0.8 | 8   | 0.02                                  |
| Methoxychlor                                 | 25,637   | 35.0              | 9.0    | 900  | N/A   | 400 | 80  | ***                                   |
| Mitotane                                     | N/A      | N/A               | N/A    | N/A  | N/A   | N/A | N/A | N/A                                   |
| Parathion                                    | 760      | 1.5               | 0.012  | 1.2  | N/A   | 500 | 8   | 1.2                                   |
| PCBs   | 17,510 * | 0.1               | 0.1    | 10.0 | 1.0   | N/A | 160 | 1.0<br>(Surface)<br>10 (sub-<br>surf) |
| Polychlorinated<br>dibenzo-<br>furans (PCDF) | N/A      | N/A               | N/A    | N/A  | N/A   | N/A | N/A | N/A                                   |
| Silvex                                       | 2,600    | 0.26              | 0.007  | 0.7  | N/A   | 600 | 330 | 0.7                                   |
| 2,4,5-T                                      | 53       | 35                | 0.019  | 1.9  | N/A   | 200 | 330 | 1.9                                   |

- a. Allowable Soil Concentration  $C_s = f \times C_w \times K_{oc}$   
b. Soil cleanup objective =  $C_s \times$  Correction Factor (CF)  
N/A is not available

\* Partition coefficient is calculated by using the following equation:  
 $\log K_{oc} = -0.55 \log S + 3.64$ , where S is solubility in water in ppm.  
All other  $K_{oc}$  values are experimental values.

\*\* Correction Factor (CF) of 100 is used as per TAGM #4046

\*\*\* As per TAGM #4046, Total VOCs  $< 10$  ppm.

Note: Soil cleanup objectives are developed for soil organic carbon content (f) of 1% (5% for PCBs as per PCB Guidance Document), and should be adjusted for the actual soil organic carbon content if it is known.

**APPENDIX A of TAGM #4046**

**TABLE 4**  
**Recommended soil cleanup objectives (mg/kg or ppm)**  
**Heavy Metals**

Shortcut to TAGM 4046 Tables for [VOCs](#) | [SVOCs](#) | [Pesticides](#) / [PCBs](#)

| Contaminants | Protect Water Quality (ppm) | Eastern USA Background (ppm) | * CRDL (mg/kg or ppm) | ***** Rec. Soil Cleanup Objective (ppm) |
|--------------|-----------------------------|------------------------------|-----------------------|---|
| Aluminum     | N/A                         | 33,000                       | 2.0                   | SB                                      |
| Antimony     | N/A                         | N/A                          | 0.6                   | SB                                      |
| Arsenic      | N/A                         | 3-12 **                      | 0.1                   | 7.5 or SB                               |
| Barium       | N/A                         | 15-600                       | 2.0                   | 300 or SB                               |
| Beryllium    | N/A                         | 0-1.75                       | 0.05                  | 0.16 (HEAST) or SB                      |
| Cadmium      | N/A                         | 0.1-1                        | 0.05                  | 1 or SB                                 |
| Calcium      | N/A                         | 130 - 35,000 ***             | 50.0                  | SB                                      |
| Chromium     | N/A                         | 1.5 - 40 **                  | 0.1                   | 10 or SB                                |
| Cobalt       | N/A                         | 2.5 - 60 **                  | 0.5                   | 30 or SB                                |
| Copper       | N/A                         | 1 - 50                       | 0.25                  | 25 or SB                                |
| Cyanide      | N/A                         | N/A                          | 0.1                   | *****                                   |
| Iron         | N/A                         | 2,000 - 550,000              | 1.0                   | 2,000 or SB                             |
| Lead         | N/A                         | *****                        | 0.03                  | SB *****                                |
| Magnesium    | N/A                         | 100 - 5,000                  | 50.0                  | SB                                      |
| Manganese    | N/A                         | 50 - 5,000                   | 0.15                  | SB                                      |
| Mercury      | N/A                         | 0.001 - 0.2                  | 0.002                 | 0.1                                     |
| Nickel       | N/A                         | 0.5 -25                      | 0.4                   | 13 or SB                                |
| Potassium    | N/A                         | 8,500 - 43,000 **            | 50.0                  | SB                                      |
| Selenium     | N/A                         | 0.1 - 3.9                    | 0.05                  | 2 or SB                                 |
| Silver       | N/A                         | N/A                          | 0.1                   | SB                                      |
| Sodium       | N/A                         | 6,000 - 8,000                | 50.0                  | SB                                      |
| Thallium     | N/A                         | N/A                          | 0.1                   | SB                                      |
| Vanadium     | N/A                         | 1-300                        | 0.5                   | 150 or SB                               |
| Zinc         | N/A                         | 9-50                         | 0.2                   | 20 or SB                                |

Note: Some forms of metal salts such as Aluminum Phosphide, Calcium Cyanide, Potassium Cyanide, Copper cyanide, Silver cyanide, Sodium cyanide, Zinc phosphide, Thallium salts, Vanadium pentoxide and Chromium (VI) compounds are more toxic in nature. Please refer to the USEPA HEASTs database to find cleanup objectives if such metals are present in soil.

SB is site background

N/A is not available

\* CRDL is contract required detection limit which is approx. 10 times the CRDL for water.

\*\* New York State background

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- \*\*\* Some forms of Cyanide are complex and very stable while other forms are pH dependent and hence are very unstable. Site-specific form(s) of Cyanide should be taken into consideration when establishing soil cleanup objective.
- \*\*\*\* Background levels for lead vary widely. Average levels in undeveloped, rural areas may range from 4-61 ppm. Average background levels in metropolitan or suburban areas or near highways are much higher and typically range from 200-500 ppm.
- \*\*\*\*\* Recommended soil cleanup objectives are average background concentrations as reported in a 1984 survey of reference material by E. Carol McGovern, NYSDEC.

REFERENCE NO. 13

### RECORD OF COMMUNICATION

TO: Tunru Yang

FROM: JANET TROTTER  
Region 2, ESAT/RSCC

DATE: Feb. 5, 2001

SUBJECT: QUALITY ASSURED DATA

=====

MESSAGE: PLEASE SIGN BELOW IN ACKNOWLEDGEMENT OF RECEIPT OF THE FOLLOWING AND RETURN ONE COPY OF THIS RECORD OF COMMUNICATION TO THE RSCC - REGION II.

① LACKAWANNA Foundry 28706 AATSHA Org-14 S/IW

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REPLY BY: \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

SIGNATURE: [Signature] DATE: 2/6/01

DATE RECEIVED BY RSCC: \_\_\_\_\_





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Functional Guidelines for Evaluating Organic Analysis

CASE No.: 28706  
LABORATORY: AATSLA

SDG No.: BNP05  
SITE: LACKAWANNA FOUNDRY

DATA ASSESSMENT

The current SOP HW-6 (Revision 11) June 1996, USEPA Region II Data Validation SOP for Statement of Work OLMO 4.2. for evaluating organic data have been applied.

All data are valid and acceptable except those analytes rejected "R" (unusable). Due to the detection of QC problems, some analytes may have the "J" (estimated), "N" (presumptive evidence for the presence of the material, "U" (non-detect) or "JN" (presumptive evidence for the presence of the material at an estimated value) flag. All action is detailed on the attached sheets.

The "R" flag means that the associated value is unusable. In other words, significant data bias is evident and the reported analyte concentration is unreliable.

Reviewer's *Andy Panayiotou*  
Signature: Andy Panayiotou Date: January 9, 2001

Verified By: *G Karas* Date: 01/10/2000 <sup>2001</sup>

SDG#BNP05

1. HOLDING TIME:

The amount of an analyte in a sample can change with time due to chemical instability, degradation, volatilization, etc. If the specified holding time is exceeded, the data may not be valid. Those analytes detected in the samples whose holding time has been exceeded will be qualified as estimated, "J". The non-detects (sample quantitation limits) will be flagged as estimated, "J", or unusable, "R", if the holding times are grossly exceeded. The following action was taken in the samples and analytes shown due to excessive holding time.

All samples met both Technical and Contractual Holding Times for all fractions.

2. SURROGATES

All samples are spiked with surrogate compounds prior to sample preparation to evaluate overall laboratory performance and efficiency of the analytical technique. If the measured surrogate concentrations were outside contract specifications, qualifications were applied to the samples and analytes as shown below.

VOA:

DC-35: The following volatile samples have system monitoring compound recoveries above the upper limit of the criteria window.  
Hits are qualified "J" and non-detects are not flagged.

BNP05MS

BNA/PEST:

No problems for this criterion.

3. MATRIX SPIKE/SPIKE DUPLICATE, MS/MSD:

The MS/MSD data are generated to determine the long term precision and accuracy of the analytical method in various matrices. The MS/MSD may be used in conjunction with other QC criteria for additional qualification of data.

No qualification is performed based on MS/MSD recovery.

4. BLANK CONTAMINATION:

Quality assurance (QA) blanks, i.e., method, trip, field, or rinse blanks are prepared to identify any contamination which may have been introduced into the samples during sample preparation or field activity. Method blanks measure laboratory contamination. Trip blanks measure cross-

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contamination of samples during shipment. Field and rinse blanks measure cross-contamination of samples during field operations. If the concentration of the analyte is less than 5 times the blank contaminant level (10 times for common contaminants), the analytes are qualified as non-detects, "U". The following analytes in the sample shown were qualified with "U" for these reasons:

A) Method blank contamination:

**DC-031:** The following volatile samples have analyte concentrations reported **above** the CRQL and less than or equal to ten times (10X) the associated method blank concentration.

**Hits are qualified "U" and non-detects are not flagged.**

BNP05, P05MSD, P12  
Acetone

**DC-200:** The following volatile samples have analyte concentrations reported **below** the CRQL and less than or equal to ten times (10X) the associated method blank concentration. Reported sample concentrations have been elevated to the CRQL.

**Hits are qualified "U" and non-detects are not flagged.**

BNP05, NP05MS, NP05MSD, NP06, NP07, NP08, NP09  
BNP10, NP11, NP12, NP13, BNQ85, Q88, Q89, Q91  
Methylene chloride

BNP07, NP08, NP09, NP11, BNQ87, Q88, Q89, Q91  
Acetone

**DC-236:** The following pesticide samples have analyte concentrations reported **below** the CRQL and less than or equal to five times (5X) the associated method blank concentration. Reported sample concentrations have been elevated to the CRQL.

**Hits are qualified "U" and non-detects are not flagged.**

BNP05  
gamma-BHC (Lindane), Endrin, Methoxychlor

BNP05MS  
Methoxychlor, Endrin aldehyde

BNP05MSD  
Methoxychlor, Endrin aldehyde

BNP06  
4,4'-DDT, Methoxychlor, Endrin aldehyde

BNP07  
Endrin, Methoxychlor, Endrin aldehyde

BNP08  
4,4'-DDT, Methoxychlor, Endrin ketone, Endrin aldehyde

BNP10  
Methoxychlor, Endrin aldehyde

BNP11  
Methoxychlor

BNP12  
4,4'-DDT

BNP13  
4,4'-DDT

BNQ85  
Endrin, 4,4'-DDT, Endrin ketone  
Endrin aldehyde

BNQ87  
Methoxychlor, Endrin aldehyde

BNQ88  
Methoxychlor, Endrin aldehyde

BNQ89  
Methoxychlor, Endrin ketone

BNQ91  
4,4'-DDT, Methoxychlor

**B) Field or rinse blank contamination:  
VOA:**

No contamination in the associated field blank.

**BNA/PEST:**

No qualification performed based on field blank contamination.

**C) Trip blank contamination for VOA aqueous samples:**

Not Applicable.

**D) Storage Blank associated with VOA samples only**

**DC-XXX:** The following volatile samples have analyte concentrations reported **above** the CRQL and less than or equal to ten times (10X) the associated **Storage** blank concentration.

Hits are qualified "U" and non-detects are not flagged.

BNP06, NP13, NP05MS  
Acetone

**E) Tics "R" rejected**

**VOA:**

The following TIC's are qualified "R". They have been identified as column bleed and Siloxane peaks.

BNP07, P08, P12, P13, BNQ85, Q88  
Unknown Siloxanes

**BNA:**

The following TIC's are qualified "R". They have been identified as Aldol condensation products.

BNP06, P07, Q87, Q88  
2-Pentanone, 4-Hydroxy-4-Methyl

The following TIC's are qualified "R". Their concentration is five times (5X) less than their associated method blank concentration.

BNP05, P06, P07, P08, P09, P10, P11, P12, P13, Q85  
BNQ87, Q88, Q89, Q91  
Unknown at various Retention times

The following TIC's are qualified "R". They have been identified as target compounds.

BNP09, P09DL  
Benzo(k)fluoranthene

**5. MASS SPECTROMETER TUNING:**

Tuning and performance criteria are established to ensure adequate mass resolution, proper identification of compounds and to some degree, sufficient instrument sensitivity. These criteria are not sample specific. Instrument performance is determined using standard materials. Therefore, these criteria should be met in all circumstances. The tuning standard for volatile organics is (BFB) Bromofluorobenzene and for semi-volatiles Decafluorotriphenyl-phosphine (DFTPP).

If the mass calibration is in error, all associated data will be classified as unusable "R".

No problems found for this qualification.

**6. CALIBRATION:**

Satisfactory instrument calibration is established to ensure that the instrument is capable of producing acceptable quantitative data. An initial calibration demonstrates that the instrument is capable of giving acceptable performance at the beginning of an experimental sequence. The continuing calibration checks document that the instrument is giving satisfactory daily performance.

A) Response Factor GC/MS:

The response factor measures the instrument's response to specific chemical compounds. The response factor for the Target Compound List (TCL) must be  $\geq 0.05$  in both initial and continuing calibrations. A value  $< 0.05$  indicates a serious detection and quantitation problem (poor sensitivity). Analytes detected in the sample will be qualified as estimated, "J". All non-detects for that compound will be rejected "R".

B) Percent Relative Standard Deviation (%RSD) and Percent Difference (%D):

Percent RSD is calculated from the initial calibration and is used to indicate the stability of the specific compound response factor over increasing concentration. Percent D compares the response factor of the continuing calibration check to the mean response factor (RRF) from the initial calibration. Percent D is a measure of the instrument's daily performance. Percent RSD must be  $< 30\%$  and %D must be  $< 25\%$ . A value outside of these limits indicates potential detection and quantitation errors. For these reasons, all positive results are flagged as estimated, "J" and non-detects are flagged "UJ". If %RSD and %D grossly exceed QC criteria, non-detects data may be qualified "R".

For the PEST/PCB fraction, if %RSD exceeds 20% for all analytes except for the two surrogates (which must not exceed 30% RSD), qualify all associated positive results "J" and non-detects "UJ".

The following analytes in the sample shown were qualified for %RSD and %D:

DC-100: The following semivolatile samples are associated with a continuing calibration percent difference (%D) outside primary criteria.

Hits are qualified "J" and non-detects are qualified "UJ".

BNP05, P05MS, P05MSD, P06, P07, P08, P09, SBLKFD  
Benzaldehyde

8. INTERNAL STANDARDS PERFORMANCE GC/MS:

Internal standards (IS) performance criteria ensure that the GC/MS sensitivity and response are stable during every experimental run. The internal standard area count must not vary by more than a factor of 2 (-50% to +100%) from the associated continuing calibration standard. The retention time of the internal standard must not vary more than  $\pm 30$  seconds from the associated continuing calibration standard. If the area count is outside the (-50% to +100%) range of the associated standard, all of the positive results for compounds quantitated using that IS are qualified as estimated, "J", and all non-detects as "UJ", or "R" if there is a severe loss of sensitivity.

If an internal standard retention time varies by more than 30 seconds, the reviewer will use professional judgement to determine either partial or total

rejection of the data for that sample fraction.

No problems with this criterion.

**9. COMPOUND IDENTIFICATION:**

**A) Volatile and Semi-Volatile Fractions:**

TCL compounds are identified on the GC/MS by using the analyte's relative retention time (RRT) and by comparison to the ion spectra obtained from known standards. For the results to be a positive hit, the sample peak must be within  $\pm 0.06$  RRT units of the standard compound and have an ion spectra which has a ratio of the primary and secondary m/e intensities within 20% of that in the standard compound. For the tentatively identified compounds (TIC) the ion spectra must match accurately. In the cases where there is not an adequate ion spectrum match, the laboratory may have provided false positive identifications.

**B) Pesticide Fraction:**

The retention times of reported compounds must fall within the calculated retention time windows for the two chromatographic columns and a GC/MS confirmation is required if the concentration exceeds 10ng/ml in the final sample extract.

**DC-422:** The following pesticide samples have Analytes for which the percent difference between column results exceeds primary criteria. Hits > CRQL are flagged "J". Or: if %D is > 50% and value is < CRQL, sample result is elevated to the CRQL and qualified "U".

**Analytes qualified "J".**

BNP05

gamma-BHC

BNP05, P12

gamma-Chlordane

BNP12

Endrin

BNP12, P05MSD

Endrin ketone

BNP13

Heptachlor, Endosulfan I, Methoxychlor, Endrin Aldehyde

BNQ85

Endosulfan II

BNP05MS

4,4'-DDD

**DC-423:** The following pesticide samples have analytes for which the percent difference between column results exceeds expanded criteria. Hits > CRQL are flagged "NJ"; or "R" when %D >100; or "NJ" when %D is between 100-200 (interference detected). Hits < CRQL are elevated to the CRQL and qualified "U".

**Analytes qualified "JN".**

BNP09  
Endrin ketone

BNP12  
Endrin Aldehyde

**Analytes qualified "R"**

BNP06  
beta-BHC

**Analytes qualified "U".**

BNP06, P08, P11, P12, P13, BNQ88  
alpha-BHC

BNP11, P13, BNQ85, Q87, Q88, Q91, Q92  
beta-BHC

BNP05, P12, Q87  
Heptachlor

BNP05, P05MS, P05MSD, P08, P09, P11, P13  
BNQ85, Q87, Q88  
Heptachlor epoxide

BNP08, P09, P10, P12, P13  
Dieldrin

BNP05MS, P05MSD, P08, P09, P12, P13  
4,4'-DDE

BNP08, P13  
Endrin

BNP08, P09, P12, Q88  
Endosulfan II

BNP06, P08, P09, P10, Q88, Q92  
4,4'-DDD

BNP05, P05MSD, P08, P09, Q85  
Endosulfan sulfate



BNP13  
Endrin ketone

BNP09  
Endrin Aldehyde

BNP08, P09, Q88  
alpha-Chlordane

BNP06, P07, P08, P09, P11  
gamma-Chlordane

**10. CONTRACT PROBLEMS NON-COMPLIANCE:**

**BNA:**

BNP08DL, P09DL - Analytes were diluted below the upper half of the calibration range. As per OLM04.2, sec. 10.6.5.3, p. D-43/SVOA, "The dilution factor chosen should keep the response of the largest peak for a target compound in the upper half of the calibration range of the Instrument."

**11. FIELD DOCUMENTATION:**

As per the Traffic reports, none of the field blanks collected required a volatile analysis.

**12. OTHER PROBLEMS**

**13. This package contains reextractions, reanalyses or dilutions. Upon reviewing the QA results, the following Form 1(s) are identified not to be used.**

**BNA:**

BNP08DL, P09DL

DPO:  ACTION  FYI

REGION 2

ORGANIC REGIONAL DATA ASSESSMENT SUMMARY

CASE NO. 28706 LABORATORY AATSLA  
SDG NO. B04AN DATA USER EPA/Region II  
SOW OLM04.2 REVIEW COMPLETION DATE January 9, 2001  
NO. OF SAMPLES 1 WATER 14 SOIL \_\_\_\_\_ OTHER \_\_\_\_\_  
REVIEWER:  ESD  ESAT  OTHER, CONTRACTOR \_\_\_\_\_

| QC ITEM                 | VOA | BNA | PEST |  |  |
|-------------------------|-----|-----|------|--|--|
| HOLDING TIMES           | O   | O   | O    |  |  |
| GC-MS PERFORMANCE       | O   | O   | F    |  |  |
| INITIAL CALIBRATIONS    | O   | O   | O    |  |  |
| CONTINUING CALIBRATIONS | O   | X   | O    |  |  |
| FIELD BLANKS (F = N/A)  | O   | O   | O    |  |  |
| LABORATORY BLANKS       | O   | O   | O    |  |  |
| SURROGATES              | X   | O   | O    |  |  |
| MATRIX SPIKE/DUPLICATES | O   | O   | O    |  |  |
| % MOISTURE              | O   | O   | O    |  |  |
| INTERNAL STANDARDS      | O   | O   | F    |  |  |
| COMPOUND IDENTIFICATION | O   | O   | X    |  |  |
| COMPOUND QUANTITATION   | O   | O   | O    |  |  |
| SYSTEM PERFORMANCE      | O   | O   | O    |  |  |
| OVERALL ASSESSMENT      | O   | O   | O    |  |  |

O = No problems or minor problems that do not affect data usability.

X = No more than about 5% of the data points are qualified as either estimated or unusable.

M = More than about 5% of the data points are qualified as either estimated or unusable.

Z = More than about 5% of the data points are qualified as unusable.

DPO ACTION ITEMS:

AREAS OF CONCERN:

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DATA REJECTION SUMMARY

Type of Review: Organic Date : January 9, 2001 Case : 28706, SDG# BNP05  
 Site Name: Zschiegner Refining Lab Name: AATSLA Reviewer's Initial : AP  
 Number of Samples: 1 Waters, 14 Soils, + QC +Dilutions/Reanalyses

Analytes Rejected Due To Exceeding Review Criteria For: Number of Compounds /Number of Fractions (Samples)

|           | Surrogates | Holding Times | Calibration | Contamination | ID | Internal Standard | Other* | Total # Samples | Total # REJECTED/<br>Total Analytes in samples<br>Percent |
|-----------|------------|---------------|-------------|---------------|----|-------------------|--------|-----------------|---|
| VOA (48)  | 0          | 0             | 0           | 0             | 0  | 0                 | 0      | 22              | 0 1056 0  |
| AE (14)   | 0          | 0             | 0           | 0             | 0  | 0                 | 0      | 21              | 0 294 0   |
| BN (51)   | 0          | 0             | 0           | 0             | 0  | 0                 | 0      | 21              | 0 1071 0  |
| PEST (21) | 0          | 0             | 0           | 0             | 1  | 0                 | 0      | 19              | 1 399 0   |
| PCB (7)   | 0          | 0             | 0           | 0             | 0  | 0                 | 0      | 19              | 0 133 0   |

Analytes Estimated Due To Exceeding Review Criteria For: Number of Compounds /Number of Fractions (Samples)

|           | Surrogates | Holding Times | Calibration | Contamination | ID | Internal Standard | Other* | Total # Samples | Total # ESTIMATED/<br>Total Analytes in samples<br>Percent |
|-----------|------------|---------------|-------------|---------------|----|-------------------|--------|-----------------|--|
| VOA (48)  | 9          | 0             | 0           | 0             | 0  | 0                 | 0      | 22              | 9 1056 1   |
| AE (14)   | 0          | 0             | 0           | 0             | 0  | 0                 | 0      | 21              | 0 294 0  |
| BN (51)   | 0          | 0             | 8           | 0             | 0  | 0                 | 0      | 21              | 8 1071 1   |
| PEST (21) | 0          | 0             | 0           | 0             | 14 | 0                 | 0      | 19              | 14 399 4   |
| PCB (7)   | 0          | 0             | 0           | 0             | 0  | 0                 | 0      | 19              | 0 133 0  |

\* % MOISTURE

STANDARD OPERATING PROCEDURE

Region II 4  
 Mod: CLP/SOW OLMOZ.2

Date: June 1996  
 SOP HW-6, Rev. 11

YES NO N/A

PACKAGE COMPLETENESS AND DELIVERABLES

CASE NUMBER: 28706 LABORATORY: AATSLA  
 SITE NAME: Lackawanna Foundry SDG Number(s): BNP05

1.0 Chain of Custody and Sampling Trip Reports

1.1 Are the Traffic Reports/Chain-of-Custody Records present for all samples? [✓]          

ACTION: If no, contact RSCC, or contact the WAM to obtain replacement of missing or illegible copies from the lab.

1.2 Is the Sampling Trip Report present for all samples and all fractions? [ ]          

ACTION: If no, contact either RSCC or ask the WAM to obtain this information from the prime contractor.

2.0 Data Completeness and Deliverables

2.1 Have any missing deliverables been received and added to the data package? [✓]          

NOTE: The lab is required to submit data for only two analyses, for each fraction. (i.e., the original sample and one dilution, or the most concentrated dilution analyzed and one further dilution.)

ACTION: Contact the WAM to obtain an explanation or resubmittal of any missing deliverables from the lab. If lab cannot provide them, note the effect on the review of the package in the Contract Problems/Non-compliance section of the Data Assessment and the Organic Regional Data Assessment Summary form.

2.2 Was CLASS CCS checklist included with package? [ ]          

2.3 Are there any discrepancies between the Traffic Reports/Chain-of-Custody Records, Sampling Report and Sample Tags?      [✓]



STANDARD OPERATING PROCEDURE

IS EPA Region II  
 Method: CLP/SOW OLMO3.2

Date: June 1996  
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YES NO N/A

ACTION: If yes, contact the WAM to obtain an explanation or resubmittal of any missing deliverables from the laboratory.

3.0 Cover Letter SDG Narrative

- |   |  |   |     |     |
|---|--|---|-----|-----|
| 3.1   | Is the Narrative or Cover Letter Present?  | [ <input checked="" type="checkbox"/> ] | ___ | ___ |
| 3.2   | Are case number, SDG number and contract number contained in the SDG Narrative or cover letter (see SOW, Exhibit B, section 2.6.1)?  | [ <input checked="" type="checkbox"/> ] | ___ | ___ |
| 3.3   | Does the narrative contain the following information:  |   |     |     |
|   | VOA: description of trap and columns used during sample analyses?  | [ <input checked="" type="checkbox"/> ] | ___ | ___ |
|   | BNA: description of columns used during sample analyses?   | [ <input checked="" type="checkbox"/> ] | ___ | ___ |
|   | Pest: description of columns used during sample analyses?  | [ <input checked="" type="checkbox"/> ] | ___ | ___ |
| NOTE: As per section 6.23.3.1 SOW/p. D-11/Pest, Packed columns are not permitted. |  |   |     |     |
| 3.4   | Does the narrative, VOA and BNA sections, contain a list of all TICs identified as alkanes and their estimated concentrations?   | [ <input checked="" type="checkbox"/> ] | ___ | ___ |
| 3.5   | Does the narrative contain a record of all cooler temperatures? If the temperature of a cooler was exceeded, > 10° C, the lab must list by fraction and sample number, all affected samples. | [ <input checked="" type="checkbox"/> ] | ___ | ___ |
| 3.6   | Does the narrative contain a list of the pH values determined for each water sample submitted for volatile analysis (SOW Exhibit B, section 2.6.1.2)?  | [ <input checked="" type="checkbox"/> ] | ___ | ___ |
| 3.7   | Does the Case Narrative contain the statement, "verbatim", as required in Section B of the SOW?  | [ <input checked="" type="checkbox"/> ] | ___ | ___ |

ACTION: If "No", to any question in this section, contact the WAM to obtain all necessary resubmittals. If information is not available, document in the Data Assessment under Contract Problems/Non-Compliance section.

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Date: June 1996  
SOP HW-6, Rev. 11

YES NO N/A

4.0 Data Validation Checklist

4.1 Check the package for the following discrepancies:

- |        |   |                                     |                                     |     |
|--------|---|-------------------------------------|-------------------------------------|-----|
| a.     | Is the package paginated in ascending order starting from the SDG narrative?              | <input checked="" type="checkbox"/> | ___                                 | ___ |
| b.     | Are all forms and copies legible?   | <input checked="" type="checkbox"/> | ___                                 | ___ |
| * 1 c. | Is each fraction assembled in the order set forth in the SOW?                             | <input checked="" type="checkbox"/> | ___                                 | ___ |
| d.     | Is a Sample Data Summary Package submitted immediately preceding the Sample Data Package? | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | ___ |

The following checklist is divided into three parts. Part A is for any VOA analyses, Part B is for BNAs and Part C is Pesticide/PCBs.

Does this package contain:

- |                     |                                     |     |
|---------------------|-------------------------------------|-----|
| VOA Data?           | <input checked="" type="checkbox"/> | ___ |
| BNA Data?           | <input checked="" type="checkbox"/> | ___ |
| Pesticide/PCB data? | <input checked="" type="checkbox"/> | ___ |

ACTION: Complete corresponding parts of checklist.

\* c1. Most of package is in the proper order except for the chromatograms. They are placed after the Quant reports while ~~st~~ they should be placed in front instead.

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US EPA Region II  
Method: CLP/SOW, OLMO3.2

Date: June 1996  
SOP HW-6, Rev. 11

YES NO N/A

PART A: VOA ANALYSES

..0 Sample Conditions/Problems

1.1 Do the Traffic Reports/Chain-of-Custody Records, Sampling Report or Lab Narrative indicate any problems with sample receipt, condition of samples, analytical problems or special circumstances affecting the quality of the data?     rV    

ACTION: If any sample analyzed as a soil, other than TCLP, contains 50% - 90% water, all data should be flagged as estimated (J). If a soil sample other than TCLP contains more than 90% water, all data should be qualified as unusable (R).

ACTION: If samples were not iced or the ice was melted upon arrival at the laboratory and the cooler temperature was elevated (> 10° C), then flag all positive results with a "J" and all non-detects "UJ".

ACTION: If both VOA vials for a sample have air bubbles or the VOA vial analyzed had air bubbles, flag all positive results "J" and all non-detects "R".

ACTION: The smallest soil size permitted is 0.5g. If any soil sample is smaller than 0.5g, document in the Data Assessment under Contract Problems/Non-Compliance.

.0 Holding Times

2.1 Have any VOA technical holding times, determined from date of collection to date of analysis, been exceeded?     rV    

Technical Holding Times: If unpreserved, aqueous samples, maintained at 4° C for aromatic hydrocarbons analysis must be analyzed within 7 days of collection. If preserved with HCl (pH < 2) and stored at 4° C, then aqueous samples must be analyzed within 14 days of collection. If uncertain about preservation, contact sampler to determine whether or not samples were preserved. The holding time for soils is 10 days from date of collection.



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YES NO N/A

Table of Holding Time Violations  
(See Chain-of-Custody Records)

| Sample ID | Sample Matrix | Was Sample Preserved? | Date Sampled | Date Lab Received | Date Analyzed |
|-----------|---------------|-----------------------|--------------|-------------------|---------------|
| _____     | _____         | _____                 | _____        | _____             | _____         |
| _____     | _____         | _____                 | _____        | _____             | _____         |
| _____     | _____         | _____                 | _____        | _____             | _____         |
| _____     | _____         | _____                 | _____        | _____             | _____         |
| _____     | _____         | _____                 | _____        | _____             | _____         |
| _____     | _____         | _____                 | _____        | _____             | _____         |
| _____     | _____         | _____                 | _____        | _____             | _____         |

ACTION: If technical holding times are exceeded, flag all positive results as estimated "J" and sample quantitation limits as estimated "UJ", and document in the Data Assessment that holding times were exceeded. If analyses were done more than 14 days beyond holding time, either on the first analysis or upon re-analysis, the reviewer must use professional judgement to determine the reliability of the data and the effects of additional storage on the sample results. At a minimum, all results must be qualified "J", but the reviewer may determine that non-detect data are unusable "R". If holding times are exceeded by more than 28 days, all non detect data are unusable "R".

NOTE: Contractual Holding Times: Analysis of water and soil/sediment samples must be completed within 10 days of Validated Time of Sample Receipt (VTSR). This requirement does not apply to Performance Evaluation (PE) samples.

ACTION: If contractual holding times are exceeded, document in the Data Assessment and on the Organic Regional Data Assessment Summary form.

NOTE: The data reviewer must note in the Data Assessment whether or not technical and contractual holding times were met.

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US EPA Region II  
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YES NO N/A

3.0 System Monitoring Compound (SMC) Recovery (Form II)

3.1 Are the VOA SMC Recovery Summaries (Form II) present for each of the following matrices:

a. Low Water?

b. Low Soil?

c. Med Soil?

3.2 Are all the VOA samples listed on the appropriate System Monitoring Compound Recovery Summary for each of the following matrices:

a. Low Water?

b. Low Soil?

c. Med Soil?

ACTION: Contact the WAM to obtain an explanation or resubmittal of any missing deliverables from the laboratory. If missing deliverables are unavailable, document the effect in the Data Assessment.

3.3 Were outliers marked correctly with an asterisk?

ACTION: Circle all outliers with red pencil.

3.4 Was one or more VOA system monitoring compound recovery outside of contract specifications for any sample or method blank?

If yes, were samples re-analyzed? *MS not required*

Were method blanks re-analyzed?

ACTION: If recoveries are  $\geq 10\%$ , but 1 or more compounds fail to meet SOW specifications:

1. All positive results are qualified as estimated "J".

2. Flag all non-detects as estimated detection limits "UJ" where recovery is less than the lower acceptance limit.

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US EPA Region II  
Method: CLP/SOW OLMO3.2

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YES NO N/A

3. If SMC recoveries are above allowable levels, do not qualify non-detects.

If any system monitoring compound recovery is < 10%:

- 1. Flag all positive results as estimated "J".
- 2. Flag all non-detects as unusable "R".

Professional judgement should be used to qualify data that only have method blank SMC recoveries out of specification in both original and re-analyses. Check the internal standard areas.

NOTE: Contractual requirements state that if any SMC fails the acceptance criteria, the sample must be re-analyzed. If the affected sample was not re-analyzed, document in the Data Assessment under Contract Problems/Non-Compliance and in the Organic Regional Data Assessment Summary.

NOTE: The laboratory must submit the following data:

1. If SMC recoveries and internal standard responses meet the acceptance criteria in the re-analyzed sample, then the laboratory must submit only the re-analysis.

2. If an SMC recovery and/or internal standard response fails to meet the acceptance criteria upon re-analysis, then submit data from both analyses.

(Refer to section 11.4.3.2, page D-46/VOA of the SOW for more information.)

3.5 Are there any transcription/calculation errors between raw data and Form II?

ACTION: If large errors exist, contact the WAM to obtain an explanation or resubmittal of corrected deliverables from the laboratory. Make any necessary corrections and note the effect in the Data Assessment.

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YES NO N/A

4.0 Matrix Spikes (Form III)

4.1 Is the Matrix Spike/Matrix Spike Duplicate Recovery Form (Form III) present?  YES  NO  N/A

4.2 Were matrix spikes analyzed at the required frequency for each of the following matrices:

a. Low Water?  YES  NO  N/A

b. Low Soil?  YES  NO  N/A

c. Med Soil?  YES  NO  N/A

ACTION: If any matrix spike data are missing, take the action specified in section 3.2 above.

4.3 How many VOA spike recoveries are outside QC limits?

|                      |                    |
|----------------------|--------------------|
| <u>Water</u>         | <u>Soils</u>       |
| <u>N/A</u> out of 10 | <u>4</u> out of 10 |

4.4 How many RPDs for matrix spike and matrix spike duplicate recoveries are outside QC limits?

|                     |                   |
|---------------------|-------------------|
| <u>Water</u>        | <u>Soils</u>      |
| <u>N/A</u> out of 5 | <u>0</u> out of 5 |

ACTION: No action is taken based upon MS/MSD data alone. However, using informed professional judgement, the MS/MSD results may be used in conjunction with other QC criteria to determine the need for qualification of the data.

ACTION: Circle all outliers with red pencil.

5.0 Blanks (Form IV)

5.1 Is the Method Blank Summary (Form IV) present?  YES  NO  N/A

5.2 Frequency of Analysis: for the analysis of VOA TCL compounds, has a reagent/method blank been analyzed for each SDG or every 20 samples of similar matrix (low water, low soil or medium soil), whichever is more frequent?  YES  NO  N/A

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|   | YES                                 | NO                       | N/A                                 |
|---|-------------------------------------|--------------------------|-------------------------------------|
| 5.3 Has a VOA method blank been analyzed at least once every twelve hours for each concentration level and GC/MS system used?                     | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/>            |
| 5.4 Was a VOA instrument blank analyzed after each sample/dilution which contained a target compound that exceeded the initial calibration range? | <input type="checkbox"/>            | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
| 5.5 Was a VOA storage blank analyzed at the end of all samples for each SDG in a case?  | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/>            |

**ACTION:** If any method/instrument blank data are missing, contact the WAM to obtain any missing deliverables from the laboratory. If method blank data are not available, reject "R" all associated positive data. However, using professional judgement, the data reviewer may substitute field blank or trip blank data for missing method blank data.

If any instrument blank analyzed after a sample with high concentration is missing, contact the WAM to obtain any missing deliverables from the laboratory. If the instrument blank was not analyzed or not available, inspect the chromatogram of the sample analyzed immediately after this analysis for possible carryover. Use professional judgement to determine if any contamination occurred and qualify analyte(s) accordingly.

If storage blank data is missing, contact the WAM to obtain any missing deliverables from the laboratory. If unavailable, note in the Contract Problems/Non-Compliance section of the Data Assessment.

5.6 The validator should verify that the correct identification scheme for the EPA Blank samples were used. See page B-33, section 3.3.7.3 of the SOW for further information.

Was the correct identification scheme used for all VOA blanks?

**ACTION:** Contact the WAM to obtain missing deliverables from the lab, or make the required corrections on the forms. Document in the Data Assessment under Contract Problems/Non-compliance if corrections were made by the validator.

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YES NO N/A

5.7 Chromatography: review the blank raw data-chromatograms (RICs), quant. reports or data system printouts and spectra. Is the chromatographic performance (baseline stability) for each instrument acceptable for VOAs?

ACTION: Use professional judgement to determine the effect on the data.

5.8 Are all detected hits for target compounds in method, instrument and storage blanks less than the CRQL for that analyte?

Exception: Acetone and 2-butanone must be less than 5 times the CRQL, and methylene chloride must be less than 2.5 times its CRQL.

ACTION: If no, an explanation and laboratory's corrective actions must be addressed in the case narrative. If the narrative contains no explanation, then make a note in the Contract Problems/Non-Compliance section of the Data Assessment.

6.0 Contamination

NOTE: "Water blanks", "drill blanks", and "distilled water blanks" are validated like any other sample, and are not used to qualify data. Do not confuse them with the other QC blanks discussed below.

6.1 Do any method/instrument/reagent/storage blanks have positive results (TCL and/or TIC) for VOAs?

NOTE: When applied as directed in the table below, the contaminant concentration in these blanks are multiplied by the sample dilution factor and corrected for %moisture when necessary.

NOTE: A contaminated instrument blank is not allowable under this SOW. See page D-48/VOA, section 12.1.2.4 for additional information. Document in the Data Assessment under Contract Problems/Non-Compliance if contaminated instrument blank was submitted.

6.2 Do any field/trip/rinse blanks have positive VOA results (TCL and/or TIC)?

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YES NO N/A

ACTION: Prepare a list of the samples associated with each of the contaminated blanks. (Attach a separate sheet.)

NOTE: All field blank results associated with a particular group of samples (may exceed one per case) must be used to qualify data. Trip blanks are used to qualify only those samples with which they were shipped and are not required for non-aqueous matrices. Blanks may not be qualified because of contamination in another blank. Field Blanks & Trip Blanks must be qualified for system monitoring compound, instrument performance criteria, spectral or calibration QC problems.

ACTION: Follow the directions in the table below to qualify TCL results due to contamination. Use the largest value from all the associated blanks. If any blanks are grossly contaminated, all associated data should be qualified as unusable "R".

| For:               | Flag sample result with a "U" when:            | Report CRQL & qualify "U" when:               | No qualification is needed when:              |
|--------------------|--|---|---|
| Methylene Chloride | Sample conc. is > CRQL, but ≤ 10× blank value. | Sample conc. is < CRQL and ≤ 10× blank value. | Sample conc. is > CRQL and > 10× blank value. |
| Acetone            |  |   |   |
| Toluene            |  |   |   |
| 2-Butanone         |  |   |   |
| Other Contaminants | Sample conc. is > CRQL, but ≤ 5× blank value.  | Sample conc. is < CRQL and ≤ 5× blank value.  | Sample conc. is > CRQL and > 5× blank value.  |

NOTE: Analytes qualified "U" for blank contamination are still considered as "hits" when qualifying for calibration criteria.

ACTION: For TIC compounds, if the concentration in the sample is less than five times the concentration in the most contaminated associated blank, flag the sample data "R".

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YES NO N/A

6.3 Are there field/rinse/equipment blanks associated with every sample?

ACTION: For low level samples, note in the Data Assessment that there is no associated field/rinse/equipment blank. For samples with high concentrations of suspected blank contaminants, use professional judgement to qualify these values and make a note in the Data Assessment.

Exception: samples taken from a drinking water tap do not have associated field blanks.

7.0 GC/MS Instrument Performance Check (Form V)

7.1 Are the GC/MS Instrument Performance Check Forms (Form V) present for Bromofluorobenzene (BFB)?

7.2 Are the enhanced bar graph spectrum and mass/charge (m/z) listing for the BFB provided for each twelve hour shift?

7.3 Has an instrument performance check been analyzed for every analytical sequence on each instrument?

ACTION: List date, time, instrument ID, and sample numbers for which associated GC/MS tuning data are unavailable.

| DATE  | TIME  | INSTRUMENT | SAMPLE NUMBERS |
|-------|-------|------------|----------------|
| _____ | _____ | _____      | _____          |
| _____ | _____ | _____      | _____          |
| _____ | _____ | _____      | _____          |

ACTION: Notify the WAM to obtain missing data, if possible. If the lab cannot provide the missing data, reject, "R", all data generated outside an acceptable twelve hour calibration interval.

7.4 Have the ion abundances been normalized to m/z 95 as specified in Exhibit D, page D-56/VOA?

NOTE: All ion abundance ratios must be normalized to m/z 96, the nominal base peak, even though the



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YES NO N/A

ion abundance of m/z 174 may be up to 120% that of m/z 95.

ACTION: If mass assignment is in error, qualify all associated data as unusable "R".

7.5 Have the ion abundance criteria been met for each instrument used?  YES  NO  N/A

ACTION: List all data which do not meet ion abundance criteria (attach a separate sheet).

ACTION: If ion abundance criteria are not met, the Region II TPO must be notified.

7.6 Are there any transcription/calculation errors between mass lists and Form Vs? (Check at least two values, but if errors are found check more.)  YES  NO  N/A

7.7 Is the number of significant figures for the reported relative abundances consistent with the number given for each ion in the ion abundance criteria column?  YES  NO  N/A

ACTION: If large errors exist, take action as specified in section 3.5 above.

7.8 Are the spectra of the mass calibration compound acceptable?  YES  NO  N/A

ACTION: Use professional judgement to determine whether associated data should be accepted, qualified, or rejected.

8.0 Target Compound List (TCL) Analytes (FORM I VOA)

8.1 Are the Organic Analysis Data Sheets (Form I VOA) present with required header information on each page, for each of the following:

a. Samples and/or fractions as appropriate?  YES  NO  N/A

b. Matrix spikes and matrix spike duplicates?  YES  NO  N/A

c. Blanks?  YES  NO  N/A

8.2 Are the VOA Reconstructed Ion Chromatograms, the mass spectra for the identified compounds, and the data system printouts (quant. reports) included in the sample package for each of the

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YES NO N/A

following:

- |  |                                     |     |     |
|--|-------------------------------------|-----|-----|
| a. Samples and/or fractions as appropriate?                                  | <input checked="" type="checkbox"/> | ___ | ___ |
| b. Matrix spikes and matrix spike duplicates<br>(mass spectra not required)? | <input checked="" type="checkbox"/> | ___ | ___ |
| c. Blanks?   | <input checked="" type="checkbox"/> | ___ | ___ |

ACTION: If any data are missing, take action specified in 3.2 above.

- |  |                                     |                                     |                                     |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| 8.3 Are the response factors shown in the quant. report?       | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | ___                                 |
| 8.4 Is chromatographic performance acceptable with respect to: |                                     |                                     |                                     |
| a. Baseline stability?   | <input checked="" type="checkbox"/> | ___                                 | ___                                 |
| b. Resolution?   | <input checked="" type="checkbox"/> | ___                                 | ___                                 |
| c. Peak shape?   | <input checked="" type="checkbox"/> | ___                                 | ___                                 |
| d. Full-scale graph (attenuation)?                             | <input checked="" type="checkbox"/> | ___                                 | ___                                 |
| e. Other: _____?   | <input type="checkbox"/>            | ___                                 | <input checked="" type="checkbox"/> |

ACTION: Use professional judgement to determine the acceptability of the data.

- |  |                                     |     |     |
|--|-------------------------------------|-----|-----|
| 8.5 Are the lab-generated standard mass spectra of the identified VOA compounds present for each sample? | <input checked="" type="checkbox"/> | ___ | ___ |
|--|-------------------------------------|-----|-----|

ACTION: If any mass spectra are missing, take action as specified in 3.2 above. If the lab does not generate its own standard spectra, document in the Contract Problems/Non-compliance section of the Data Assessment and the Organic Regional Data Assessment Summary.

- |   |                                     |     |     |
|---|-------------------------------------|-----|-----|
| 8.6 Is the RRT of each reported compound within 0.06 RRT units of the standard RRT in the continuing calibration?                         | <input checked="" type="checkbox"/> | ___ | ___ |
| 8.7 Are all ions present in the standard mass spectrum at a relative intensity greater than 10% also present in the sample mass spectrum? | <input checked="" type="checkbox"/> | ___ | ___ |

0.27

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YES NO N/A

8.8 Do sample and standard relative ion intensities agree within  $\pm 20\%$ ?         

ACTION: Use professional judgement to determine acceptability of data. If it is determined that incorrect identifications were made, all such data should be rejected "R", flagged "N" (presumptive evidence of the presence of the compound) or changed to not detected "U" at the calculated detection limit. In order to be positively identified, the data must comply with the criteria listed in 8.6, 8.7, and 8.8.

ACTION: When sample carry-over is suspected, use professional judgement determine if instrument cross-contamination has affected positive compound identifications.

9.0 Tentatively Identified Compounds (TIC)

9.1 Are all Tentatively Identified Compound Forms (Form I Part B) present; and do listed TICs include scan number or retention time, estimated concentration and "JN" qualifier?         

9.2 Are the mass spectra for the tentatively identified compounds and associated "best match" spectra included in the sample package for each of the following:

a. Samples and/or fractions as appropriate?         

b. Blanks?         

c. Alkanes listed for each sample?         

ACTION: If any TIC data are missing, take action specified in 3.2 above.

ACTION: Add "JN" qualifier to all chemically named TICs, if missing.

9.3 Are any TCL compounds (from any fraction) listed as TIC compounds? (Example: 1,2-dimethylbenzene is xylene, a VOA TCL analyte, and should not be reported as a TIC.)         

ACTION: Flag with "R" any TCL compound listed as a TIC.

9.4 Are all ions present in the reference mass

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YES NO N/A

spectrum with a relative intensity greater than 10% also present in the sample mass spectrum?

9.5 Do TIC and "best match" standard relative ion intensities agree within ±20%?

ACTION: Use professional judgement to determine the acceptability of TIC identifications. If it is determined an incorrect identification was made, change the identification to "unknown," or to some less specific identification as appropriate. (Example: "C3 substituted benzene.")

Also, when a compound is not found in any blank, but is detected in a sample and is a suspected artifact of a common laboratory contaminant, the result should be qualified as unusable "R". (E.g., Common Lab Contaminants: CO<sub>2</sub> (M/E 44), siloxanes (M/E 73) hexane, aldol condensation products, solvent preservatives, and related by-products - see the National Functional Guidelines for further guidance.)

9.6 Are TICs with responses < 10% of the internal standard (as determined by inspection of the peak areas or height) reported?

ACTION: If yes, cross out questionable TIC(s).

10.0 Compound Quantitation and Reported Detection Limits

10.1 Are there any transcription/calculation errors in Form I results? (Check at least two positive values. Verify that the correct internal standards, quantitation ions, and RRF were used to calculate Form I results.)

10.2 Are the CRQLs adjusted to reflect sample dilutions and, for soils, sample moisture?

ACTION: If errors are large, take action as specified in section 3.5 above.

ACTION: When a sample is analyzed at more than one dilution, the lowest CRQLs are used (unless a QC exceedance dictates the use of the higher CRQL data from the diluted sample). Replace concentrations that exceeded the calibration range in the original analysis by crossing out

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YES NO N/A

the "E" and its corresponding value on the original Form I and substituting the data from the diluted sample. Specify which Form I is to be used, then draw a red "X" across the entire page of all Form Is not to be used, including any in the data summary package.

11.0 Standards Data (GC/MS)

11.1 Are the Reconstructed Ion Chromatograms, and data system printouts (quant. reports) present for each initial and continuing calibration? [✓] \_\_\_ \_\_\_

ACTION: If any calibration standard data are missing, take action specified in 3.2 above.

12.0 GC/MS Initial Calibration (Form VI)

12.1 Are the Initial Calibration Forms (Form VI) present and complete at concentrations of 10, 20, 50, 100, 200ng for separate calibrations of low water/med soils (unheated purge) and low soils (heated purge)? [✓] \_\_\_ \_\_\_

ACTION: If any calibration standard forms are missing, take action specified in 3.2 above.

12.2 Were all low level soil standards, blanks and samples analyzed by heated purge? [✓] \_\_\_ \_\_\_

ACTION: If low level soil samples were not heated during purge, qualify positive hits "J" (estimated) and non-detects "R".

12.3 Are the % relative standard deviation (%RSD) values for VOAs  $\leq$  30% over the concentration range of the calibration? [✓] \_\_\_ \_\_\_

NOTE: Although 11 VOA compounds have a contractual minimum RRF and no maximum %RSD, the technical acceptance criteria are the same for all analytes.

ACTION: Circle all outliers with red pencil.

ACTION: If %RSD is  $>$  30.0%, qualify associated positive results for that analyte "J" (estimated) and non-detects using professional judgement. When %RSD is  $>$  90%, flag all non-detects for that analyte "R" (unusable) and positive hits "J".

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YES NO N/A

NOTE: Analytes previously qualified "U" for blank contamination are still considered as "hits" when qualifying for initial calibration criteria.

12.4 Are any average RRFs < 0.05?         

ACTION: Circle all outliers with red pencil.

ACTION: If the average RRF is < 0.05, then qualify associated non-detects with an "R" and flag associated positive data as estimated "J".

NOTE: Contract Requirement: The SOW allows up to two of the required analytes to fail contractual %RSD or RRF criteria, provided the %RSD is ≤ 40% and RRF is ≥ 0.010. (See Table 5, page D-59/VOA and analytes marked with a "\*" on Form VI for required analytes and contractual criteria.) Technical criteria, however, are the same for all analytes.

ACTION: If more than two analytes failed %RSD or RRF criteria, document in the Data Assessment under Contract Problems/Non-Compliance and the Organic Regional Data Assessment Summary.

12.5 Are there any transcription/calculation errors in the reporting of average relative response factors (RRF) or %RSD? (Check at least 2 values, but if errors are found, check more.)         

ACTION: Circle errors with red pencil.

ACTION: If errors are large, contact the WAM to obtain an explanation/resubmittal from the lab, document in the Data Assessment under Contract Problems/Non-Compliance and in the Organic Regional Data Assessment Summary.

13.0 GC/MS Continuing Calibration (Form VII)

13.1 Are the Continuing Calibration Forms (Form VII) present and complete for separate calibration of low water/med soil and low soil samples?         

13.2 Has a continuing calibration standard been analyzed for every twelve hours of sample analysis per instrument?         

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YES NO N/A

ACTION: If any forms are missing or no continuing calibration standard has been analyzed within twelve hours of every sample analysis, contact the WAM to request an explanation/resubmittal from the lab. If continuing calibration data are not available, flag all associated sample data as unusable "R".

ACTION: List below all sample(s) that were not analyzed within twelve hours of the previous continuing calibration.

\_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

13.3 Do any volatile compounds have a percent difference (%D) between the initial and continuing RRF which exceeds the ±25% criteria?               

NOTE: Although 11 VOA compounds have a contractual minimum RRF and no maximum %D, the technical acceptance criteria are the same for all analytes.

ACTION: Circle all outliers with red pencil.

ACTION: Qualify both positive results and non-detects for the outlier compound(s) as estimated. When %D is > 90%, qualify all non-detects for that analyte unusable (R) and positive results estimated (J).

13.4 Are any continuing calibration RRFs < 0.05?               

ACTION: Circle all outliers with red pencil.

ACTION: If the RRF is < 0.05, qualify the associated non-detects as unusable "R" and the associated positive values "J".

NOTE: Contract Requirement: The SOW allows up to two of the required analytes to fail contractual %D and RRF criteria, provided that the %D is ≤ 40% and the RRF is ≥ 0.010. (See Table 5 pg. D-59/VOA or analytes marked with a "\*" on Form VI for required analytes.) Technical criteria, however, are the same for all analytes.

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YES NO N/A

ACTION: If more than two analytes failed %D and RRF, criteria document in the Data Assessment under contract Problems/Non-Compliance and on the Organic Regional Data Assessment Summary form.

13.5 Are there any transcription/calculation errors in the reporting of RRF or %D between initial and continuing RRFs? (Check at least two values, but if errors are found, check more.)

\_\_\_\_\_ RV \_\_\_\_\_

ACTION: Circle errors with red pencil.

ACTION: If errors are large, contact the WAM to obtain an explanation/resubmittal from the lab, document in the Data Assessment under Contract Problems/Non-Compliance.

14.0 Internal Standard (Form VIII)

14.1 Are the internal standard areas (Form VIII) of every sample and blank within the upper and lower limits (-50% to +100%) for each continuing calibration?

RV \_\_\_\_\_

If no, was the sample re-analyzed?

1 \_\_\_\_\_

ACTION: 1. Circle all outliers with red pencil.

2. List all the outliers below.

| Sample # | Internal Std. | Area  | Lower/Upper Limit |
|----------|---------------|-------|-------------------|
| _____    | _____         | _____ | _____ / _____     |
| _____    | _____         | _____ | _____ / _____     |
| _____    | _____         | _____ | _____ / _____     |
| _____    | _____         | _____ | _____ / _____     |

(Attach additional sheets if necessary, or attach copies of Form VIIIs.)

ACTION: If any sample was not re-analyzed, document in the Data Assessment under Contract Problems/Non-Compliance.

ACTION: 1. If the internal standard area count is outside the upper or lower limit, flag with "J" all positive results quantitated with this

0.33



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YES NO N/A

internal standard.

2. Do not qualify non-detects when associated IS area counts are > 100%.

3. If the IS area in the sample is below the "lower limit," < 50%, qualify all analytes associated with that IS estimated, "J". If the area counts are extremely low, < 25% of the area in the 12 hour standard, or if performance exhibits a major abrupt drop-off, flag all associated non-detects as unusable, "R", and positive hits estimated, "J".

14.2 Are the retention times of the internal standards within 30 seconds of the associated calibration standard?

ACTION: Professional judgement should be used to qualify data if the retention times differ by more than 30 seconds.

NOTE: Contractual requirements state that if any internal standard fails the acceptance criteria, the sample must be re-analyzed. If the affected sample was not re-analyzed, document in the Data Assessment under Contract Problems/Non-Compliance.

NOTE: See Notes in section 3.4, page 7 for a description of sample data the laboratory must submit.

15.0 Field Duplicates

15.1 Were any field duplicates submitted for VOA analysis?

ACTION: Compare the reported results for field duplicates and calculate the relative percent difference.

ACTION: Any gross variation between duplicate results must be addressed in the reviewer narrative. However, if large differences exist, identification of field duplicates should be confirmed by contacting the sampler.

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YES NO N/A

PART B: BNA ANALYSES

1.0 Sample Conditions/Problems

1.1 Do the Traffic Reports/Chain-of-Custody records or laboratory SDG Narrative indicate any problems with sample receipt, condition of samples, analytical problems or special notations affecting the quality of the data?

ACTION: If any sample analyzed as a soil, other than TCLP, contains 50% - 90% water, all data should be flagged as estimated "J". If a soil sample, other than TCLP, contains more than 90% water, all data should be qualified as unusable "R".

ACTION: If samples were not iced or if the ice was melted upon arrival at the laboratory and the temperature of the cooler was elevated (> 10° C), flag all positive results "J" and all non-detects "UJ".

2.0 Holding Times

2.1 Have any BNA technical holding times, determined from date of collection to date of extraction, been exceeded?

Technical Holding Time: Continuous extraction of water samples for BNA analysis must be started within seven days of the date of collection. Soil/sediment samples must be extracted within 7 days of collection. Extracts must be analyzed within 40 days of the date of extraction.

Table of Holding Time Violations  
(See Chain-of-Custody Records)

| Sample Analyzed | Sample Matrix | Date Sampled | Date Lab Received | Date Extracted | Date Analyzed |
|-----------------|---------------|--------------|-------------------|----------------|---------------|
| _____           | _____         | _____        | _____             | _____          | _____         |
| _____           | _____         | _____        | _____             | _____          | _____         |
| _____           | _____         | _____        | _____             | _____          | _____         |
| _____           | _____         | _____        | _____             | _____          | _____         |

0.35

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YES NO N/A

ACTION: If technical holding times were exceeded, flag all positive results as estimated (J) and sample quantitation limits as estimated (UJ), and document in the narrative that holding times were exceeded. If analyses were done more than 14 days beyond holding time, either on the first analysis or upon reanalysis, the reviewer must use professional judgement to determine the reliability of the data and the effects of additional storage on sample results. At a minimum, all results should be qualified "J", but the reviewer may determine that non-detect data are unusable "R". If holding times were exceeded by more than 28 days, all non-detect data must be qualified "R", unusable.

NOTE: Contractual Holding Times: Extraction of water samples must be started within 5 days VTSR. Soil/sediment samples must be extracted within 10 days of VTSR. This requirement does not apply to Performance Evaluation (PE) samples. Water and soil/sediment extracts must be analyzed within 40 days following extraction.

ACTION: If contractual holding times are exceeded, document in the Data Assessment and on the Organic Regional Data Assessment Summary form.

NOTE: The data reviewer must note in the Data Assessment whether or not technical and contractual holding times were met.

3.0 Surrogate Recovery (Form II)

3.1 Are BNA Surrogate Recovery Summaries (Form II) present for each of the following matrices:

- |               |                                     |     |                                     |
|---------------|-------------------------------------|-----|-------------------------------------|
| a. Low Water? | <input checked="" type="checkbox"/> | ___ | ___                                 |
| b. Low Soil?  | <input checked="" type="checkbox"/> | ___ | ___                                 |
| c. Med Soil?  | <input type="checkbox"/>            | ___ | <input checked="" type="checkbox"/> |

3.2 Are all the BNA samples listed on the appropriate Surrogate Recovery Summaries for each of the following matrices:

- |               |                                     |     |     |
|---------------|-------------------------------------|-----|-----|
| a. Low Water? | <input checked="" type="checkbox"/> | ___ | ___ |
| b. Low Soil?  | <input checked="" type="checkbox"/> | ___ | ___ |

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YES NO N/A

c. Med Soil?

ACTION: Contact the WAM to request an explanation or resubmittal of any missing deliverables from the laboratory. If missing deliverables are unavailable, document the effect in the Data Assessment.

3.3 Were outliers marked correctly with an asterisk?

ACTION: Circle all outliers with red pencil.

3.4 Were two or more base-neutral OR acid surrogate recoveries out of specification for any sample or method blank?

If yes, were samples reanalyzed?

Were method blanks reanalyzed?

ACTION: If all BNA surrogate recoveries are  $\geq 10\%$ , but two within the base-neutral or acid fraction do not meet SOW specifications, for the affected fraction only (i.e. acid or base-neutral compounds):

1. Flag all positive results as estimated (J).

2. Flag all non-detects as estimated detection limits ("UJ") when recoveries are less than the lower acceptance limit.

3. Do not qualify non-detects if recoveries are greater than the upper acceptance limit.

If any base-neutral or acid surrogate has a recovery of  $< 10\%$ :

1. Qualify positive results for that fraction as estimated (J).

2. Qualify non-detects for that fraction as unusable (R).

Professional judgement should be used to qualify data that have method blank surrogate recoveries out of specification in both original and reanalyses. Check the internal standard areas.

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YES NO N/A

NOTE: Contractual requirements state that if any surrogate fails acceptance criteria, the sample must be re-analyzed. If sample was not re-analyzed, document in the Data Assessment under Contract Problems/Non-Compliance.

NOTE: The laboratory must submit the following data:

1. If surrogate recoveries and internal standard responses meet the acceptance criteria in the re-analyzed sample, then the laboratory must submit only the re-analysis.

2. If surrogate recoveries and/or internal standard responses fail to meet the acceptance criteria upon re-analysis, then submit data from both analyses.

3.5 Are there any transcription/calculation errors between raw data and Form II?       

ACTION: If large errors exist, contact the WAM to request an explanation or resubmittal of corrected deliverables from the laboratory. Make necessary corrections and note errors in the Data Assessment.

4.0 Matrix Spikes (Form III)

4.1 Is the Matrix Spike/Matrix Spike Duplicate Recovery Form (Form III) present?       

4.2 Were matrix spikes analyzed at the required frequency for each of the following matrices:

a. Low Water?      

b. Low Soil?       

c. Med Soil?      

ACTION: If any matrix spike data are missing, take the action specified in 3.2 above.

4.3 How many BNA spike recoveries are outside QC limits?

Water

Soils

N/A out of 22

0 out of 22

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YES NO N/A

4.4 How many RPDs for matrix spike and matrix spike duplicate recoveries are outside QC limits?

|                     |                           |   |
|---------------------|---------------------------|---|
| <u>Water</u>        | <u>Soils</u>              | 9 |
| <u>MA</u> out of 11 | <u>0</u> out of <u>11</u> |   |

ACTION: No action is taken based upon MS/MSD data alone. However, using informed professional judgement, the data reviewer may use the matrix spike and matrix spike duplicate results in conjunction with other QC criteria and determine the need for some qualification of the data.

ACTION: Circle all outliers with red pencil.

5.0 Blanks (Form IV)

5.1 Is the Method Blank Summary (Form IV) present?

5.2 Frequency of Analysis: Has a reagent/method blank analysis been reported per 20 samples of similar matrix, or concentration level, and for each extraction batch?

5.3 Has a BNA method blank been analyzed for each GC/MS system used? (See SOW pg. D-54/SVOA, Section 12.1.2.)

ACTION: If any method blank data are missing, contact the WAM to obtain an explanation/resubmittal from the lab. If resubmittals are unavailable, use professional judgement to determine if the associated sample data should be qualified.

5.4 The validator should verify that the correct identification scheme for the EPA Blank samples were used. See page B-33, sec. 3.3.7.3 of the SOW for further information.

Was the correct identification scheme used for all BNA blanks?

ACTION: Contact the WAM to obtain resubmittals from the lab or make the required corrections on the forms. Document all corrections made by the validator in the Data Assessment under Contract Problems/Non-Compliance.

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YES NO N/A

5.5 Chromatography: review the blank raw data - chromatograms (RICs), quant. reports or data system printouts and spectra. Is the chromatographic performance (baseline stability) acceptable for each instrument?

ACTION: Use professional judgement to determine the effect on the data.

5.6 Are all detected hits for target compounds less than the CRQL for that analyte in all method blanks?

Exception: Phthalate esters must be less than five times (5x) the CRQL.

6.0 Contamination

NOTE: "Water blanks", "drill blanks" and "distilled water blanks" are validated like any other sample and are not used to qualify data. Do not confuse them with the other QC blanks discussed below.

6.1 Do any method/reagent blanks have positive results (TCL and/or TIC)?

NOTE: Water: When applied as directed in the table below (page 29), the contaminant concentration in method/instrument/reagent blanks is multiplied by the sample dilution factor, where necessary.

Soil: If the lab has not already done so, the contaminant concentration in soil blanks is multiplied by 33 times the sample dilution factor and corrected for %moisture (fraction of solid) where necessary. 30 grams of sodium sulfate (1 gram for medium level soils) are used to prepare the soil reagent/method blank as instructed on page D-54/SVOA, section 12.1.3. Contact the WAM to obtain resubmittals if the soil blanks are not reported in soil units ( $\mu\text{g}/\text{kg}$ ).

6.2 Do any field/rinse blanks have positive BNA results (TCL and/or TIC)?

ACTION: Prepare a list of samples associated with each contaminated blank. (Attach a separate sheet.)

NOTE: All field blank results associated to a particular group of samples (may exceed one per case) must be

0.40

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YES NO N/A

used to qualify sample data. Do not convert field blank results to account for the difference in soil CRQLs. Blanks may not be qualified because of contamination in another blank. Field blanks must be qualified for surrogate, spectral, instrument performance or calibration QC problems.

ACTION: Follow the directions in the table below to qualify TCL results due to contamination. Use the largest value from all the associated blanks. If gross contamination exists, all data in the associated samples should be qualified as unusable "R".

| For:                    | Flag sample result with a "U" when:            | Report CRQL & qualify "U" when:               | No qualification is needed when:              |
|-------------------------|--|---|---|
| Common Phthalate-Esters | Sample conc. is > CRQL, but ≤ 10× blank value. | Sample conc. is < CRQL and ≤ 10× blank value. | Sample conc. is > CRQL and > 10× blank value. |
| Other Contaminants      | Sample conc. is > CRQL, but ≤ 5× blank value.  | Sample conc. is < CRQL and ≤ 5× blank value.  | Sample conc. is > CRQL and > 5× blank value.  |

NOTE: Analytes qualified "U" for blank contamination are still treated as "hits" when qualifying for calibration criteria.

ACTION: For TIC compounds, if the concentration in the sample is less than five times the concentration in the most contaminated associated blank, flag the sample data "R" (unusable).

6.3 Are there field/rinse/equipment blanks associated with every sample? ✓

ACTION: For low level samples, note in the Data Assessment that there is no associated field/rinse/equipment blank. For analytes with high concentration, use professional judgement on qualification of these values and make a note in the Data Assessment.

Exception: samples taken from a drinking water tap do not have associated field blanks.

7.41



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YES NO N/A

7.0 GC/MS Instrument Performance Check

7.1 Are the GC/MS Instrument Performance Check Forms (Form V) present for Decafluorotriphenylphosphine (DFTPP)?

7.2 Are the enhanced bar graph spectrum and mass/charge (m/z) listing for the DFTPP provided for each twelve hour shift?

7.3 Has an instrument performance check solution been analyzed for every twelve hours of sample analysis per instrument?

ACTION: List date, time, instrument ID, and sample number for which no associated GC/MS tuning data are valid.

| SAMPLE NUMBERS | DATE  | TIME  | INSTRUMENT ID |
|----------------|-------|-------|---------------|
| _____          | _____ | _____ | _____         |
| _____          | _____ | _____ | _____         |
| _____          | _____ | _____ | _____         |
| _____          | _____ | _____ | _____         |

ACTION: If the WAM cannot obtain missing data from the lab, reject "R" all data generated outside an acceptable twelve hour calibration interval.

7.4 Have the ion abundances been normalized to m/z 198 (see SOW, page D-61/SVOA)?

NOTE: All ion abundance ratios must be normalized to m/z 198, the nominal base peak, even though the ion abundance of m/z 442 may up to 110% that of m/z 198.

ACTION: If mass assignment is in error, flag all associated sample data as unusable "R".

7.5 Have the ion abundance criteria been met for each instrument used?

ACTION: List all data which do not meet ion abundance criteria (attach a separate sheet).

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YES NO N/A

ACTION: If ion abundance criteria are not met, the Region II TPO must be notified.

7.6 Are there any transcription/calculation errors between mass lists and Form Vs? (Check at least two values, but if errors are found check more.)

7.7 Is the number of significant figures for the reported relative abundances consistent with the number given for each ion in the ion abundance criteria column?

ACTION: If large errors exist, take action as specified in section 3.5 above.

7.8 Are the spectra of the mass calibration compound acceptable?

ACTION: Use professional judgement to determine whether associated data should be accepted, qualified, or rejected.

3.0 Target Compound List (TCL) Analytes (FORM I SV)

8.1 Are the Organic Analysis Data Sheets (Form I SV) present with required header information on each page, for each of the following:

a. Samples and/or fractions as appropriate?

b. Matrix spikes and matrix spike duplicates?

c. Blanks?

8.2 Has GPC cleanup been performed on all soil/sediment sample extracts?

ACTION: If data suggests that GPC was not performed, use professional judgement. Make note in Contract Problems/Non-Compliance section of the Data Assessment and the Organic Regional Data Assessment Summary.

8.3 Are the BNA Reconstructed Ion Chromatograms, the mass spectra for the identified compounds, and the data system printouts (quant. reports) included in the sample package for each of the following:

a. Samples and/or fractions as appropriate?

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|  | YES                                 | NO                                  | N/A                                 |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| b. Matrix spikes and matrix spike duplicates (mass spectra not required)?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| c. Blanks?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| ACTION: If any data are missing, take action specified in 3.2 above.   |                                     |                                     |                                     |
| 8.4 Are the response factors shown in the quant. report?   | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/>            |
| 8.5 Is chromatographic performance acceptable with respect to:   |                                     |                                     |                                     |
| Baseline stability?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| Resolution?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| Peak shape?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| Full-scale graph (attenuation)?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| Other: _____?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| ACTION: Use professional judgement to determine the acceptability of the data.   |                                     |                                     |                                     |
| 8.6 Are lab-generated standard mass spectra of identified BNA compounds present for each sample?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| ACTION: If any mass spectra are missing, take action specified in 3.2 above. Note under Contract Non-compliance if lab does not generate their own standard spectra. If spectra are missing, reject all positive data. |                                     |                                     |                                     |
| 8.7 Is the RRT of each reported compound within 0.06 RRT units of the standard RRT in the continuing calibration?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 8.8 Are all ions present in the standard mass spectrum at a relative intensity greater than 10% also present in the sample mass spectrum?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 8.9 Do sample and standard relative ion intensities agree within ±20%?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| ACTION: Use professional judgement to determine acceptability of data. If it is determined that incorrect identifications were made, all   |                                     |                                     |                                     |

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YES NO N/A

such data should be rejected "R", flagged "N" (presumptive evidence of the presence of the compound) or changed to not detected "U" at the calculated detection limit. In order to be positively identified, the data must comply with the criteria listed in 8.7, 8.8, and 8.9.

ACTION: When sample carry-over is a possibility, professional judgement should be used to determine if instrument cross-contamination has affected any positive compound identification.

9.0 Tentatively Identified Compounds (TIC)

9.1 Are all Tentatively Identified Compound Forms (Form I, Part B) present; and do listed TICs include scan number or retention time, estimated concentration and "JN" qualifier?

9.2 Are the mass spectra for the tentatively identified compounds and associated "best match" spectra included in the sample package for each of the following:

a. Samples and/or fractions as appropriate?

b. Blanks?

c. Alkanes listed for each sample?

ACTION: If any TIC data are missing, take action specified in 3.2 above.

ACTION: Add "N" qualifier to all chemically named TICs, if missing.

9.3 Are any TCL compounds (from any fraction) listed as TIC compounds? (Example: 1,2-dimethylbenzene is xylene - a VOA TCL - and should not be reported as a TIC.)

ACTION: Flag with "R" any TCL compound listed as a TIC.

9.4 Are all ions present in the reference mass spectrum with a relative intensity greater than 10% also present in the sample mass spectrum?

9.5 Do TIC and "best match" standard relative ion intensities agree within ±20%?

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YES NO N/A

ACTION: Use professional judgement to determine the acceptability of TIC identifications. If it is determined that an incorrect identification was made, change the identification to "unknown," or to some less specific identification (example: "C3 substituted benzene") as appropriate. Also, when a compound is not found in any blank, but is a suspected artifact of a common laboratory contaminant, the result should be qualified as unusable, "R".

9.6 Are any TICs with responses < 10% of the internal standard (as determined by inspection of the peak areas or height) reported? \_\_\_  \_\_\_

ACTION: If yes, cross out questionable TIC(s).

10.0 Compound Quantitation and Reported Detection Limits

10.1 Are there any transcription/calculation errors in Form I results? (Check at least two positive values. Verify that the correct internal standard, quantitation ion, and RRF were used to calculate Form I result.) \_\_\_  \_\_\_

10.2 Are the CRQLs adjusted to reflect sample dilutions and, for soils, sample moisture? \_\_\_  \_\_\_

ACTION: If errors are large, take action as specified in section 3.5 above.

ACTION: When a sample is analyzed at more than one dilution, the lowest CRQLs are used (unless a QC exceedance dictates the use of the higher CRQL data from the diluted sample analysis). Replace concentrations that exceed the calibration range in the original analysis by crossing out the "E" and its associated value on the original Form I and substituting the data from the analysis of the diluted sample. Specify which Form I is to be used, then draw a red "X" across the entire page of all Form Is that should not be used, including any in the summary package.

11.0 Standards Data (GC/MS)

11.1 Are the Reconstructed Ion Chromatograms, and data system printouts (quant. reports) present for initial and continuing calibration? \_\_\_  \_\_\_

0,46

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YES NO N/A

ACTION: If any calibration standard data are missing, take action specified in 3.2 above.

12.0 GC/MS Initial Calibration (Form VI)

12.1 Are the Initial Calibration Forms (Form VI) present and complete for the BNA fraction?

ACTION: If any calibration standard forms are missing, take action specified in 3.2 above.

12.2 Are the % relative standard deviation (%RSD) values for BNAs  $\leq 30\%$  over the concentration range of the calibration?

ACTION: Circle all outliers with red pencil.

NOTE: Although 21 BNA compounds have a contractual minimum RRF and no maximum %RSD, the technical criteria are the same for all analytes.

NOTE: Eight BNA compounds do not require a 20ng standard. Refer to SOW section 7.2.4.5.1, page D-15/SVOA for a list of required compounds and contractual criteria.

ACTION: If the %RSD is  $> 30.0\%$ , qualify positive results for that analyte "J" and non-detects using professional judgement. When %RSD is  $> 90\%$ , flag all non-detect results for that analyte "R" (unusable) and all positive results "J" (estimated).

NOTE: Analytes previously qualified "U" due to blank contamination are still considered as "hits" when qualifying for calibration criteria.

12.3 Are any average RRFs  $< 0.05$ ?

ACTION: Circle all outliers with red pencil.

ACTION: If the average RRF is  $< 0.05$  then:

1. "R" all non-detects.
2. "J" all positive results.

12.4 Are there any transcription/calculation errors in the reporting of RRFs and/or %RSDs? (Check at least two values; if errors are found check more.)

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YES NO N/A

ACTION: Circle errors with red pencil.

ACTION: If errors are large, take action as specified in section 3.5 above.

NOTE: Contract Requirement: The SOW allows up to four of the required analytes to fail contractual %RSD or RRF criteria provided the %RSD is  $\leq 40\%$  or RRF is  $\geq 0.010$ . (See Table 5, page D-66/SVOA and analytes marked with a "\*" on Form VI for a list of required analytes and contractual criteria.) Technical criteria, however, are the same for all analytes.

ACTION: If more than four analytes fail %RSD or RRF criteria, document in the Data Assessment under Contract Problems/Non-Compliance and on the Organic Regional Data Assessment Summary form.

13.0 GC/MS Continuing Calibration (Form VII)

13.1 Are the Continuing Calibration Forms (Form VII) present and complete for the BNA fraction?

13.2 Has a continuing calibration standard been analyzed for every twelve hours of sample analysis per instrument?

ACTION: List below all sample analyses that were not analyzed within twelve hours of a continuing calibration standard for each instrument used.

\_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

ACTION: If any forms are missing, or no continuing calibration standard has been analyzed within twelve hours of every sample analysis, contact the WAM to obtain an explanation/resubmittal from the lab. If continuing calibration data are unavailable, flag all associated sample data as unusable "R".

13.3 Does any BNA compound have a percent difference (%D) between the initial and continuing calibration RRFs which exceeds the  $\pm 25.0\%$  criteria?

0.48

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YES NO N/A

ACTION: Circle all outliers with red pencil.

ACTION: Qualify both positive results and non-detects for the outlier compound(s) as estimated "J". When %D is > 90%, reject all non-detects for that analyte, "R", and qualify positive results "J" (estimated).

13.4 Are any continuing RRFs < 0.05?         

ACTION: Circle all outliers with red pencil.

ACTION: If the RRF is < 0.05, qualify as unusable (R) associated non-detects and "J" associated positive values.

NOTE: Contract Requirement: The SOW allows up to four of the required analytes to fail contractual %D and RRF criteria, provided that the %D is ≤ 40% and the RRF is ≥ 0.010. (See Table 5 page D-66/SVOA or analytes marked with a "\*" on Form VI for a list of the required analytes.) Technical criteria, however, are the same for all analytes.

ACTION: If more than four analytes failed %D and RRF criteria, document in the Data Assessment under Contract Problems/Non-Compliance and on the Organic Regional Data Summary Form.

13.5 Are there any transcription/calculation errors in the reporting of average relative response factors (RRF) or %difference (%D) between initial and continuing RRFs? (Check at least two values, but if errors are found, check more.)         

ACTION: Circle errors with red pencil.

ACTION: If errors are large, take action as specified in section 3.5 above.

14.0 Internal Standards (Form VIII)

14.1 Are the internal standard areas (Form VIII) of every sample and blank within the upper and lower limits (-50% to +100%) for each continuing calibration?         

If no, was sample re-analyzed?         

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YES NO N/A

- ACTION: 1. Circle all outliers with red pencil.  
2. List all the outliers below.

ACTION: If sample was not reanalyzed, document in Data Assessment in Contract Problems/Non-Compliance.

| Sample # | Internal Std. | Area  | Lower/Upper Limit |
|----------|---------------|-------|-------------------|
| _____    | _____         | _____ | _____/_____       |
| _____    | _____         | _____ | _____/_____       |
| _____    | _____         | _____ | _____/_____       |
| _____    | _____         | _____ | _____/_____       |

(Attach additional sheets if necessary.)  
(or attach copies of Form VIIIs)

ACTION: 1. If the internal standard area count is outside the "upper" or "lower" limit, flag with "J" all positive results and non-detects quantitated with this internal standard.

2. Do not qualify non-detects associated with IS area > 100%.

3. If the IS area in the sample is < 50%, qualify all analytes associated with that IS estimated (J). If area counts are extremely low (< 25% of the area in the 12 hour standard), or if performance exhibits a major abrupt drop-off, flag all associated non-detects as unusable (R) and positive hits estimated (J).

14.2 Are the retention times of the internal standards within 30 seconds of the associated calibration standard? ✓

ACTION: Professional judgement should be used to qualify data if the retention times differ by more than 30 seconds.

NOTE: Contractual requirements state that if any internal standard fails the acceptance criteria, the sample must be re-analyzed. If the affected sample was not re-analyzed, document in the Data Assessment under Contract Problems/Non-Compliance.

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YES NO N/A

NOTE: See Notes in section 3.4, page 24 for a description of sample data the laboratory must submit.

15.0 Field Duplicates

15.1 Were any field duplicates submitted for BNA analysis?

ACTION: Compare the reported results for field duplicates and calculate the relative percent difference.

ACTION: Any gross variation between field duplicate results must be addressed in the reviewer narrative. However, if large differences exist, identification of field duplicates should be confirmed by contacting the sampler.

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YES NO N/A

PART C: PESTICIDE/PCB ANALYSIS

1.0 Sample Conditions/Problems

- 1.1 Do the Traffic Reports/Chain-of-Custody Records or SDG Narrative indicate any problems with sample receipt, condition of the samples, analytical problems or special circumstances affecting the quality of the data? \_\_\_\_    \_\_\_\_

ACTION: If any sample analyzed as a soil, other than TCLP, contains 50% - 90% water, all data should be qualified as estimated "J". If a soil sample, other than TCLP, contains more than 90% water, all data should be qualified as unusable "R".

ACTION: If samples were not iced, or if the ice was melted upon arrival at the laboratory, and the temperature of the cooler was elevated > 10° C, flag all positive results "J" and all non-detects "UJ".

ACTION: Check aqueous extraction log for sample pH, if adjustment was needed, it should have been noted in the SDG Narrative. If more information is needed, notify the WAM to contact the lab.

2.0 Holding Times

- 2.1 Have any PEST/PCB technical holding times, determined from date of collection to date of extraction, been exceeded? \_\_\_\_    \_\_\_\_

NOTE: Technical Holding Times: Water and soil samples for PEST/PCB analysis must be extracted within 7 days of the date of collection. Extracts must be analyzed within 40 days of the date extraction.

ACTION: If technical holding times are exceeded, flag all positive results as estimated "J" and sample quantitation limits "UJ" and document in the narrative that holding times were exceeded. If analyses were done more than 14 days beyond holding time, either on the first analysis or upon re-analysis, the reviewer must use professional judgement to determine the reliability of the data and the effects of

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YES NO N/A

additional storage on the sample results. At a minimum, all the data should at least be qualified "J", but the reviewer may determine that non-detects are unusable "R".

Table of Holding Time Violations  
(See Chain-of-Custody Records)

| Sample Analyzed | Sample Matrix | Date Sampled | Date Lab Received | Date Extracted | Date Analyzed |
|-----------------|---------------|--------------|-------------------|----------------|---------------|
| _____           | _____         | _____        | _____             | _____          | _____         |
| _____           | _____         | _____        | _____             | _____          | _____         |
| _____           | _____         | _____        | _____             | _____          | _____         |
| _____           | _____         | _____        | _____             | _____          | _____         |

NOTE: Contractual Holding Times: Extraction of water samples must be completed within 5 days VTSR. Soil/sediment samples must be extracted within 10 days of VTSR. This requirement does not apply to Performance Evaluation (PE) samples. Extracts of water and soil/sediment samples must be analyzed within 40 days following start of extraction.

ACTION: If contractual holding times are exceeded, document in the Data Assessment and Organic Regional Data Assessment Summary form.

NOTE: The data reviewer must note in the Data Assessment whether or not technical and contractual holding times were met.

3.0 Surrogate Recovery (Form II)

3.1 Are the PEST/PCB Surrogate Recovery Summaries (Form II) present for each of the following matrices:

a. Low Water?

\_\_\_\_\_

b. Soil?

\_\_\_\_\_

3.2 Are all the PEST/PCB samples listed on the appropriate Surrogate Recovery Summary for each of the following matrices:

0.53

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YES NO N/A

a. Low Water?

b. Soil?

ACTION: Contact the WAM to obtain an explanation or resubmittal of any missing deliverables from the laboratory. If missing deliverables are unavailable, document the effect in the Data Assessment.

3.3 Were outliers marked correctly with an asterisk?

ACTION: Circle all outliers with red pencil.

3.4 Were surrogate recoveries of TCX or DCB outside of the contract specification for any sample, method blank or sulfur clean-up blank (30-150%)?

ACTION: In the absence of matrix interference, qualification of the data is not required in the following three situations:

1. When surrogates on both columns are diluted out.
2. When one surrogate on one column was outside (either above or below) the contract limits but above 10%.
3. When the same surrogate on both columns is above the contract limit.

If the same surrogate on both columns is below the contract limit but above 10%, check chromatograms for interference. The reviewer may use professional judgement, and qualify only those analytes which elute in the region of the GC chromatogram where interference was observed.

If the same surrogate on both columns is below the contract limit but above 10% (with no interference), qualify non-detects and positive hits "J" (estimated).

If recoveries for both surrogates on both columns are below the contract limit but above 10%, flag positive results and non-detects for that sample "J".

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YES NO N/A

If recoveries are above the contract limit for both surrogates on both columns, then qualify positive values "J".

If both surrogates on one column are below the contract limit but above 10%, then use the data from the other column, providing both surrogates on that column are within contract limits. The validator must check from which column the concentration is reported for each analyte. If the value is reported from the failed column, then cross it out and use the value from the other column. Document this change in the Data Assessment.

If recovery is below 10% for either surrogate on any column, qualify positive results "J" and flag non-detects "R".

3.5 Were surrogate retention times (RT) within the windows established during the initial 3-point analysis of Individual Standard Mixture A (see Form VI Pest-1)?

ACTION: If the RT limits are not met, positive results and non-detects for that sample may be qualified unusable, "R", based on professional judgement.

3.6 Are there any transcription/calculation errors between raw data and Form II?

ACTION: If large errors exist, contact the WAM to obtain an explanation or resubmittal of corrected deliverables from the laboratory. Make any necessary corrections and document the effect in the Data Assessment.

4.0 Matrix Spikes (Form III)

4.1 Is the Matrix Spike/Matrix Spike Duplicate Recovery Form (Form III) present?

4.2 Were matrix spikes analyzed at the required frequency for each of the following matrices (one MS/MSD must be performed for every 20 samples of similar matrix or concentration level):

a. Low Water? *field HLC*

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YES NO N/A

b. Soil?

ACTION: If any matrix spike data are missing, take the action specified in 3.2 above.

ACTION: Circle all outliers with red pencil.

4.3 How many PEST/PCB spike recoveries are outside QC limits?

Water

Soil

11/12 out of 12

0 out of 12

4.4 How many RPDs for matrix spike and matrix spike duplicate recoveries are outside QC limits?

Water

Soil

1/1 out of 6

0 out of 6

ACTION: No action is taken on MS/MSD data alone. However, using informed professional judgement, the data reviewer may use the matrix spike and matrix spike duplicate results in conjunction with other QC criteria and determine the need for some qualification of the data.

5.0 Blanks (Form IV)

5.1 Is the Method Blank Summary (Form IV) present?

5.2 Frequency of Analysis: Has a reagent/method blank been analyzed for each SDG, every 20 samples of similar matrix and concentration level or each extraction batch, whichever is more frequent?

ACTION: If any blank data are missing, take action as specified above in section 3.2. If blank data is not available, reject "R" all associated positive data. However, using professional judgement, the data reviewer may substitute field blank data for missing method blank data.

5.3 A separate Form IV should be present if part of an extraction batch required sulfur removal. In such cases some samples will be listed on two blank summary forms - once under the method

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YES NO N/A

blank, and once under the sulfur clean-up blank (PCBLK). Was this additional blank raw data and Form IV submitted when required?

ACTION: If sulfur clean-up blank data and Form IV are missing, take action as specified in 3.2 above.

5.4 Has a PEST/PCB instrument blank been analyzed at the beginning of every 12 hr. period following the initial calibration sequence (minimum contract requirement)?

ACTION: If any blank data are missing, take action as specified in section 3.2 above.

5.5 Was the correct identification scheme used for all Pest/PCB blanks? (See page B-33, sec. 3.3.7.3 of the SOW for further information.)

ACTION: Contact the WAM to obtain resubmittals or make the required corrections on the forms. Document in the Data Assessment under Contract Problems/Non-Compliance all corrections made by the validator.

5.6 Chromatography: review the blank raw data - chromatograms, quant. reports and data system printouts. Is the chromatographic performance (baseline stability) for each instrument acceptable?

ACTION: Use professional judgement to determine the effect on the data.

6.0 Contamination

NOTE: "Water blanks", "distilled water blanks" and "drilling water blanks" are validated like any other sample and are not used to qualify the data. Do not confuse them with the other QC blanks discussed below.

6.1 Do any method/reagent, instrument, or cleanup blanks show positive hits for pest/PCBs?

6.2 If any method blanks and/or sulfur clean-up blanks contain "hits" for target compounds, are these hits greater than the CRQL for that



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|   | YES   | NO                                  | N/A   |
|---|-------|-------------------------------------|-------|
| analyte?  | _____ | <input checked="" type="checkbox"/> | _____ |
| 6.3 In any instrument blanks, is the concentration of any target hit > 0.5 times CRQL for that analyte (see SOW, section 12.1.4.4.2, page D-77/PEST)? | _____ | <input checked="" type="checkbox"/> | _____ |

NOTE: Most labs will report 0.5 times CRQLs on the instrument blank Form I instead of the actual method CRQLs. If the lab reported the actual CRQLs, then check if any detected hits are above 0.5 times the CRQLs reported on the Form I.

ACTION: If yes to any of the above questions: note in the Data Assessment under Contract Problems/Non-Compliance if any method or clean-up blanks contain hits > the CRQL, or of instrument blank contained hits > 0.5 times CRQL for that analyte.

|   |                                     |                          |       |
|---|-------------------------------------|--------------------------|-------|
| 6.4 Do any field/rinse blanks have positive pest/PCB results? | <input checked="" type="checkbox"/> | <input type="checkbox"/> | _____ |
|---|-------------------------------------|--------------------------|-------|

ACTION: Prepare a list of the samples associated with each contaminated blank. (Attach a separate sheet)

NOTE: All field blank results associated to a particular group of samples (may exceed one per case or one per day) may be used to qualify data. Do not convert field blank results to account for the difference in soil CRQLs. Blanks may not be qualified because of contamination in another blank. Field blanks must be qualified for surrogate, and/or calibration QC problems.

ACTION: Follow the directions in the table below to qualify TCL results due to contamination. Use the largest value from all the associated blanks.

NOTE: When applied as directed in the table below, the contaminant concentration in method/instrument/reagent/cleanup blanks is multiplied by the sample dilution factor, where necessary.

If the laboratory has not already done so, the contaminant concentration in soil blanks is multiplied by 33 times the sample dilution factor and corrected for %moisture (fraction of solid) where necessary. 30 grams of sodium sulfate are used to prepare each soil reagent/method blank as instructed on page D-72/PEST, section 12.1.2.3.1. Ask the WAM

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to contact the laboratory if the soil blanks are not reported in soil units ( $\mu\text{g}/\text{kg}$ ).

| Flag sample result with a "U":                  | Report CRQL & qualify "U":                            | No qualification is needed:                     |
|---|---|---|
| Sample conc. > CRQL, but $\leq 5 \times$ blank. | Sample conc. < CRQL & is $\leq 5 \times$ blank value. | Sample conc. > CRQL & > $5 \times$ blank value. |

NOTE: If gross blank contamination exists, all data in the associated samples should be qualified as "R", unusable.

6.5 Are there field/rinse/equipment blanks associated with every sample?

ACTION: For low level samples, note in the Data Assessment that there is no associated field/rinse/equipment blank. For analytes with high concentrations, use professional judgement to qualify these values and document in the Data Assessment.

Exception: samples taken from a drinking water tap do not have associated field blanks.

7.0 Calibration and GC Performance

- 7.1 Are the following Gas Chromatograms and Data Systems Printouts for both columns present for all samples, blanks and MS/MSD:
- a. Peak resolution check?
  - b. Performance evaluation mixtures?
  - c. Aroclor 1016/1260?
  - d. Aroclors 1221, 1232, 1242, 1248, 1254?
  - e. Toxaphene?
  - f. Low points individual mixtures A & B?
  - g. Med points individual mixtures A & B?
  - h. High points individual mixtures A & B?

0.59

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|  | YES                                 | NO                                  | N/A                      |
|--|-------------------------------------|-------------------------------------|--------------------------|
| i. Instrument blanks?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/> |
| j. Were the appropriate GC columns used as specified on pg. D-11/PEST, sections 6.23.3.1 to 6.23.3.7, in the SOW?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/> |
| 7.2 Do the chromatograms for all Individual Standard Mixtures and PEM analyses display single component analytes at > 10% but < 100% of full scale (see sections 9.3.5.8.1 thru 9.3.5.8.4, pages D-32 & 33/PEST)?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/> |
| Have chromatograms for Individual Standard Mixtures and PEM analyses been replotted, showing scaling factor(s), to meet the above requirements when necessary?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/> |
| NOTE: All standard chromatograms must clearly display all peaks at > 10% but < 100% of full scale, and replotted if necessary to accommodate peaks not properly scaled in the initial chromatogram(s). Both the initial and replotted chromatograms must be submitted with the data package. |                                     |                                     |                          |
| ACTION: If all single component peaks are not clearly displayed on chromatograms for all Individual Standard Mixtures and PEM analyses, notify the WAM to obtain resubmittal of the necessary data.  |                                     |                                     |                          |
| 7.3 Are Forms VI PEST 1-7 present and complete for each column and each analytical sequence?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/> |
| ACTION: If no, take action as specified in 3.2 above.  |                                     |                                     |                          |
| 7.4 Are there any transcription/ calculation errors between raw data and Forms VI?   | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| ACTION: If large errors exist, take action as specified in section 3.6 above.  |                                     |                                     |                          |
| 7.5 Do all standard retention times, including each pesticide in each level of Individual Mixtures A & B, fall within the windows established during the Initial Calibration (see Form VI PEST-1)?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/> |
| ACTION: If no, all samples in the entire analytical sequence are potentially affected. Check to see if the chromatograms contain peaks within an expanded window surrounding the expected  |                                     |                                     |                          |

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YES NO N/A

retention times. If no peaks are found and the surrogates are visible, non-detects are valid. If peaks are present and cannot be identified through pattern recognition or using a revised RT window, qualify all positive results "JN" and non-detects as unusable (R). For aroclors, the RT may be outside the window, but the aroclor may still be identified from its distinctive pattern.

- 7.6 Are the linearity criteria for the initial analyses of Individual Standards A & B within limits for both columns? (%RSD must be  $\leq 25.0$  for alpha and delta BHC,  $\leq 30.0$  for the two surrogates and  $\leq 20\%$  for all other analytes.)

Y      

NOTE: Contractual requirements allow up to two single component TCL compounds, but not surrogates, on each column to exceed the criteria provided the %RSD is  $\leq 30\%$ . (See page D-28/Pest, sec. 9.2.5.7 in the SOW.) Technical criteria, however, are the same for all analytes.

ACTION: If technical criteria were not met, qualify all associated positive results generated during the entire analytical sequence "J" and all non-detects "UJ". When %RSD  $> 90\%$ , flag all non-detect results for that analyte "R" (unusable).

ACTION: If more than two analytes failed %RSD, document in the Data Assessment Contract Problems/Non-Compliance section and Organic Regional Data Assessment Summary form.

- 7.7 Is the resolution between each pair of adjacent peaks in the Resolution Check Mixture  $\geq 60.0\%$  for both columns? (See Form VI PEST-4.)

Y      

ACTION: If no, qualify positive results for compounds that were not adequately resolved "J". Use professional judgement to determine if non-detects which elute in areas affected by co-eluting peaks should be qualified "N" as presumptive evidence of presence or unusable (R).

- 7.8 Is Form VI PEST-5 present and complete for each Performance Evaluation Mixture (PEM) standard used for both initial and continuing calibrations (see SOW section 3.12.4.4, page B-52)?

Y      

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YES NO N/A

ACTION: If no, take action as specified in section 3.2 above.

7.9 For each PEM standard, was the resolution between each pair of adjacent peaks  $\geq 90.0\%$  on both columns?

ACTION: Qualify positive results for compounds not adequately resolved estimated (J). Qualify non-detects based on professional judgement.

7.10 Have Forms VI PEST-6 & PEST-7 been completed for all midpoint Individual Standards A and B used for initial calibration?

For each standard, was the resolution between each pair of adjacent peaks  $\geq 90.0\%$  on both columns?

ACTION: If no, qualify positive results for compounds that were not adequately resolved estimated (J). Use professional judgement to determine if non-detects which elute in areas affected by co-eluting peaks should be qualified "N" as presumptive evidence of presence or unusable "R".

7.11 Is Form VII Pest-1 present and complete for each PEM standard analyzed during the analytical sequence for both columns?

Was the %Breakdown of DDT and Endrin calculated using the equations given on page D-26/PEST, sec. 9.2.4.8 in the SOW?

Were all pesticides and surrogates in each PEM standard within the RT windows established during the Initial Calibration?

ACTION: If no, take action as specified in 3.2 above.

7.12 Has the individual percent breakdown for DDT/Endrin exceeded 20.0% in any PEM on either column? (See Form VII PEST-1.)

- for 4,4'-DDT?

- for Endrin?

Has the combined percent breakdown for DDT/Endrin

0.62

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YES NO N/A

exceeded 30.0% in any PEM on either column  
(required for all PEM analyses)?

  ✓  

- ACTION: 1. If any percent breakdown has failed the QC criteria in either PEM in steps 2 and 17 in the initial calibration sequence (page D-28/Pest, sec. 9.2.5.6 in the SOW), qualify all samples in the entire analytical sequence as described in sections 2.a, b and c below.
2. If any percent breakdown failed the QC criteria in a PEM calibration verification analysis, review data beginning with the samples which followed the last in-control standard until the next acceptable PEM and qualify the data as described below.
- a. 4,4'-DDT Breakdown: If DDT breakdown was > 20.0%:
- i. Qualify all positive results for DDT with "J". If DDT was not detected, but DDD and DDE are positive, then qualify the quantitation limit for DDT unusable, "R".
  - ii. Qualify positive results for DDD and/or DDE as presumptively present at an approximated quantity "JN".
- b. Endrin Breakdown: If endrin breakdown was > 20.0%:
- i. Qualify all positive results for endrin with "J". If endrin was not detected, but endrin aldehyde and endrin ketone are positive, then qualify the quantitation limit for Endrin as unusable "R".
  - ii. Qualify positive results for endrin ketone and endrin aldehyde as presumptively present at an approximated quantity "JN".
- c. Combined Breakdown: If the combined 4,4'-DDT and endrin breakdown is greater than 30.0%:
- i. Qualify all positive results for DDT and Endrin with "J". If endrin was not detected, but endrin aldehyde and endrin ketone are positive, then qualify the quantitation limit for endrin as unusable

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"R". If DDT was not detected, but DDD and DDE are positive, then qualify the quantitation limit for DDT as unusable "R".

- ii. Qualify positive results for endrin ketone and endrin aldehyde as presumptively present at an approximated quantity "JN". Qualify positive results for DDD and/or DDE as presumptively present at an approximated quantity "JN".

- 7.13 Are all percent difference (%D) values for PEM analytes and surrogates on both columns  $\geq -25\%$  and  $\leq +25.0\%$ ? (See Form VII PEST-1.)

YES  NO  N/A

ACTION: If no, qualify all associated positive results generated during the analytical sequence "J" and sample quantitation limits "UJ".

NOTE: If the failing PEM is part of the initial calibration, all samples are potentially affected. If the offending standard is a calibration verification, the associated samples are those which followed the last in-control standard until the next passing standard.

- 7.14 Is Form VII Pest-2 present and complete for each INDA and INDB calibration verification analyzed?

YES  NO  N/A

ACTION: If no, take action specified in 3.2 above.

- 7.15 Are there any transcription/calculation errors between raw data and Form VII Pest-2?

YES  NO  N/A

ACTION: If large errors exists, take action as specified in section 3.6 above.

- 7.16 Do all standard retention times for each INDA and INDB calibration verification fall within the RT windows established during the initial calibration sequence? (See Form VII PEST-2.)

YES  NO  N/A

ACTION: If no, beginning with the samples which followed the last in-control standard, check to see if the chromatograms contain peaks within an expanded window surrounding the expected retention times. If no peaks are found and the surrogates are visible, non-detects are valid. If peaks are present and cannot be identified through pattern recognition or using a revised

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YES NO N/A

RT window, qualify all positive results and non-detects as unusable (R).

7.17 Are all %D values for INDA and INDB calibration verification compounds  $\geq -25.0\%$  and  $\leq +25.0\%$ ?

ACTION: If the %D is outside the  $\pm 25.0\%$  range for any compound(s), qualify associated positive results for that compound "J" and non-detects "UJ". The "associated samples" are those which followed the last in-control standard up to the next passing standard containing the analyte(s) in question. If the %D is  $> 90\%$ , flag all non-detects for that analyte "R" (unusable).

8.0 Analytical Sequence Check (Form VIII-PEST)

8.1 Is Form VIII present and complete for each column and each period of analyses?

ACTION: If no, take action specified in 3.2 above.

8.2 Was the proper analytical sequence followed for each initial calibration and subsequent analyses, and all standards analyzed at the required frequency for each GC/EC instrument used? (See SOW pages D-23 & D-58/PEST.)

ACTION: If no, use professional judgement to determine the severity of the effect on the data and qualify accordingly. Generally, the effect is negligible unless the sequence was grossly altered and/or the calibration was out of QC limits.

8.3 Were all samples analyzed within a 12 hour time period beginning with the injection of an instrument blank and bracketed by acceptable analyses of the proper standards?

ACTION: If no, use professional judgement to determine the severity of the effect on the data and qualify accordingly. Document in the Data Assessment under Contract Problems/Non-Compliance and Organic Regional Data Assessment Summary.

8.4 If a multi-component analyte was detected in a sample, was a matching multi-component standard analyzed within 72 hours of the injection of the

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YES NO N/A

sample and within a valid 12 hour sequence?

NOTE: This additional standard is for identification purposes only. Positive results for Aroclors and Toxaphene are quantitated from the initial calibration.

ACTION: If no, document in the Data Assessment under Contract Problems/Non-Compliance and on the Organic Regional Data Assessment Summary form.

9.0 Cleanup Efficiency Verification (Form IX)

9.1 Is Form IX PEST-1 present and complete for each lot of Florisil Cartridges used? (Florisil Cleanup is required for all Pest/PCB extracts.)

Are all samples listed on the Pesticide Florisil Cartridge Check Form?

ACTION: If no, take action specified in 3.2 above. If data suggests florisil clean-up was not performed, document in the Data Assessment under the Contract Non-compliance section.

9.2 Are percent recoveries (%REC) of the pesticide and surrogate compounds used to check the efficiency of the florisil clean-up procedure within QC limits of 80 - 120%?

ACTION: Qualify only the analyte(s) which failed the recovery criteria as follows:

If %REC is < 80%, qualify positive results "J" and non-detects "UJ".

If any pesticide %REC was zero, flag non-detects "R" for that compound.

Use professional judgement to qualify positive results if any recoveries are > 120%.

NOTE: Sample data should be evaluated for potential interferences if recovery of 2,4,5-trichlorophenol was > 5% in the Florisil Cartridge Performance Check analysis. Document any problems found in the Data Assessment under the Contract Problems/Non-Compliance section.

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YES NO N/A

9.3 If GPC Cleanup was performed (mandatory for all soil sample extracts), is Form IX Pest-2 present?

Are all soil samples listed on Form IX Pest-2?

ACTION: If no, take action specified in 3.2 above. If data suggests GPC clean-up was not performed when required, document in the Data Assessment under the Contract Problems/Non-Compliance section and Organic Regional Data Assessment Summary.

Are the %REC values for all pesticides in the GPC calibration solution between 80 - 110%?

ACTION: Qualify only those analytes which failed the recovery criteria as follows:

If %REC are < 80%, qualify positive results "J" and non-detects "UJ".

If any pesticide %REC was zero, flag non-detects "R" for that compound.

Use professional judgement to qualify positive results if any recoveries are > 110%.

NOTE: An Aroclor mixture containing Aroclors 1016 and 1260 is also analyzed during GPC calibration; however, Aroclor data is not listed on Form IX PEST-2. The raw GPC data for Aroclors 1016/1260 must be evaluated for pattern similarity with previously analyzed Aroclor standards.

9.4 The validator should verify that the correct identification scheme for the EPA Blank samples were used. See page B-35, sec. 3.3.7.8 and 3.3.7.9 of the SOW for further information.

Was the correct identification scheme used for GPC and Florisil blanks?

10.0 Pesticide/PCB Identification

10.1 Is Form X complete for every sample in which a pesticide or PCB was detected?

ACTION: If no, take action specified in 3.2 above.

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|  | YES                                 | NO                                  | N/A                                 |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| 10.2 Are all sample chromatograms properly scaled, attenuated, etc. as required for proper identification of single and multi-component analytes? (Refer to SOW sections 11.3.7.1 thru 11.3.7.8, page D-70/Pest for specific details.)   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| NOTE: Proper verification of Pest/PCB results depends on clear, legible presentation of the raw data. Single component pesticides and all peaks chosen for quantitation of multi-component analytes must appear at less than full scale. Toxaphene and PCB patterns must be clearly visible to enable comparison with standard chromatograms.  |                                     |                                     |                                     |
| ACTION: If retention times or apex of peaks cannot be verified, or if multi-component peak patterns cannot be discerned, contact the WAM to obtain rescaled chromatograms from the lab.  |                                     |                                     |                                     |
| 10.3 Are there any transcription/calculation errors between raw data and Forms 10A and 10B?  | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/>            |
| ACTION: If large errors exist, take action as specified in section 3.6 above.  |                                     |                                     |                                     |
| 10.4 Are RTs of sample compounds within the established RT windows for analyses on both columns?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| Was GC/MS confirmation provided when required (when compound concentration is > 10 ug/ml in the final extract)?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| ACTION: Use professional judgement to qualify positive results which were not confirmed by GC/MS analysis. Qualify as unusable (R) all positive results which were not confirmed on a second GC column. Also qualify as unusable (R) all positive results which do not meet RT window criteria, unless associated standard compounds are similarly biased. Use professional judgement to assign an appropriate quantitation limit. |                                     |                                     |                                     |
| 10.5 Is the percent difference (%D) calculated for the positive sample results on both columns > 25.0%?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| ACTION: If the reviewer finds neither column shows interference for the positive hits, the data should be flagged as follows:  |                                     |                                     |                                     |

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YES NO N/A

| <u>% Difference</u>                 | <u>Qualifier</u> |
|-------------------------------------|------------------|
| 0 - 25%                             | None             |
| 25 - 70%                            | "J"              |
| 70 - 100%                           | "JN"             |
| > 100% (No interference)            | "R"              |
| 100 - 200% (Interference detected)* | "JN"             |
| > 50% (Pesticide value is < CRQL)** | "U"              |
| > 200%                              | "R"              |

\* When the reported %D is 100 - 200%, but interference is detected on either column, qualify the data with "J".

\*\* When the reported pesticide value is lower than the CRQL, and the %D is > 50%, raise the value to the CRQL and qualify "U", undetected.

NOTE: For Aroclors, if the %D is > 50%, but the pattern of GC peaks on both columns indicates a specific Aroclor is present, qualify that Aroclor "J".

NOTE: The lower of the two values is reported on Form I. If using professional judgement, the reviewer determines that the higher result was more acceptable, the reviewer should replace the value and indicate the reason for the change in the Data Assessment.

10.6 Check chromatograms for false negatives, especially the multiple-peak compounds (Toxaphene and the PCBs). Were there any false negatives?

ACTION: Use professional judgement to decide if the compound should be reported. If the appropriate PCB standards were not analyzed within 72 hrs. of the sample(s) in question, qualify the data unusable "R".

Also note in Data Assessment under Contract Problems/Non-Compliance if the lab failed to analyze Aroclor standards when required.

11.0 Target Compound List (TCL) Analytes

11.1 Are the Organic Analysis Data Sheets (Form I Pest) present with required header information on each page, for each of the following:

- a. Samples and/or fractions as appropriate?
- b. Matrix spikes and matrix spike duplicates?

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|  | YES                                 | NO                                  | N/A                                 |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| c. Blanks?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| d. Instrument Blanks (per column & analysis)?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 11.2 Are the Pest chromatograms and quant. reports included in the sample data package for each of the following:  |                                     |                                     |                                     |
| a. Samples and/or fractions as appropriate?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| b. Matrix spikes and matrix spike duplicates?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| c. Blanks?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| d. Instrument Blanks (per column & analysis)?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| ACTION: If any data are missing, take action specified in 3.2 above.   |                                     |                                     |                                     |
| 11.3 Are the calibration factors shown in the quant. reports?  | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/>            |
| 11.4 Is chromatographic performance acceptable with respect to:  |                                     |                                     |                                     |
| a. Baseline stability?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| b. Resolution?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| c. Peak shape?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| d. Full-scale graph attenuation?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| e. Other: _____?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| 11.5 Were any electropositive displacement (negative peaks) or unusual peaks seen?   | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/>            |
| ACTION: Use professional judgement to determine the acceptability of the data. Address comments under System Performance section of the Data Assessment. |                                     |                                     |                                     |

12.0 Compound Quantitation and Reported Detection Limits

|  |                          |                                     |                          |
|--|--------------------------|-------------------------------------|--------------------------|
| 12.1 Are there any transcription/calculation errors in Form I results? Check at least two positive results. Were any errors found? | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
|--|--------------------------|-------------------------------------|--------------------------|

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YES NO N/A

NOTE: Single-peak pesticide results can be checked for rough agreement between quantitative results obtained on the two GC columns. Use professional judgement to decide whether a large discrepancy indicates the presence of an interfering compound. If an interfering compound is visible on the chromatogram, the lower of the two values should be reported and qualified as presumptively present at an approximated quantity "JN". This necessitates a determination of an estimated concentration on the confirmation column. The narrative should indicate that the presence of interferences has interfered with the evaluation of the second column confirmation.

12.2 Are the CRQLs adjusted to reflect sample dilutions?

ACTION: If large errors exist, take action as specified in section 3.6 above.

ACTION: When a sample is analyzed at more than one dilution, the lowest CRQLs are used (unless a QC exceedance dictates the use of the higher CRQLs from the diluted sample). Replace concentrations which exceed the calibration range in the original analysis by crossing out the "E" value on the original Form I and substituting it with the result from the diluted sample. Specify which Form I is to be used, then draw a red "X" across the entire page of all Form I's that should not be used, including those in the data summary package.

ACTION: Quantitation limits affected by large, off-scale peaks should be qualified as unusable (R). If the interference is on-scale, the reviewer may offer an approximated quantitation limit (UJ) for each affected compound.

NOTE: If a sample required greater than a 10 times dilution, then a 10 times more concentrated analysis must also be performed and submitted (see SOW, page D-60/PEST, section 10.2.3.5).

ACTION: If a more concentrated analysis is unavailable, document in the Contract Problems/Non-Compliance section of the Data Assessment. Use professional judgement to qualify non-detects and positive hits below the CRQL.

STANDARD OPERATING PROCEDURE

US EPA Region II  
Method: CLP/SOW OLMO3.2

Date: June 1996  
SOP HW-6, Rev. 11

YES NO N/A

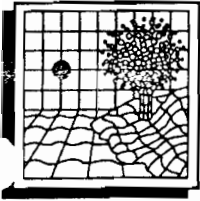
13.0 Field Duplicates

13.1 Were any field duplicates submitted?

ACTION: Compare the reported results for field duplicates and calculate the relative percent difference.

ACTION: Any gross variation between field duplicate results must be addressed in the reviewer narrative. However, if large differences exist, identification of field duplicates should be confirmed by contacting the sampler.

NOV 30 2000



# AMERICAN ANALYTICAL & TECHNICAL SERVICES, INC.

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000001

## SDG NARRATIVE

Laboratory Name: **AATSLA**, Baton Rouge, LA  
Case No.: 28706  
SDG No.: BNP05  
Contract: 68-W0-0081

Samples were received November 8<sup>th</sup>, 2000, for organic analyses following the USEPA Contract Laboratory Program (CLP) Statement of Work (SOW) OLM04.2 and the terms of contract 68W00081. The samples are as listed below in Table I. Also included in this table are pH values obtained from water sample volatile containers measured just following the analysis of each sample (where applicable).

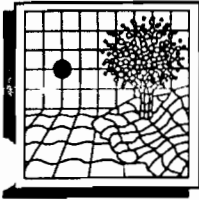
### TABLE I

**EPA Sample ID:**

BNP05  
BNP05MS  
BNP05MSD  
BNP06  
BNP07  
BNP08  
BNP09  
BNP10  
BNP11  
BNP12  
BNP13  
BNQ85  
BNQ87  
BNQ88  
BNQ89  
BNQ91  
BNQ92

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000002

**General:** The samples of this SDG were received intact and on ice. The temperature of the samples at receipt was recorded as 5 to 6°C as determined by a non contact thermometer (Raytek Raynger model ST-2). Cooler temperature indicators were not supplied with the coolers. No problems were recorded with the receipt of the samples.

## **Volatile Organic Analysis:**

All calibration, surrogate, and internal standard control criteria were within contractual limits with the exception of a surrogate for the MS of sample BNP05. A reanalysis is not required for a matrix spike sample. Recovery outliers were also observed for the MS/MSD set of sample BNP05. All volatile holding times were met.

No additional billables are submitted for VOAs:

A 75m x 0.53mm ID, 3 $\mu$ m film DB-624 J&W Bonded, fused silica capillary column was utilized for analysis. Also a 30.5 cm Supelco Purge Trap "I" (VOCARB 4000) consisting of Carboxen 1000 & 1001

## **Semi-Volatile Organic Analysis:**

All calibration, internal standard, matrix spike, and surrogate control criteria were within normal limits. All semivolatile holding times were met.

Two (2) additional billables are submitted for BNAs as follows:

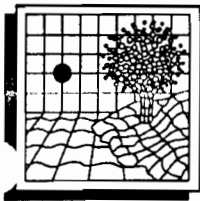
BNP08DL (Dilution)  
BNP09DL (Dilution)

Dilutions were necessary for these two samples to bring target analyte(s) into the instrument calibration range. The original is submitted for added value.

Column used for semivolatile analysis:

Restek XTI-5 (Bonded fused silica, 5% Phenyl, 95% Methyl silicone), 0.25mm internal diameter, 0.25 $\mu$ m film thickness, 30m length.

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000003

## Pesticide Organic Analysis:

The Columns used for Pesticide analysis are:

- a. CLPESTICII - Restek Proprietary 0.53 mm ID, 30 m length.
- b. CLPESTIC - Restek Proprietary 0.53 mm ID, 30 m length

## Qualifications:

Volatile and semivolatile target analytes were evaluated and reported to a quantification report limit of 1.0. Compounds that were detected below this level were assessed only if they otherwise were detected as a TIC.

Note that results for xylenes as presented on quantitation reports are not calculated in accordance to the requirements of the SOW. Corrections in compliance with the SOW are manually generated and are documented on applicable quantitation reports.

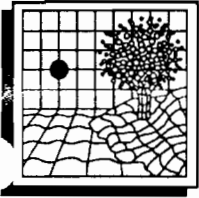
Dilutions are generally performed to bring target analytes into instrument calibration range. A second level of dilution may have been reported (if performed) to provide added value. Dilutions performed for other reasons are explained elsewhere in this narrative.

Volatile soil samples analyzed using the Method 5035 technique are indicated by the word "Encore" (or like notation) on the header of the quantification report. Those performed using soil from a jar are those generally analyzed with exact sample weights of 5.0g and the word "jar" written on the quantification report header.

Quantification limits for Encore samples containing more than 5.0 grams are arbitrarily cut off at the CRQL of 10 as required by the SOW.

Manual integrations, if necessary, were performed to correct errors of automated software routines. They are represented by an "m" flag on the quantification reports and are each accompanied by hardcopy graphic displays initialed and dated by the analyst or reviewer performing the manual integration.

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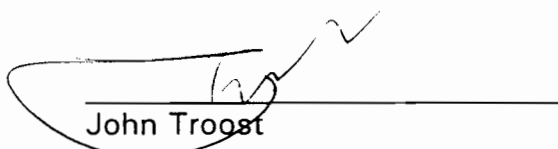


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000004

"I certify that this data package is in compliance with the terms and conditions of the contract, both technically and for completeness, for other than the conditions detailed above. Release of the data contained in this hardcopy data package and in the computer-readable data submitted on diskette has been authorized by the laboratory manager or his/her designee, as verified by the following signature."

  
\_\_\_\_\_  
John Troost  
GC/MS Supervisor  
November 28<sup>th</sup>, 2000  
JT/jj

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000005

AATSLA  
GC/MS LABORATORY : ALKANE REPORT

REPORT DATE: 11/27/00 DATA FILE: 319F08.D  
SAMPLE ID: BNP05 MATRIX: SOIL  
LAB ID: 45798.01 DATE ANALYZED: 11/14/00

| CAS #          | COMPOUND | RT    | EST. CONC<br>(UG/KG ) |
|----------------|----------|-------|-----------------------|
| 1) 000112-95-8 | EICOSANE | 12.02 | 175                   |
| 2) 000112-95-8 | EICOSANE | 12.76 | 212                   |

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000006

AATSLA  
GC/MS LABORATORY : ALKANE REPORT

REPORT DATE: 11/27/00

DATA FILE: 319F11.D

SAMPLE ID: BNP06

MATRIX: SOIL

LAB ID: 45798.04

DATE ANALYZED: 11/14/00

| CAS #          | COMPOUND         | RT    | EST. CONC<br>(UG/KG ) |
|----------------|------------------|-------|-----------------------|
| 1) 000112-95-8 | EICOSANE         | 12.03 | 181                   |
| 2) 000630-07-9 | PENTATRIACONTANE | 12.76 | 208                   |

000007

AATSLA  
GC/MS LABORATORY : ALKANE REPORT

REPORT DATE: 11/27/00

DATA FILE: 319F14.D

SAMPLE ID: BNP09

MATRIX: SOIL

LAB ID: 45798.07

DATE ANALYZED: 11/14/00

| CAS #           | COMPOUND                     | RT    | EST. CONC<br>(UG/KG ) |
|-----------------|------------------------------|-------|-----------------------|
| 1) 000544-76-3  | HEXADECANE                   | 5.67  | 146                   |
| 2) 055045-07-3  | DODECANE, 2-METHYL-8-PROPYL- | 5.96  | 93                    |
| 3) 000629-97-0  | DOCOSANE                     | 9.02  | 104                   |
| 4) 000638-67-5  | TRICOSANE                    | 9.50  | 149                   |
| 5) 000646-31-1  | TETRACOSANE                  | 9.98  | 233                   |
| 6) 000630-06-8  | HEXATRIACONTANE              | 10.43 | 401                   |
| 7) 000629-94-7  | HENEICOSANE                  | 10.86 | 642                   |
| 8) 000629-78-7  | HEPTADECANE                  | 11.27 | 570                   |
| 9) 000630-01-3  | HEXACOSANE                   | 11.68 | 563                   |
| 10) 000629-92-5 | NONADECANE                   | 12.06 | 591                   |
| 11) 000112-95-8 | EICOSANE                     | 12.79 | 411                   |
| 12) 013475-77-9 | EICOSANE, 9-OCTYL-           | 13.14 | 355                   |
| 13) 000629-94-7 | HENEICOSANE                  | 13.50 | 217                   |
| 1) 000593-49-7  | HEPTACOSANE                  | 14.33 | 90                    |

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000008

AATSLA  
GC/MS LABORATORY : ALKANE REPORT

REPORT DATE: 11/27/00 DATA FILE: 322F04.D  
SAMPLE ID: BNP09DL MATRIX: SOIL  
LAB ID: 45798.07DL DATE ANALYZED: 11/17/00

| CAS #          | COMPOUND              | RT    | EST. CONC<br>(UG/KG ) |
|----------------|-----------------------|-------|-----------------------|
| 1) 000646-31-1 | TETRACOSANE           | 9.94  | 504                   |
| 2) 013475-75-7 | PENTADECANE, 8-HEXYL- | 10.39 | 895                   |
| 3) 000629-97-0 | DOCOSANE              | 10.82 | 1381                  |
| 4) 000629-78-7 | HEPTADECANE           | 11.24 | 1222                  |
| 5) 000629-94-7 | HENEICOSANE           | 11.64 | 1282                  |
| 6) 000593-45-3 | OCTADECANE            | 12.03 | 1498                  |
| 7) 000630-06-8 | HEXATRIACONTANE       | 12.77 | 1054                  |
| 8) 000112-95-8 | EICOSANE              | 13.12 | 841                   |
| 9) 000112-95-8 | EICOSANE              | 13.47 | 570                   |

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AATSLA  
GC/MS LABORATORY : ALKANE REPORT

REPORT DATE: 11/27/00

DATA FILE: 322F05.D

SAMPLE ID: BNP13

MATRIX: SOIL

LAB ID: 45798.11

DATE ANALYZED: 11/17/00

| CAS #          | COMPOUND                | RT   | EST. CONC<br>(UG/KG ) |
|----------------|-------------------------|------|-----------------------|
| 1) 017301-33-6 | UNDECANE, 4,8-DIMETHYL- | 4.86 | 100                   |



AATSLA

Sample Delivery Group (SDG)  
Cover Sheet

SDG Number: BNP05

Laboratory Name: American Analytical & Technical Services, Incorporated Laboratory Code AATSLA

Contract No. 68-W0-0081 Case No. 28706

Analysis Price \_\_\_\_\_ SDG Turnaround 21 Days

EPA Sample Numbers in SDG (Listed in Numerical Order)

|    |       |     |       |     |       |     |
|----|-------|-----|-------|-----|-------|-----|
| 1) | BNP05 | 7)  | BNP11 | 13) | BNQ89 | 19) |
| 2) | BNP06 | 8)  | BNP12 | 14) | BNQ91 | 20) |
| 3) | BNP07 | 9)  | BNP13 | 15) | BNQ92 | 21) |
| 4) | BNP08 | 10) | BNQ85 | 16) |       | 22) |
| 5) | BNP09 | 11) | BNQ87 | 17) |       | 23) |
| 6) | BNP10 | 12) | BNQ88 | 18) |       | 24) |

BNP05

First Sample in SDG

BNQ92

Last Sample in SDG

11/08/00

First Sample Receipt Date

11/08/00

Last Sample Receipt Date

**Note:** There are a maximum of 20 field samples (excluding PE samples) in an SDG. Attach TRs to this form in alphanumeric order (the order listed above on this form).

Signature Nancy LeBlanc

Date 11/09/00

000029

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP05

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) SOIL Lab Sample ID: 45798.01  
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: 315A10.D  
 Level: (low/med) LOW Date Received: 11/08/00  
 % Moisture: not dec. 27<sup>v</sup> Date Analyzed: 11/10/00  
 GC Column: DB-624 ID: 0.53 (mm) Dilution Factor: 1.0  
 Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

CAS NO. COMPOUND

| CAS NO.   | COMPOUND                              | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) UG/KG | Q   |
|-----------|---------------------------------------|---|-----|
| 75-71-8   | Dichlorodifluoromethane               | 14  | U   |
| 74-87-3   | Chloromethane                         | 14  | U   |
| 75-01-4   | Vinyl Chloride                        | 14  | U   |
| 74-83-9   | Bromomethane                          | 14  | U   |
| 75-00-3   | Chloroethane                          | 14  | U   |
| 75-69-4   | Trichlorofluoromethane                | 14  | U   |
| 75-35-4   | 1,1-Dichloroethene                    | 14  | U   |
| 76-13-1   | 1,1,2-Trichloro-1,2,2-trifluoroethane | 14  | U   |
| 67-64-1   | Acetone                               | 24  | U B |
| 75-15-0   | Carbon Disulfide                      | 2   | J   |
| 79-20-9   | Methyl Acetate                        | 14  | U   |
| 75-09-2   | Methylene Chloride                    | 14  | U   |
| 156-60-5  | trans-1,2-Dichloroethene              | 14  | U   |
| 1634-04-4 | Methyl tert-Butyl Ether               | 14  | U   |
| 75-34-3   | 1,1-Dichloroethane                    | 14  | U   |
| 156-59-2  | cis-1,2-Dichloroethene                | 14  | U   |
| 78-93-3   | 2-Butanone                            | 11  | J   |
| 67-66-3   | Chloroform                            | 14  | U   |
| 71-55-6   | 1,1,1-Trichloroethane                 | 14  | U   |
| 110-82-7  | Cyclohexane                           | 14  | U   |
| 56-23-5   | Carbon Tetrachloride                  | 14  | U   |
| 71-43-2   | Benzene                               | 14  | U   |
| 107-06-2  | 1,2-Dichloroethane                    | 14  | U   |

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000030

1B  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP05

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) SOIL Lab Sample ID: 45798.01  
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: 315A10.D  
 Level: (low/med) LOW Date Received: 11/08/00  
 % Moisture: not dec. 27 Date Analyzed: 11/10/00  
 GC Column: DB-624 ID: 0.53 (mm) Dilution Factor: 1.0  
 Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO. COMPOUND

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

Q

| CAS NO.    | COMPOUND                    | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) UG/KG | Q |
|------------|-----------------------------|---|---|
| 79-01-6    | Trichloroethene             | 14  | U |
| 108-87-2   | Methylcyclohexane           | 14  | U |
| 78-87-5    | 1,2-Dichloropropane         | 14  | U |
| 75-27-4    | Bromodichloromethane        | 14  | U |
| 10061-01-5 | cis-1,3-Dichloropropene     | 14  | U |
| 108-10-1   | 4-Methyl-2-Pentanone        | 14  | U |
| 108-88-3   | Toluene                     | 14  | U |
| 10061-02-6 | trans-1,3-Dichloropropene   | 14  | U |
| 79-00-5    | 1,1,2-Trichloroethane       | 14  | U |
| 127-18-4   | Tetrachloroethene           | 14  | U |
| 591-78-6   | 2-Hexanone                  | 14  | U |
| 124-48-1   | Dibromochloromethane        | 14  | U |
| 106-93-4   | 1,2-Dibromoethane           | 14  | U |
| 108-90-7   | Chlorobenzene               | 14  | U |
| 100-41-4   | Ethylbenzene                | 14  | U |
| 1330-20-7  | Xylene (total)              | 14  | U |
| 100-42-5   | Styrene                     | 14  | U |
| 75-25-2    | Bromoform                   | 14  | U |
| 98-82-8    | Isopropylbenzene            | 14  | U |
| 79-34-5    | 1,1,2,2-Tetrachloroethane   | 14  | U |
| 541-73-1   | 1,3-Dichlorobenzene         | 14  | U |
| 106-46-7   | 1,4-Dichlorobenzene         | 14  | U |
| 95-50-1    | 1,2-Dichlorobenzene         | 14  | U |
| 96-12-8    | 1,2-Dibromo-3-chloropropane | 14  | U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 14  | U |

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1F  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

BNP05

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.01

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: 315A10.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: not dec. 27

Date Analyzed: 11/10/00

GC Column: DB-624 ID: 0.53 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs found: 0

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

| CAS NUMBER | COMPOUND NAME | RT | EST. CONC. | Q |
|------------|---------------|----|------------|---|
| 1.         |               |    |            |   |
| 2.         |               |    |            |   |
| 3.         |               |    |            |   |
| 4.         |               |    |            |   |
| 5.         |               |    |            |   |
| 6.         |               |    |            |   |
| 7.         |               |    |            |   |
| 8.         |               |    |            |   |
| 9.         |               |    |            |   |
| 10.        |               |    |            |   |
| 11.        |               |    |            |   |
| 12.        |               |    |            |   |
| 13.        |               |    |            |   |
| 14.        |               |    |            |   |
| 15.        |               |    |            |   |
| 16.        |               |    |            |   |
| 17.        |               |    |            |   |
| 18.        |               |    |            |   |
| 19.        |               |    |            |   |
| 20.        |               |    |            |   |
| 21.        |               |    |            |   |
| 22.        |               |    |            |   |
| 23.        |               |    |            |   |
| 24.        |               |    |            |   |
| 25.        |               |    |            |   |
| 26.        |               |    |            |   |
| 27.        |               |    |            |   |
| 28.        |               |    |            |   |
| 29.        |               |    |            |   |
| 30.        |               |    |            |   |

2,85

000037

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP06

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) SOIL Lab Sample ID: 45798.04  
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: 315A11.D  
 Level: (low/med) LOW Date Received: 11/08/00  
 % Moisture: not dec. 28 ✓ Date Analyzed: 11/10/00  
 GC Column: DB-624 ID: 0.53 (mm) Dilution Factor: 1.0  
 Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NO. COMPOUND

| CAS NO.   | COMPOUND                              | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) UG/KG | Q |
|-----------|---------------------------------------|---|---|
| 75-71-8   | Dichlorodifluoromethane               | 14  | U |
| 74-87-3   | Chloromethane                         | 14  | U |
| 75-01-4   | Vinyl Chloride                        | 14  | U |
| 74-83-9   | Bromomethane                          | 14  | U |
| 75-00-3   | Chloroethane                          | 14  | U |
| 75-69-4   | Trichlorofluoromethane                | 14  | U |
| 75-35-4   | 1,1-Dichloroethene                    | 14  | U |
| 76-13-1   | 1,1,2-Trichloro-1,2,2-trifluoroethane | 14  | U |
| 67-64-1   | Acetone                               | 59  | U |
| 75-15-0   | Carbon Disulfide                      | 14  | U |
| 79-20-9   | Methyl Acetate                        | 14  | U |
| 75-09-2   | Methylene Chloride                    | 14  | U |
| 156-60-5  | trans-1,2-Dichloroethene              | 14  | U |
| 1634-04-4 | Methyl tert-Butyl Ether               | 14  | U |
| 75-34-3   | 1,1-Dichloroethane                    | 14  | U |
| 156-59-2  | cis-1,2-Dichloroethene                | 14  | U |
| 78-93-3   | 2-Butanone                            | 29  | U |
| 67-66-3   | Chloroform                            | 14  | U |
| 71-55-6   | 1,1,1-Trichloroethane                 | 14  | U |
| 110-82-7  | Cyclohexane                           | 14  | U |
| 56-23-5   | Carbon Tetrachloride                  | 14  | U |
| 71-43-2   | Benzene                               | 14  | U |
| 107-06-2  | 1,2-Dichloroethane                    | 14  | U |

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000038

1B  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP06

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) SOIL Lab Sample ID: 45798.04  
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: 315A11.D  
 Level: (low/med) LOW Date Received: 11/08/00  
 % Moisture: not dec. 28 Date Analyzed: 11/10/00  
 GC Column: DB-624 ID: 0.53 (mm) Dilution Factor: 1.0  
 Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO. COMPOUND

 CONCENTRATION UNITS:  
 (ug/L or ug/Kg) UG/KG Q

|            |                             |    |   |
|------------|-----------------------------|----|---|
| 79-01-6    | Trichloroethene             | 14 | U |
| 108-87-2   | Methylcyclohexane           | 14 | U |
| 78-87-5    | 1,2-Dichloropropane         | 14 | U |
| 75-27-4    | Bromodichloromethane        | 14 | U |
| 10061-01-5 | cis-1,3-Dichloropropene     | 14 | U |
| 108-10-1   | 4-Methyl-2-Pentanone        | 14 | U |
| 108-88-3   | Toluene                     | 14 | U |
| 10061-02-6 | trans-1,3-Dichloropropene   | 14 | U |
| 79-00-5    | 1,1,2-Trichloroethane       | 14 | U |
| 127-18-4   | Tetrachloroethene           | 14 | U |
| 591-78-6   | 2-Hexanone                  | 14 | U |
| 124-48-1   | Dibromochloromethane        | 14 | U |
| 106-93-4   | 1,2-Dibromoethane           | 14 | U |
| 108-90-7   | Chlorobenzene               | 14 | U |
| 100-41-4   | Ethylbenzene                | 14 | U |
| 1330-20-7  | Xylene (total)              | 14 | U |
| 100-42-5   | Styrene                     | 14 | U |
| 75-25-2    | Bromoform                   | 14 | U |
| 98-82-8    | Isopropylbenzene            | 14 | U |
| 79-34-5    | 1,1,2,2-Tetrachloroethane   | 14 | U |
| 541-73-1   | 1,3-Dichlorobenzene         | 14 | U |
| 106-46-7   | 1,4-Dichlorobenzene         | 14 | U |
| 95-50-1    | 1,2-Dichlorobenzene         | 14 | U |
| 96-12-8    | 1,2-Dibromo-3-chloropropane | 14 | U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 14 | U |

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1F  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

BNP06

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) SOIL Lab Sample ID: 45798.04  
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: 315A11.D  
 Level: (low/med) LOW Date Received: 11/08/00  
 % Moisture: not dec. 28 Date Analyzed: 11/10/00  
 GC Column: DB-624 ID: 0.53 (mm) Dilution Factor: 1.0  
 Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 Number TICs found: 0 CONCENTRATION UNITS:  
 (ug/L or ug/Kg) UG/KG

| CAS NUMBER | COMPOUND NAME | RT | EST. CONC. | Q |
|------------|---------------|----|------------|---|
| 1.         |               |    |            |   |
| 2.         |               |    |            |   |
| 3.         |               |    |            |   |
| 4.         |               |    |            |   |
| 5.         |               |    |            |   |
| 6.         |               |    |            |   |
| 7.         |               |    |            |   |
| 8.         |               |    |            |   |
| 9.         |               |    |            |   |
| 10.        |               |    |            |   |
| 11.        |               |    |            |   |
| 12.        |               |    |            |   |
| 13.        |               |    |            |   |
| 14.        |               |    |            |   |
| 15.        |               |    |            |   |
| 16.        |               |    |            |   |
| 17.        |               |    |            |   |
| 18.        |               |    |            |   |
| 19.        |               |    |            |   |
| 20.        |               |    |            |   |
| 21.        |               |    |            |   |
| 22.        |               |    |            |   |
| 23.        |               |    |            |   |
| 24.        |               |    |            |   |
| 25.        |               |    |            |   |
| 26.        |               |    |            |   |
| 27.        |               |    |            |   |
| 28.        |               |    |            |   |
| 29.        |               |    |            |   |
| 30.        |               |    |            |   |

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1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP07

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) SOIL Lab Sample ID: 45798.05  
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: 315A12.D  
 Level: (low/med) LOW Date Received: 11/08/00  
 % Moisture: not dec. 21 ✓ Date Analyzed: 11/10/00  
 GC Column: DB-624 ID: 0.53 (mm) Dilution Factor: 1.0  
 Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

| CAS NO.   | COMPOUND                              | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) UG/KG | Q                |
|-----------|---------------------------------------|---|------------------|
| 75-71-8   | Dichlorodifluoromethane               | 13  | U                |
| 74-87-3   | Chloromethane                         | 13  | U                |
| 75-01-4   | Vinyl Chloride                        | 13  | U                |
| 74-83-9   | Bromomethane                          | 13  | U                |
| 75-00-3   | Chloroethane                          | 13  | U                |
| 75-69-4   | Trichlorofluoromethane                | 13  | U                |
| 75-35-4   | 1,1-Dichloroethene                    | 13  | U                |
| 76-13-1   | 1,1,2-Trichloro-1,2,2-trifluoroethane | 13  | U                |
| 67-64-1   | Acetone                               | 13 4  | U B <del>+</del> |
| 75-15-0   | Carbon Disulfide                      | 13  | U                |
| 79-20-9   | Methyl Acetate                        | 13  | U                |
| 75-09-2   | Methylene Chloride                    | 13 3  | U B <del>+</del> |
| 156-60-5  | trans-1,2-Dichloroethene              | 13  | U                |
| 1634-04-4 | Methyl tert-Butyl Ether               | 13  | U                |
| 75-34-3   | 1,1-Dichloroethane                    | 13  | U                |
| 156-59-2  | cis-1,2-Dichloroethene                | 13  | U                |
| 78-93-3   | 2-Butanone                            | 13  | U                |
| 67-66-3   | Chloroform                            | 13  | U                |
| 71-55-6   | 1,1,1-Trichloroethane                 | 13  | U                |
| 110-82-7  | Cyclohexane                           | 13  | U                |
| 56-23-5   | Carbon Tetrachloride                  | 13  | U                |
| 71-43-2   | Benzene                               | 13  | U                |
| 107-06-2  | 1,2-Dichloroethane                    | 13  | U                |

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1B  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP07

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) SOIL Lab Sample ID: 45798.05  
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: 315A12.D  
 Level: (low/med) LOW Date Received: 11/08/00  
 % Moisture: not dec. 21 Date Analyzed: 11/10/00  
 GC Column: DB-624 ID: 0.53 (mm) Dilution Factor: 1.0  
 Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NO. COMPOUND

Q

| CAS NO.    | COMPOUND                    | UG/KG | Q |
|------------|-----------------------------|-------|---|
| 79-01-6    | Trichloroethene             | 13    | U |
| 108-87-2   | Methylcyclohexane           | 13    | U |
| 78-87-5    | 1,2-Dichloropropane         | 13    | U |
| 75-27-4    | Bromodichloromethane        | 13    | U |
| 10061-01-5 | cis-1,3-Dichloropropene     | 13    | U |
| 108-10-1   | 4-Methyl-2-Pentanone        | 13    | U |
| 108-88-3   | Toluene                     | 13    | U |
| 10061-02-6 | trans-1,3-Dichloropropene   | 13    | U |
| 79-00-5    | 1,1,2-Trichloroethane       | 13    | U |
| 127-18-4   | Tetrachloroethene           | 13    | U |
| 591-78-6   | 2-Hexanone                  | 13    | U |
| 124-48-1   | Dibromochloromethane        | 13    | U |
| 106-93-4   | 1,2-Dibromoethane           | 13    | U |
| 108-90-7   | Chlorobenzene               | 13    | U |
| 100-41-4   | Ethylbenzene                | 13    | U |
| 1330-20-7  | Xylene (total)              | 13    | U |
| 100-42-5   | Styrene                     | 13    | U |
| 75-25-2    | Bromoform                   | 13    | U |
| 98-82-8    | Isopropylbenzene            | 13    | U |
| 79-34-5    | 1,1,2,2-Tetrachloroethane   | 13    | U |
| 541-73-1   | 1,3-Dichlorobenzene         | 13    | U |
| 106-46-7   | 1,4-Dichlorobenzene         | 13    | U |
| 95-50-1    | 1,2-Dichlorobenzene         | 13    | U |
| 96-12-8    | 1,2-Dibromo-3-chloropropane | 13    | U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 13    | U |

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1F  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

BNP07

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.05

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: 315A12.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: not dec. 21

Date Analyzed: 11/10/00

GC Column: DB-624 ID: 0.53 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs found: 2

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

| CAS NUMBER  | COMPOUND NAME                | RT    | EST. CONC. | Q  |
|-------------|------------------------------|-------|------------|----|
| 1. 556-67-2 | CYCLOTETRASILOXANE, OCTAMETH | 14.11 | 10         | JN |
| 2.          | UNKNOWN                      | 16.75 | 10         | J  |
| 3.          |                              |       |            |    |
| 4.          |                              |       |            |    |
| 5.          |                              |       |            |    |
| 6.          |                              |       |            |    |
| 7.          |                              |       |            |    |
| 8.          |                              |       |            |    |
| 9.          |                              |       |            |    |
| 10.         |                              |       |            |    |
| 11.         |                              |       |            |    |
| 12.         |                              |       |            |    |
| 13.         |                              |       |            |    |
| 14.         |                              |       |            |    |
| 15.         |                              |       |            |    |
| 16.         |                              |       |            |    |
| 17.         |                              |       |            |    |
| 18.         |                              |       |            |    |
| 19.         |                              |       |            |    |
| 20.         |                              |       |            |    |
| 21.         |                              |       |            |    |
| 22.         |                              |       |            |    |
| 23.         |                              |       |            |    |
| 24.         |                              |       |            |    |
| 25.         |                              |       |            |    |
| 26.         |                              |       |            |    |
| 27.         |                              |       |            |    |
| 28.         |                              |       |            |    |
| 29.         |                              |       |            |    |
| 30.         |                              |       |            |    |

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1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP08

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.06

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: 315A13.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: not dec. 3 ✓

Date Analyzed: 11/10/00

GC Column: DB-624 ID: 0.53 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

| CAS NO.   | COMPOUND                              | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) UG/KG | Q               |
|-----------|---------------------------------------|---|-----------------|
| 75-71-8   | Dichlorodifluoromethane               | 10  | U               |
| 74-87-3   | Chloromethane                         | 10  | U               |
| 75-01-4   | Vinyl Chloride                        | 10  | U               |
| 74-83-9   | Bromomethane                          | 10  | U               |
| 75-00-3   | Chloroethane                          | 10  | U               |
| 75-69-4   | Trichlorofluoromethane                | 10  | U               |
| 75-35-4   | 1,1-Dichloroethene                    | 10  | U               |
| 76-13-1   | 1,1,2-Trichloro-1,2,2-trifluoroethane | 10  | U               |
| 67-64-1   | Acetone                               | 10 <del>±</del>                               | U <del>B+</del> |
| 75-15-0   | Carbon Disulfide                      | 10  | U               |
| 79-20-9   | Methyl Acetate                        | 10  | U               |
| 75-09-2   | Methylene Chloride                    | 10 <del>±</del>                               | U <del>B+</del> |
| 156-60-5  | trans-1,2-Dichloroethene              | 10  | U               |
| 1634-04-4 | Methyl tert-Butyl Ether               | 10  | U               |
| 75-34-3   | 1,1-Dichloroethane                    | 10  | U               |
| 156-59-2  | cis-1,2-Dichloroethene                | 10  | U               |
| 78-93-3   | 2-Butanone                            | 10  | U               |
| 67-66-3   | Chloroform                            | 10  | U               |
| 71-55-6   | 1,1,1-Trichloroethane                 | 10  | U               |
| 110-82-7  | Cyclohexane                           | 10  | U               |
| 56-23-5   | Carbon Tetrachloride                  | 10  | U               |
| 71-43-2   | Benzene                               | 10  | U               |
| 107-06-2  | 1,2-Dichloroethane                    | 10  | U               |

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1B  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP08

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) SOIL Lab Sample ID: 45798.06  
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: 315A13.D  
 Level: (low/med) LOW Date Received: 11/08/00  
 % Moisture: not dec. 3 Date Analyzed: 11/10/00  
 GC Column: DB-624 ID: 0.53 (mm) Dilution Factor: 1.0  
 Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO. COMPOUND

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

Q

| CAS NO.    | COMPOUND                    | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) UG/KG | Q |
|------------|-----------------------------|---|---|
| 79-01-6    | Trichloroethene             | 10  | U |
| 108-87-2   | Methylcyclohexane           | 10  | U |
| 78-87-5    | 1,2-Dichloropropane         | 10  | U |
| 75-27-4    | Bromodichloromethane        | 10  | U |
| 10061-01-5 | cis-1,3-Dichloropropene     | 10  | U |
| 108-10-1   | 4-Methyl-2-Pentanone        | 10  | U |
| 108-88-3   | Toluene                     | 10  | U |
| 10061-02-6 | trans-1,3-Dichloropropene   | 10  | U |
| 79-00-5    | 1,1,2-Trichloroethane       | 10  | U |
| 127-18-4   | Tetrachloroethene           | 10  | U |
| 591-78-6   | 2-Hexanone                  | 10  | U |
| 124-48-1   | Dibromochloromethane        | 10  | U |
| 106-93-4   | 1,2-Dibromoethane           | 10  | U |
| 108-90-7   | Chlorobenzene               | 10  | U |
| 100-41-4   | Ethylbenzene                | 10  | U |
| 1330-20-7  | Xylene (total)              | 10  | U |
| 100-42-5   | Styrene                     | 10  | U |
| 75-25-2    | Bromoform                   | 10  | U |
| 98-82-8    | Isopropylbenzene            | 10  | U |
| 79-34-5    | 1,1,2,2-Tetrachloroethane   | 10  | U |
| 541-73-1   | 1,3-Dichlorobenzene         | 10  | U |
| 106-46-7   | 1,4-Dichlorobenzene         | 10  | U |
| 95-50-1    | 1,2-Dichlorobenzene         | 10  | U |
| 96-12-8    | 1,2-Dibromo-3-chloropropane | 10  | U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 10  | U |

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1F  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

BNP08

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.06

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: 315A13.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: not dec. 3

Date Analyzed: 11/10/00

GC Column: DB-624 ID: 0.53 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs found: 4

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

| CAS NUMBER  | COMPOUND NAME                | RT    | EST. CONC. | Q  |
|-------------|------------------------------|-------|------------|----|
| 1. 556-67-2 | CYCLOTETRASILOXANE, OCTAMETH | 14.10 | 29         | JN |
| 2. 526-73-8 | BENZENE, 1,2,3-TRIMETHYL-    | 14.30 | 7          | JN |
| 3. 95-63-6  | BENZENE, 1,2,4-TRIMETHYL-    | 14.82 | 10         | JN |
| 4. _____    | UNKNOWN                      | 16.73 | 7          |    |
| 5. _____    |                              |       |            |    |
| 6. _____    |                              |       |            |    |
| 7. _____    |                              |       |            |    |
| 8. _____    |                              |       |            |    |
| 9. _____    |                              |       |            |    |
| 10. _____   |                              |       |            |    |
| 11. _____   |                              |       |            |    |
| 12. _____   |                              |       |            |    |
| 13. _____   |                              |       |            |    |
| 14. _____   |                              |       |            |    |
| 15. _____   |                              |       |            |    |
| 16. _____   |                              |       |            |    |
| 17. _____   |                              |       |            |    |
| 18. _____   |                              |       |            |    |
| 19. _____   |                              |       |            |    |
| 20. _____   |                              |       |            |    |
| 21. _____   |                              |       |            |    |
| 22. _____   |                              |       |            |    |
| 23. _____   |                              |       |            |    |
| 24. _____   |                              |       |            |    |
| 25. _____   |                              |       |            |    |
| 26. _____   |                              |       |            |    |
| 27. _____   |                              |       |            |    |
| 28. _____   |                              |       |            |    |
| 29. _____   |                              |       |            |    |
| 30. _____   |                              |       |            |    |

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1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP09

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) SOIL Lab Sample ID: 45798.07  
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: 315A14.D  
 Level: (low/med) LOW Date Received: 11/08/00  
 % Moisture: not dec. 10 ✓ Date Analyzed: 11/10/00  
 GC Column: DB-624 ID: 0.53 (mm) Dilution Factor: 1.0  
 Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

| CAS NO.   | COMPOUND                              | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) UG/KG | Q |
|-----------|---------------------------------------|---|---|
| 75-71-8   | Dichlorodifluoromethane               | 11  | U |
| 74-87-3   | Chloromethane                         | 11  | U |
| 75-01-4   | Vinyl Chloride                        | 11  | U |
| 74-83-9   | Bromomethane                          | 11  | U |
| 75-00-3   | Chloroethane                          | 11  | U |
| 75-69-4   | Trichlorofluoromethane                | 11  | U |
| 75-35-4   | 1,1-Dichloroethene                    | 11  | U |
| 76-13-1   | 1,1,2-Trichloro-1,2,2-trifluoroethane | 11  | U |
| 67-64-1   | Acetone                               | 11  | U |
| 75-15-0   | Carbon Disulfide                      | 11  | U |
| 79-20-9   | Methyl Acetate                        | 11  | U |
| 75-09-2   | Methylene Chloride                    | 11  | U |
| 156-60-5  | trans-1,2-Dichloroethene              | 11  | U |
| 1634-04-4 | Methyl tert-Butyl Ether               | 11  | U |
| 75-34-3   | 1,1-Dichloroethane                    | 11  | U |
| 156-59-2  | cis-1,2-Dichloroethene                | 11  | U |
| 78-93-3   | 2-Butanone                            | 11  | U |
| 67-66-3   | Chloroform                            | 11  | U |
| 71-55-6   | 1,1,1-Trichloroethane                 | 11  | U |
| 110-82-7  | Cyclohexane                           | 11  | U |
| 56-23-5   | Carbon Tetrachloride                  | 11  | U |
| 71-43-2   | Benzene                               | 11  | U |
| 107-06-2  | 1,2-Dichloroethane                    | 11  | U |

0.45

000064

1B  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP09

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) SOIL Lab Sample ID: 45798.07  
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: 315A14.D  
 Level: (low/med) LOW Date Received: 11/08/00  
 % Moisture: not dec. 10 Date Analyzed: 11/10/00  
 GC Column: DB-624 ID: 0.53 (mm) Dilution Factor: 1.0  
 Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO. COMPOUND

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

Q

|            |                             |    |   |
|------------|-----------------------------|----|---|
| 79-01-6    | Trichloroethene             | 11 | U |
| 108-87-2   | Methylcyclohexane           | 11 | U |
| 78-87-5    | 1,2-Dichloropropane         | 11 | U |
| 75-27-4    | Bromodichloromethane        | 11 | U |
| 10061-01-5 | cis-1,3-Dichloropropene     | 11 | U |
| 108-10-1   | 4-Methyl-2-Pentanone        | 11 | U |
| 108-88-3   | Toluene                     | 11 | U |
| 10061-02-6 | trans-1,3-Dichloropropene   | 11 | U |
| 79-00-5    | 1,1,2-Trichloroethane       | 11 | U |
| 127-18-4   | Tetrachloroethene           | 11 | U |
| 591-78-6   | 2-Hexanone                  | 11 | U |
| 124-48-1   | Dibromochloromethane        | 11 | U |
| 106-93-4   | 1,2-Dibromoethane           | 11 | U |
| 108-90-7   | Chlorobenzene               | 11 | U |
| 100-41-4   | Ethylbenzene                | 11 | U |
| 1330-20-7  | Xylene (total)              | 11 | U |
| 100-42-5   | Styrene                     | 11 | U |
| 75-25-2    | Bromoform                   | 11 | U |
| 98-82-8    | Isopropylbenzene            | 11 | U |
| 79-34-5    | 1,1,2,2-Tetrachloroethane   | 11 | U |
| 541-73-1   | 1,3-Dichlorobenzene         | 11 | U |
| 106-46-7   | 1,4-Dichlorobenzene         | 11 | U |
| 95-50-1    | 1,2-Dichlorobenzene         | 11 | U |
| 96-12-8    | 1,2-Dibromo-3-chloropropane | 11 | U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 11 | U |

p 90

000065

EPA SAMPLE NO.

1F  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

BNP09

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.07

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: 315A14.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: not dec. 10

Date Analyzed: 11/10/00

GC Column: DB-624 ID: 0.53 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs found: 4

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

| CAS NUMBER  | COMPOUND NAME             | RT    | EST. CONC. | Q  |
|-------------|---------------------------|-------|------------|----|
| 1.          | UNKNOWN                   | 14.11 | 88         | J  |
| 2. 95-63-6  | BENZENE, 1,2,4-TRIMETHYL- | 14.32 | 10         | JN |
| 3. 526-73-8 | BENZENE, 1,2,3-TRIMETHYL- | 14.83 | 15         | JN |
| 4.          | UNKNOWN                   | 16.75 | 23         | J  |
| 5.          |                           |       |            |    |
| 6.          |                           |       |            |    |
| 7.          |                           |       |            |    |
| 8.          |                           |       |            |    |
| 9.          |                           |       |            |    |
| 10.         |                           |       |            |    |
| 11.         |                           |       |            |    |
| 12.         |                           |       |            |    |
| 13.         |                           |       |            |    |
| 14.         |                           |       |            |    |
| 15.         |                           |       |            |    |
| 16.         |                           |       |            |    |
| 17.         |                           |       |            |    |
| 18.         |                           |       |            |    |
| 19.         |                           |       |            |    |
| 20.         |                           |       |            |    |
| 21.         |                           |       |            |    |
| 22.         |                           |       |            |    |
| 23.         |                           |       |            |    |
| 24.         |                           |       |            |    |
| 25.         |                           |       |            |    |
| 26.         |                           |       |            |    |
| 27.         |                           |       |            |    |
| 28.         |                           |       |            |    |
| 29.         |                           |       |            |    |
| 30.         |                           |       |            |    |

297



000074

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP10

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) SOIL Lab Sample ID: 45798.08  
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: 315A15.D  
 Level: (low/med) LOW Date Received: 11/08/00  
 % Moisture: not dec. 19 Date Analyzed: 11/10/00  
 GC Column: DB-624 ID: 0.53 (mm) Dilution Factor: 1.0  
 Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

| CAS NO.   | COMPOUND                              | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) UG/KG | Q |
|-----------|---------------------------------------|---|---|
| 75-71-8   | Dichlorodifluoromethane               | 12  | U |
| 74-87-3   | Chloromethane                         | 12  | U |
| 75-01-4   | Vinyl Chloride                        | 12  | U |
| 74-83-9   | Bromomethane                          | 12  | U |
| 75-00-3   | Chloroethane                          | 12  | U |
| 75-69-4   | Trichlorofluoromethane                | 12  | U |
| 75-35-4   | 1,1-Dichloroethene                    | 12  | U |
| 76-13-1   | 1,1,2-Trichloro-1,2,2-trifluoroethane | 12  | U |
| 67-64-1   | Acetone                               | 12  | U |
| 75-15-0   | Carbon Disulfide                      | 12  | U |
| 79-20-9   | Methyl Acetate                        | 12  | U |
| 75-09-2   | Methylene Chloride                    | 12  | U |
| 156-60-5  | trans-1,2-Dichloroethene              | 12  | U |
| 1634-04-4 | Methyl tert-Butyl Ether               | 12  | U |
| 75-34-3   | 1,1-Dichloroethane                    | 12  | U |
| 156-59-2  | cis-1,2-Dichloroethene                | 12  | U |
| 78-93-3   | 2-Butanone                            | 12  | U |
| 67-66-3   | Chloroform                            | 12  | U |
| 71-55-6   | 1,1,1-Trichloroethane                 | 12  | U |
| 110-82-7  | Cyclohexane                           | 12  | U |
| 56-23-5   | Carbon Tetrachloride                  | 12  | U |
| 71-43-2   | Benzene                               | 12  | U |
| 107-06-2  | 1,2-Dichloroethane                    | 12  | U |

000075

1B  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP10

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) SOIL Lab Sample ID: 45798.08  
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: 315A15.D  
 Level: (low/med) LOW Date Received: 11/08/00  
 % Moisture: not dec. 19 Date Analyzed: 11/10/00  
 GC Column: DB-624 ID: 0.53 (mm) Dilution Factor: 1.0  
 Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

CAS NO. COMPOUND

| CAS NO.    | COMPOUND                    | UG/KG | Q |
|------------|-----------------------------|-------|---|
| 79-01-6    | Trichloroethene             | 12    | U |
| 108-87-2   | Methylcyclohexane           | 12    | U |
| 78-87-5    | 1,2-Dichloropropane         | 12    | U |
| 75-27-4    | Bromodichloromethane        | 12    | U |
| 10061-01-5 | cis-1,3-Dichloropropene     | 12    | U |
| 108-10-1   | 4-Methyl-2-Pentanone        | 12    | U |
| 108-88-3   | Toluene                     | 12    | U |
| 10061-02-6 | trans-1,3-Dichloropropene   | 12    | U |
| 79-00-5    | 1,1,2-Trichloroethane       | 12    | U |
| 127-18-4   | Tetrachloroethene           | 12    | U |
| 591-78-6   | 2-Hexanone                  | 12    | U |
| 124-48-1   | Dibromochloromethane        | 12    | U |
| 106-93-4   | 1,2-Dibromoethane           | 12    | U |
| 108-90-7   | Chlorobenzene               | 12    | U |
| 100-41-4   | Ethylbenzene                | 12    | U |
| 1330-20-7  | Xylene (total)              | 12    | U |
| 100-42-5   | Styrene                     | 12    | U |
| 75-25-2    | Bromoform                   | 12    | U |
| 98-82-8    | Isopropylbenzene            | 12    | U |
| 79-34-5    | 1,1,2,2-Tetrachloroethane   | 12    | U |
| 541-73-1   | 1,3-Dichlorobenzene         | 12    | U |
| 106-46-7   | 1,4-Dichlorobenzene         | 12    | U |
| 95-50-1    | 1,2-Dichlorobenzene         | 12    | U |
| 96-12-8    | 1,2-Dibromo-3-chloropropane | 12    | U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 12    | U |

000076

1F  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

BNP10

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.08

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: 315A15.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: not dec. 19

Date Analyzed: 11/10/00

GC Column: DB-624 ID: 0.53 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs found: 2

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

| CAS NUMBER | COMPOUND NAME | RT    | EST. CONC. | Q |
|------------|---------------|-------|------------|---|
| 1.         | UNKNOWN       | 14.10 | 20         | J |
| 2.         | UNKNOWN       | 16.74 | 14         | J |
| 3.         |               |       |            |   |
| 4.         |               |       |            |   |
| 5.         |               |       |            |   |
| 6.         |               |       |            |   |
| 7.         |               |       |            |   |
| 8.         |               |       |            |   |
| 9.         |               |       |            |   |
| 10.        |               |       |            |   |
| 11.        |               |       |            |   |
| 12.        |               |       |            |   |
| 13.        |               |       |            |   |
| 14.        |               |       |            |   |
| 15.        |               |       |            |   |
| 16.        |               |       |            |   |
| 17.        |               |       |            |   |
| 18.        |               |       |            |   |
| 19.        |               |       |            |   |
| 20.        |               |       |            |   |
| 21.        |               |       |            |   |
| 22.        |               |       |            |   |
| 23.        |               |       |            |   |
| 24.        |               |       |            |   |
| 25.        |               |       |            |   |
| 26.        |               |       |            |   |
| 27.        |               |       |            |   |
| 28.        |               |       |            |   |
| 29.        |               |       |            |   |
| 30.        |               |       |            |   |

000082

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP11

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) SOIL Lab Sample ID: 45798.09  
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: 315A16.D  
 Level: (low/med) LOW Date Received: 11/08/00  
 % Moisture: not dec. 20 Date Analyzed: 11/10/00  
 GC Column: DB-624 ID: 0.53 (mm) Dilution Factor: 1.0  
 Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

CAS NO. COMPOUND

| CAS NO.   | COMPOUND                              | CONCENTRATION UNITS | UG/KG | Q    |
|-----------|---------------------------------------|---------------------|-------|------|
| 75-71-8   | Dichlorodifluoromethane               | 13                  |       | U    |
| 74-87-3   | Chloromethane                         | 13                  |       | U    |
| 75-01-4   | Vinyl Chloride                        | 13                  |       | U    |
| 74-83-9   | Bromomethane                          | 13                  |       | U    |
| 75-00-3   | Chloroethane                          | 13                  |       | U    |
| 75-69-4   | Trichlorofluoromethane                | 13                  |       | U    |
| 75-35-4   | 1,1-Dichloroethene                    | 13                  |       | U    |
| 76-13-1   | 1,1,2-Trichloro-1,2,2-trifluoroethane | 13                  |       | U    |
| 67-64-1   | Acetone                               | 13                  | 3     | U BF |
| 75-15-0   | Carbon Disulfide                      | 13                  |       | U    |
| 79-20-9   | Methyl Acetate                        | 13                  |       | U    |
| 75-09-2   | Methylene Chloride                    | 13                  | 3     | U BF |
| 156-60-5  | trans-1,2-Dichloroethene              | 13                  |       | U    |
| 1634-04-4 | Methyl tert-Butyl Ether               | 13                  |       | U    |
| 75-34-3   | 1,1-Dichloroethane                    | 13                  |       | U    |
| 156-59-2  | cis-1,2-Dichloroethene                | 13                  |       | U    |
| 78-93-3   | 2-Butanone                            | 13                  |       | U    |
| 67-66-3   | Chloroform                            | 13                  |       | U    |
| 71-55-6   | 1,1,1-Trichloroethane                 | 13                  |       | U    |
| 110-82-7  | Cyclohexane                           | 13                  |       | U    |
| 56-23-5   | Carbon Tetrachloride                  | 13                  |       | U    |
| 71-43-2   | Benzene                               | 13                  |       | U    |
| 107-06-2  | 1,2-Dichloroethane                    | 13                  |       | U    |

11/01

000083

1B  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP11

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) SOIL Lab Sample ID: 45798.09  
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: 315A16.D  
 Level: (low/med) LOW Date Received: 11/08/00  
 % Moisture: not dec. 20 Date Analyzed: 11/10/00  
 GC Column: DB-624 ID: 0.53 (mm) Dilution Factor: 1.0  
 Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NO. COMPOUND

| CAS NO.    | COMPOUND                    | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) UG/KG | Q |
|------------|-----------------------------|---|---|
| 79-01-6    | Trichloroethene             | 13  | U |
| 108-87-2   | Methylcyclohexane           | 13  | U |
| 78-87-5    | 1,2-Dichloropropane         | 13  | U |
| 75-27-4    | Bromodichloromethane        | 13  | U |
| 10061-01-5 | cis-1,3-Dichloropropene     | 13  | U |
| 108-10-1   | 4-Methyl-2-Pentanone        | 13  | U |
| 108-88-3   | Toluene                     | 13  | U |
| 10061-02-6 | trans-1,3-Dichloropropene   | 13  | U |
| 79-00-5    | 1,1,2-Trichloroethane       | 13  | U |
| 127-18-4   | Tetrachloroethene           | 13  | U |
| 591-78-6   | 2-Hexanone                  | 13  | U |
| 124-48-1   | Dibromochloromethane        | 13  | U |
| 106-93-4   | 1,2-Dibromoethane           | 13  | U |
| 108-90-7   | Chlorobenzene               | 13  | U |
| 100-41-4   | Ethylbenzene                | 13  | U |
| 1330-20-7  | Xylene (total)              | 13  | U |
| 100-42-5   | Styrene                     | 13  | U |
| 75-25-2    | Bromoform                   | 13  | U |
| 98-82-8    | Isopropylbenzene            | 13  | U |
| 79-34-5    | 1,1,2,2-Tetrachloroethane   | 13  | U |
| 541-73-1   | 1,3-Dichlorobenzene         | 13  | U |
| 106-46-7   | 1,4-Dichlorobenzene         | 13  | U |
| 95-50-1    | 1,2-Dichlorobenzene         | 13  | U |
| 96-12-8    | 1,2-Dibromo-3-chloropropane | 13  | U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 13  | U |

000084

1F  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

BNP11

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) SOIL Lab Sample ID: 45798.09  
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: 315A16.D  
 Level: (low/med) LOW Date Received: 11/08/00  
 % Moisture: not dec. 20 Date Analyzed: 11/10/00  
 GC Column: DB-624 ID: 0.53 (mm) Dilution Factor: 1.0  
 Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 Number TICs found: 0 CONCENTRATION UNITS:  
 (ug/L or ug/Kg) UG/KG

| CAS NUMBER | COMPOUND NAME | RT | EST. CONC. | Q |
|------------|---------------|----|------------|---|
| 1.         |               |    |            |   |
| 2.         |               |    |            |   |
| 3.         |               |    |            |   |
| 4.         |               |    |            |   |
| 5.         |               |    |            |   |
| 6.         |               |    |            |   |
| 7.         |               |    |            |   |
| 8.         |               |    |            |   |
| 9.         |               |    |            |   |
| 10.        |               |    |            |   |
| 11.        |               |    |            |   |
| 12.        |               |    |            |   |
| 13.        |               |    |            |   |
| 14.        |               |    |            |   |
| 15.        |               |    |            |   |
| 16.        |               |    |            |   |
| 17.        |               |    |            |   |
| 18.        |               |    |            |   |
| 19.        |               |    |            |   |
| 20.        |               |    |            |   |
| 21.        |               |    |            |   |
| 22.        |               |    |            |   |
| 23.        |               |    |            |   |
| 24.        |               |    |            |   |
| 25.        |               |    |            |   |
| 26.        |               |    |            |   |
| 27.        |               |    |            |   |
| 28.        |               |    |            |   |
| 29.        |               |    |            |   |
| 30.        |               |    |            |   |

2103

000089

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP12

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) SOIL Lab Sample ID: 45798.10  
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: 315A17.D  
 Level: (low/med) LOW Date Received: 11/08/00  
 % Moisture: not dec. 21 Date Analyzed: 11/10/00  
 GC Column: DB-624 ID: 0.53 (mm) Dilution Factor: 1.0  
 Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

| CAS NO.   | COMPOUND                              |      |     |
|-----------|---------------------------------------|------|-----|
| 75-71-8   | Dichlorodifluoromethane               | 13   | U   |
| 74-87-3   | Chloromethane                         | 13   | U   |
| 75-01-4   | Vinyl Chloride                        | 13   | U   |
| 74-83-9   | Bromomethane                          | 13   | U   |
| 75-00-3   | Chloroethane                          | 13   | U   |
| 75-69-4   | Trichlorofluoromethane                | 13   | U   |
| 75-35-4   | 1,1-Dichloroethene                    | 13   | U   |
| 76-13-1   | 1,1,2-Trichloro-1,2,2-trifluoroethane | 13   | U   |
| 67-64-1   | Acetone                               | 27   | U B |
| 75-15-0   | Carbon Disulfide                      | 13   | U   |
| 79-20-9   | Methyl Acetate                        | 13   | U   |
| 75-09-2   | Methylene Chloride                    | 13 2 | U B |
| 156-60-5  | trans-1,2-Dichloroethene              | 13   | U   |
| 1634-04-4 | Methyl tert-Butyl Ether               | 13   | U   |
| 75-34-3   | 1,1-Dichloroethane                    | 13   | U   |
| 156-59-2  | cis-1,2-Dichloroethene                | 13   | U   |
| 78-93-3   | 2-Butanone                            | 9    | J   |
| 67-66-3   | Chloroform                            | 13   | U   |
| 71-55-6   | 1,1,1-Trichloroethane                 | 13   | U   |
| 110-82-7  | Cyclohexane                           | 13   | U   |
| 56-23-5   | Carbon Tetrachloride                  | 13   | U   |
| 71-43-2   | Benzene                               | 13   | U   |
| 107-06-2  | 1,2-Dichloroethane                    | 13   | U   |

000090

EPA SAMPLE NO.

1B  
VOLATILE ORGANICS ANALYSIS DATA SHEET

BNP12

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.10

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: 315A17.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: not dec. 21

Date Analyzed: 11/10/00

GC Column: DB-624 ID: 0.53 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

|            |                             |    |   |
|------------|-----------------------------|----|---|
| 79-01-6    | Trichloroethene             | 13 | U |
| 108-87-2   | Methylcyclohexane           | 13 | U |
| 78-87-5    | 1,2-Dichloropropane         | 13 | U |
| 75-27-4    | Bromodichloromethane        | 13 | U |
| 10061-01-5 | cis-1,3-Dichloropropene     | 13 | U |
| 108-10-1   | 4-Methyl-2-Pentanone        | 13 | U |
| 108-88-3   | Toluene                     | 13 | U |
| 10061-02-6 | trans-1,3-Dichloropropene   | 13 | U |
| 79-00-5    | 1,1,2-Trichloroethane       | 13 | U |
| 127-18-4   | Tetrachloroethene           | 13 | U |
| 591-78-6   | 2-Hexanone                  | 13 | U |
| 124-48-1   | Dibromochloromethane        | 13 | U |
| 106-93-4   | 1,2-Dibromoethane           | 13 | U |
| 108-90-7   | Chlorobenzene               | 13 | U |
| 100-41-4   | Ethylbenzene                | 13 | U |
| 1330-20-7  | Xylene (total)              | 13 | U |
| 100-42-5   | Styrene                     | 13 | U |
| 75-25-2    | Bromoform                   | 13 | U |
| 98-82-8    | Isopropylbenzene            | 13 | U |
| 79-34-5    | 1,1,2,2-Tetrachloroethane   | 13 | U |
| 541-73-1   | 1,3-Dichlorobenzene         | 13 | U |
| 106-46-7   | 1,4-Dichlorobenzene         | 13 | U |
| 95-50-1    | 1,2-Dichlorobenzene         | 13 | U |
| 96-12-8    | 1,2-Dibromo-3-chloropropane | 13 | U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 13 | U |



000091

1F  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

BNP12

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) SOIL Lab Sample ID: 45798.10  
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: 315A17.D  
 Level: (low/med) LOW Date Received: 11/08/00  
 % Moisture: not dec. 21 Date Analyzed: 11/10/00  
 GC Column: DB-624 ID: 0.53 (mm) Dilution Factor: 1.0  
 Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs found: 2 CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

| CAS NUMBER  | COMPOUND NAME                | RT    | EST. CONC. | Q  |
|-------------|------------------------------|-------|------------|----|
| 1. 541-05-9 | CYCLOTRISILOXANE, HEXAMETHYL | 11.04 | 7          | JN |
| 2. 556-67-2 | CYCLOTETRAILOXANE, OCTAMETH  | 14.10 | 7          | JN |
| 3.          |                              |       |            |    |
| 4.          |                              |       |            |    |
| 5.          |                              |       |            |    |
| 6.          |                              |       |            |    |
| 7.          |                              |       |            |    |
| 8.          |                              |       |            |    |
| 9.          |                              |       |            |    |
| 10.         |                              |       |            |    |
| 11.         |                              |       |            |    |
| 12.         |                              |       |            |    |
| 13.         |                              |       |            |    |
| 14.         |                              |       |            |    |
| 15.         |                              |       |            |    |
| 16.         |                              |       |            |    |
| 17.         |                              |       |            |    |
| 18.         |                              |       |            |    |
| 19.         |                              |       |            |    |
| 20.         |                              |       |            |    |
| 21.         |                              |       |            |    |
| 22.         |                              |       |            |    |
| 23.         |                              |       |            |    |
| 24.         |                              |       |            |    |
| 25.         |                              |       |            |    |
| 26.         |                              |       |            |    |
| 27.         |                              |       |            |    |
| 28.         |                              |       |            |    |
| 29.         |                              |       |            |    |
| 30.         |                              |       |            |    |

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000098

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP13

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) SOIL Lab Sample ID: 45798.11  
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: 319A10.D  
 Level: (low/med) LOW Date Received: 11/08/00  
 % Moisture: not dec. 24 Date Analyzed: 11/14/00  
 GC Column: DB-624 ID: 0.53 (mm) Dilution Factor: 1.0  
 Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

| CAS NO.   | COMPOUND                              | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) UG/KG | Q              |
|-----------|---------------------------------------|---|----------------|
| 75-71-8   | Dichlorodifluoromethane               | 13  | U              |
| 74-87-3   | Chloromethane                         | 13  | U              |
| 75-01-4   | Vinyl Chloride                        | 13  | U              |
| 74-83-9   | Bromomethane                          | 13  | U              |
| 75-00-3   | Chloroethane                          | 13  | U              |
| 75-69-4   | Trichlorofluoromethane                | 13  | U              |
| 75-35-4   | 1,1-Dichloroethene                    | 13  | U              |
| 76-13-1   | 1,1,2-Trichloro-1,2,2-trifluoroethane | 13  | U              |
| 67-64-1   | Acetone                               | 61  | U <del>B</del> |
| 75-15-0   | Carbon Disulfide                      | 13  | U              |
| 79-20-9   | Methyl Acetate                        | 13  | U              |
| 75-09-2   | Methylene Chloride                    | 13  | U              |
| 156-60-5  | trans-1,2-Dichloroethene              | 13  | U              |
| 1634-04-4 | Methyl tert-Butyl Ether               | 13  | U              |
| 75-34-3   | 1,1-Dichloroethane                    | 13  | U              |
| 156-59-2  | cis-1,2-Dichloroethene                | 13  | U              |
| 78-93-3   | 2-Butanone                            | 13  | U              |
| 67-66-3   | Chloroform                            | 13  | U              |
| 71-55-6   | 1,1,1-Trichloroethane                 | 13  | U              |
| 110-82-7  | Cyclohexane                           | 13  | U              |
| 56-23-5   | Carbon Tetrachloride                  | 13  | U              |
| 71-43-2   | Benzene                               | 13  | U              |
| 107-06-2  | 1,2-Dichloroethane                    | 13  | U              |

000099

1B  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP13

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) SOIL Lab Sample ID: 45798.11  
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: 319A10.D  
 Level: (low/med) LOW Date Received: 11/08/00  
 % Moisture: not dec. 24 Date Analyzed: 11/14/00  
 GC Column: DB-624 ID: 0.53 (mm) Dilution Factor: 1.0  
 Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

CAS NO. COMPOUND

| CAS NO.    | COMPOUND                    | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) UG/KG | Q |
|------------|-----------------------------|---|---|
| 79-01-6    | Trichloroethene             | 13  | U |
| 108-87-2   | Methylcyclohexane           | 13  | U |
| 78-87-5    | 1,2-Dichloropropane         | 13  | U |
| 75-27-4    | Bromodichloromethane        | 13  | U |
| 10061-01-5 | cis-1,3-Dichloropropene     | 13  | U |
| 108-10-1   | 4-Methyl-2-Pentanone        | 13  | U |
| 108-88-3   | Toluene                     | 13  | U |
| 10061-02-6 | trans-1,3-Dichloropropene   | 13  | U |
| 79-00-5    | 1,1,2-Trichloroethane       | 13  | U |
| 127-18-4   | Tetrachloroethene           | 13  | U |
| 591-78-6   | 2-Hexanone                  | 13  | U |
| 124-48-1   | Dibromochloromethane        | 13  | U |
| 106-93-4   | 1,2-Dibromoethane           | 13  | U |
| 108-90-7   | Chlorobenzene               | 13  | U |
| 100-41-4   | Ethylbenzene                | 13  | U |
| 1330-20-7  | Xylene (total)              | 13  | U |
| 100-42-5   | Styrene                     | 13  | U |
| 75-25-2    | Bromoform                   | 13  | U |
| 98-82-8    | Isopropylbenzene            | 13  | U |
| 79-34-5    | 1,1,2,2-Tetrachloroethane   | 13  | U |
| 541-73-1   | 1,3-Dichlorobenzene         | 13  | U |
| 106-46-7   | 1,4-Dichlorobenzene         | 13  | U |
| 95-50-1    | 1,2-Dichlorobenzene         | 13  | U |
| 96-12-8    | 1,2-Dibromo-3-chloropropane | 13  | U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 13  | U |

1F  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

BNP13

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) SOIL Lab Sample ID: 45798.11  
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: 319A10.D  
 Level: (low/med) LOW Date Received: 11/08/00  
 % Moisture: not dec. 24 Date Analyzed: 11/14/00  
 GC Column: DB-624 ID: 0.53 (mm) Dilution Factor: 1.0  
 Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 Number TICs found: 2 CONCENTRATION UNITS:  
 (ug/L or ug/Kg) UG/KG

| CAS NUMBER  | COMPOUND NAME                | RT    | EST. CONC. | Q   |
|-------------|------------------------------|-------|------------|-----|
| 1. 556-67-2 | CYCLOTETRASILOXANE, OCTAMETH | 14.12 | 20         | BJN |
| 2.          | UNKNOWN                      | 16.76 | 19         | J   |
| 3.          |                              |       |            |     |
| 4.          |                              |       |            |     |
| 5.          |                              |       |            |     |
| 6.          |                              |       |            |     |
| 7.          |                              |       |            |     |
| 8.          |                              |       |            |     |
| 9.          |                              |       |            |     |
| 10.         |                              |       |            |     |
| 11.         |                              |       |            |     |
| 12.         |                              |       |            |     |
| 13.         |                              |       |            |     |
| 14.         |                              |       |            |     |
| 15.         |                              |       |            |     |
| 16.         |                              |       |            |     |
| 17.         |                              |       |            |     |
| 18.         |                              |       |            |     |
| 19.         |                              |       |            |     |
| 20.         |                              |       |            |     |
| 21.         |                              |       |            |     |
| 22.         |                              |       |            |     |
| 23.         |                              |       |            |     |
| 24.         |                              |       |            |     |
| 25.         |                              |       |            |     |
| 26.         |                              |       |            |     |
| 27.         |                              |       |            |     |
| 28.         |                              |       |            |     |
| 29.         |                              |       |            |     |
| 30.         |                              |       |            |     |

000107

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

|       |
|-------|
| BNQ85 |
|-------|

Lab Name: AATSLA Contract: 68-W0-0081

Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05

Matrix: (soil/water) SOIL Lab Sample ID: 45798.12

Sample wt/vol: 5.0 (g/mL) G Lab File ID: 315A19.D

Level: (low/med) LOW Date Received: 11/08/00

% Moisture: not dec. 25 ✓ Date Analyzed: 11/10/00

GC Column: DB-624 ID: 0.53 (mm) Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

| CAS NO.   | COMPOUND                              | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) UG/KG | Q   |
|-----------|---------------------------------------|---|-----|
| 75-71-8   | Dichlorodifluoromethane               | 13  | U   |
| 74-87-3   | Chloromethane                         | 13  | U   |
| 75-01-4   | Vinyl Chloride                        | 13  | U   |
| 74-83-9   | Bromomethane                          | 13  | U   |
| 75-00-3   | Chloroethane                          | 13  | U   |
| 75-69-4   | Trichlorofluoromethane                | 13  | U   |
| 75-35-4   | 1,1-Dichloroethene                    | 13  | U   |
| 76-13-1   | 1,1,2-Trichloro-1,2,2-trifluoroethane | 13  | U   |
| 67-64-1   | Acetone                               | 23  | U B |
| 75-15-0   | Carbon Disulfide                      | 13  | U   |
| 79-20-9   | Methyl Acetate                        | 13  | U   |
| 75-09-2   | Methylene Chloride                    | 13  | U   |
| 156-60-5  | trans-1,2-Dichloroethene              | 13  | U   |
| 1634-04-4 | Methyl tert-Butyl Ether               | 13  | U   |
| 75-34-3   | 1,1-Dichloroethane                    | 13  | U   |
| 156-59-2  | cis-1,2-Dichloroethene                | 13  | U   |
| 78-93-3   | 2-Butanone                            | 8   | J   |
| 67-66-3   | Chloroform                            | 13  | U   |
| 71-55-6   | 1,1,1-Trichloroethane                 | 13  | U   |
| 110-82-7  | Cyclohexane                           | 13  | U   |
| 56-23-5   | Carbon Tetrachloride                  | 13  | U   |
| 71-43-2   | Benzene                               | 13  | U   |
| 107-06-2  | 1,2-Dichloroethane                    | 13  | U   |

000108

1B  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNQ85

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) SOIL Lab Sample ID: 45798.12  
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: 315A19.D  
 Level: (low/med) LOW Date Received: 11/08/00  
 % Moisture: not dec. 25 Date Analyzed: 11/10/00  
 GC Column: DB-624 ID: 0.53 (mm) Dilution Factor: 1.0  
 Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO. COMPOUND

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

Q

|            |                             |    |   |
|------------|-----------------------------|----|---|
| 79-01-6    | Trichloroethene             | 13 | U |
| 108-87-2   | Methylcyclohexane           | 13 | U |
| 78-87-5    | 1,2-Dichloropropane         | 13 | U |
| 75-27-4    | Bromodichloromethane        | 13 | U |
| 10061-01-5 | cis-1,3-Dichloropropene     | 13 | U |
| 108-10-1   | 4-Methyl-2-Pentanone        | 13 | U |
| 108-88-3   | Toluene                     | 13 | U |
| 10061-02-6 | trans-1,3-Dichloropropene   | 13 | U |
| 79-00-5    | 1,1,2-Trichloroethane       | 13 | U |
| 127-18-4   | Tetrachloroethene           | 13 | U |
| 591-78-6   | 2-Hexanone                  | 13 | U |
| 124-48-1   | Dibromochloromethane        | 13 | U |
| 106-93-4   | 1,2-Dibromoethane           | 13 | U |
| 108-90-7   | Chlorobenzene               | 13 | U |
| 100-41-4   | Ethylbenzene                | 13 | U |
| 1330-20-7  | Xylene (total)              | 13 | U |
| 100-42-5   | Styrene                     | 13 | U |
| 75-25-2    | Bromoform                   | 13 | U |
| 98-82-8    | Isopropylbenzene            | 13 | U |
| 79-34-5    | 1,1,2,2-Tetrachloroethane   | 13 | U |
| 541-73-1   | 1,3-Dichlorobenzene         | 13 | U |
| 106-46-7   | 1,4-Dichlorobenzene         | 13 | U |
| 95-50-1    | 1,2-Dichlorobenzene         | 13 | U |
| 96-12-8    | 1,2-Dibromo-3-chloropropane | 13 | U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 13 | U |

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000109

1F  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

BNQ85

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.12

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: 315A19.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: not dec. 25

Date Analyzed: 11/10/00

GC Column: DB-624 ID: 0.53 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs found: 1

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

| CAS NUMBER  | COMPOUND NAME                | RT    | EST. CONC. | Q  |
|-------------|------------------------------|-------|------------|----|
| 1. 541-05-9 | CYCLOTRISILOXANE, HEXAMETHYL | 11.04 | 8          | JN |
| 2.          |                              |       |            |    |
| 3.          |                              |       |            |    |
| 4.          |                              |       |            |    |
| 5.          |                              |       |            |    |
| 6.          |                              |       |            |    |
| 7.          |                              |       |            |    |
| 8.          |                              |       |            |    |
| 9.          |                              |       |            |    |
| 10.         |                              |       |            |    |
| 11.         |                              |       |            |    |
| 12.         |                              |       |            |    |
| 13.         |                              |       |            |    |
| 14.         |                              |       |            |    |
| 15.         |                              |       |            |    |
| 16.         |                              |       |            |    |
| 17.         |                              |       |            |    |
| 18.         |                              |       |            |    |
| 19.         |                              |       |            |    |
| 20.         |                              |       |            |    |
| 21.         |                              |       |            |    |
| 22.         |                              |       |            |    |
| 23.         |                              |       |            |    |
| 24.         |                              |       |            |    |
| 25.         |                              |       |            |    |
| 26.         |                              |       |            |    |
| 27.         |                              |       |            |    |
| 28.         |                              |       |            |    |
| 29.         |                              |       |            |    |
| 30.         |                              |       |            |    |

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

|       |
|-------|
| BNQ87 |
|-------|

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) SOIL Lab Sample ID: 45798.13  
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: 315A20.D  
 Level: (low/med) LOW Date Received: 11/08/00  
 % Moisture: not dec. 27 ✓ Date Analyzed: 11/10/00  
 GC Column: DB-624 ID: 0.53 (mm) Dilution Factor: 1.0  
 Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

| CAS NO.   | COMPOUND                              | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) UG/KG | Q              |
|-----------|---------------------------------------|---|----------------|
| 75-71-8   | Dichlorodifluoromethane               | 14  | U              |
| 74-87-3   | Chloromethane                         | 14  | U              |
| 75-01-4   | Vinyl Chloride                        | 14  | U              |
| 74-83-9   | Bromomethane                          | 14  | U              |
| 75-00-3   | Chloroethane                          | 14  | U              |
| 75-69-4   | Trichlorofluoromethane                | 14  | U              |
| 75-35-4   | 1,1-Dichloroethene                    | 14  | U              |
| 76-13-1   | 1,1,2-Trichloro-1,2,2-trifluoroethane | 14  | U              |
| 67-64-1   | Acetone                               | 14 <del>13</del>                              | U <del>B</del> |
| 75-15-0   | Carbon Disulfide                      | 14  | U              |
| 79-20-9   | Methyl Acetate                        | 14  | U              |
| 75-09-2   | Methylene Chloride                    | 14  | U              |
| 156-60-5  | trans-1,2-Dichloroethene              | 14  | U              |
| 1634-04-4 | Methyl tert-Butyl Ether               | 14  | U              |
| 75-34-3   | 1,1-Dichloroethane                    | 14  | U              |
| 156-59-2  | cis-1,2-Dichloroethene                | 14  | U              |
| 78-93-3   | 2-Butanone                            | 5   | J              |
| 67-66-3   | Chloroform                            | 14  | U              |
| 71-55-6   | 1,1,1-Trichloroethane                 | 14  | U              |
| 110-82-7  | Cyclohexane                           | 14  | U              |
| 56-23-5   | Carbon Tetrachloride                  | 14  | U              |
| 71-43-2   | Benzene                               | 14  | U              |
| 107-06-2  | 1,2-Dichloroethane                    | 14  | U              |



000117

1B  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNQ87

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) SOIL Lab Sample ID: 45798.13  
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: 315A20.D  
 Level: (low/med) LOW Date Received: 11/08/00  
 % Moisture: not dec. 27 Date Analyzed: 11/10/00  
 GC Column: DB-624 ID: 0.53 (mm) Dilution Factor: 1.0  
 Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

| CAS NO.    | COMPOUND                    | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) UG/KG | Q |
|------------|-----------------------------|---|---|
| 79-01-6    | Trichloroethene             | 14  | U |
| 108-87-2   | Methylcyclohexane           | 14  | U |
| 78-87-5    | 1,2-Dichloropropane         | 14  | U |
| 75-27-4    | Bromodichloromethane        | 14  | U |
| 10061-01-5 | cis-1,3-Dichloropropene     | 14  | U |
| 108-10-1   | 4-Methyl-2-Pentanone        | 14  | U |
| 108-88-3   | Toluene                     | 14  | U |
| 10061-02-6 | trans-1,3-Dichloropropene   | 14  | U |
| 79-00-5    | 1,1,2-Trichloroethane       | 14  | U |
| 127-18-4   | Tetrachloroethene           | 14  | U |
| 591-78-6   | 2-Hexanone                  | 14  | U |
| 124-48-1   | Dibromochloromethane        | 14  | U |
| 106-93-4   | 1,2-Dibromoethane           | 14  | U |
| 108-90-7   | Chlorobenzene               | 14  | U |
| 100-41-4   | Ethylbenzene                | 14  | U |
| 1330-20-7  | Xylene (total)              | 14  | U |
| 100-42-5   | Styrene                     | 14  | U |
| 75-25-2    | Bromoform                   | 14  | U |
| 98-82-8    | Isopropylbenzene            | 14  | U |
| 79-34-5    | 1,1,2,2-Tetrachloroethane   | 14  | U |
| 541-73-1   | 1,3-Dichlorobenzene         | 14  | U |
| 106-46-7   | 1,4-Dichlorobenzene         | 14  | U |
| 95-50-1    | 1,2-Dichlorobenzene         | 14  | U |
| 96-12-8    | 1,2-Dibromo-3-chloropropane | 14  | U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 14  | U |

00118

1F  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

BNQ87

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.13

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: 315A20.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: not dec. 27

Date Analyzed: 11/10/00

GC Column: DB-624 ID: 0.53 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs found: 1

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

| CAS NUMBER | COMPOUND NAME | RT    | EST. CONC. | Q |
|------------|---------------|-------|------------|---|
| 1.         | UNKNOWN       | 14.08 | 8          | J |
| 2.         |               |       |            |   |
| 3.         |               |       |            |   |
| 4.         |               |       |            |   |
| 5.         |               |       |            |   |
| 6.         |               |       |            |   |
| 7.         |               |       |            |   |
| 8.         |               |       |            |   |
| 9.         |               |       |            |   |
| 10.        |               |       |            |   |
| 11.        |               |       |            |   |
| 12.        |               |       |            |   |
| 13.        |               |       |            |   |
| 14.        |               |       |            |   |
| 15.        |               |       |            |   |
| 16.        |               |       |            |   |
| 17.        |               |       |            |   |
| 18.        |               |       |            |   |
| 19.        |               |       |            |   |
| 20.        |               |       |            |   |
| 21.        |               |       |            |   |
| 22.        |               |       |            |   |
| 23.        |               |       |            |   |
| 24.        |               |       |            |   |
| 25.        |               |       |            |   |
| 26.        |               |       |            |   |
| 27.        |               |       |            |   |
| 28.        |               |       |            |   |
| 29.        |               |       |            |   |
| 30.        |               |       |            |   |

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1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNQ88

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) SOIL Lab Sample ID: 45798.14  
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: 315A21.D  
 Level: (low/med) LOW Date Received: 11/08/00  
 % Moisture: not dec. 20 Date Analyzed: 11/10/00  
 GC Column: DB-624 ID: 0.53 (mm) Dilution Factor: 1.0  
 Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

CAS NO. COMPOUND

| CAS NO.   | COMPOUND                              | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) | UG/KG | Q    |
|-----------|---------------------------------------|---|-------|------|
| 75-71-8   | Dichlorodifluoromethane               | 13                                      |       | U    |
| 74-87-3   | Chloromethane                         | 13                                      |       | U    |
| 75-01-4   | Vinyl Chloride                        | 13                                      |       | U    |
| 74-83-9   | Bromomethane                          | 13                                      |       | U    |
| 75-00-3   | Chloroethane                          | 13                                      |       | U    |
| 75-69-4   | Trichlorofluoromethane                | 13                                      |       | U    |
| 75-35-4   | 1,1-Dichloroethene                    | 13                                      |       | U    |
| 76-13-1   | 1,1,2-Trichloro-1,2,2-trifluoroethane | 13                                      |       | U    |
| 67-64-1   | Acetone                               | 13                                      | 5     | U BF |
| 75-15-0   | Carbon Disulfide                      | 13                                      |       | U    |
| 79-20-9   | Methyl Acetate                        | 13                                      |       | U    |
| 75-09-2   | Methylene Chloride                    | 13                                      | 2     | U BF |
| 156-60-5  | trans-1,2-Dichloroethene              | 13                                      |       | U    |
| 1634-04-4 | Methyl tert-Butyl Ether               | 13                                      |       | U    |
| 75-34-3   | 1,1-Dichloroethane                    | 13                                      |       | U    |
| 156-59-2  | cis-1,2-Dichloroethene                | 13                                      |       | U    |
| 78-93-3   | 2-Butanone                            | 13                                      |       | U    |
| 67-66-3   | Chloroform                            | 13                                      |       | U    |
| 71-55-6   | 1,1,1-Trichloroethane                 | 13                                      |       | U    |
| 110-82-7  | Cyclohexane                           | 13                                      |       | U    |
| 56-23-5   | Carbon Tetrachloride                  | 13                                      |       | U    |
| 71-43-2   | Benzene                               | 13                                      |       | U    |
| 107-06-2  | 1,2-Dichloroethane                    | 13                                      |       | U    |

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1B  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

|       |
|-------|
| BNQ88 |
|-------|

|  |  |
|--|--|
| Lab Name: AATSLA                       | Contract: 68-W0-0081                       |
| Lab Code: AATSLA    Case No.: 28706    | SAS No.:                    SDG No.: BNP05 |
| Matrix: (soil/water) SOIL              | Lab Sample ID: 45798.14                    |
| Sample wt/vol:            5.0 (g/mL) G | Lab File ID:    315A21.D                   |
| Level:    (low/med) LOW                | Date Received: 11/08/00                    |
| % Moisture: not dec. 20                | Date Analyzed: 11/10/00                    |
| GC Column: DB-624        ID: 0.53 (mm) | Dilution Factor: 1.0                       |
| Soil Extract Volume: _____ (uL)        | Soil Aliquot Volume: _____ (uL)            |

## CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

Q

CAS NO. COMPOUND

| CAS NO.    | COMPOUND                    | UG/KG | Q |
|------------|-----------------------------|-------|---|
| 79-01-6    | Trichloroethene             | 13    | U |
| 108-87-2   | Methylcyclohexane           | 13    | U |
| 78-87-5    | 1,2-Dichloropropane         | 13    | U |
| 75-27-4    | Bromodichloromethane        | 13    | U |
| 10061-01-5 | cis-1,3-Dichloropropene     | 13    | U |
| 108-10-1   | 4-Methyl-2-Pentanone        | 13    | U |
| 108-88-3   | Toluene                     | 13    | U |
| 10061-02-6 | trans-1,3-Dichloropropene   | 13    | U |
| 79-00-5    | 1,1,2-Trichloroethane       | 13    | U |
| 127-18-4   | Tetrachloroethene           | 13    | U |
| 591-78-6   | 2-Hexanone                  | 13    | U |
| 124-48-1   | Dibromochloromethane        | 13    | U |
| 106-93-4   | 1,2-Dibromoethane           | 13    | U |
| 108-90-7   | Chlorobenzene               | 13    | U |
| 100-41-4   | Ethylbenzene                | 13    | U |
| 1330-20-7  | Xylene (total)              | 13    | U |
| 100-42-5   | Styrene                     | 13    | U |
| 75-25-2    | Bromoform                   | 13    | U |
| 98-82-8    | Isopropylbenzene            | 13    | U |
| 79-34-5    | 1,1,2,2-Tetrachloroethane   | 13    | U |
| 541-73-1   | 1,3-Dichlorobenzene         | 13    | U |
| 106-46-7   | 1,4-Dichlorobenzene         | 13    | U |
| 95-50-1    | 1,2-Dichlorobenzene         | 13    | U |
| 96-12-8    | 1,2-Dibromo-3-chloropropane | 13    | U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 13    | U |

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1F  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

BNQ88

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) SOIL Lab Sample ID: 45798.14  
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: 315A21.D  
 Level: (low/med) LOW Date Received: 11/08/00  
 % Moisture: not dec. 20 Date Analyzed: 11/10/00  
 GC Column: DB-624 ID: 0.53 (mm) Dilution Factor: 1.0  
 Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 Number TICs found: 4 CONCENTRATION UNITS:  
 (ug/L or ug/Kg) UG/KG

| CAS NUMBER  | COMPOUND NAME                | RT    | EST. CONC. | Q  |
|-------------|------------------------------|-------|------------|----|
| 1.          | UNKNOWN                      | 11.06 | 7          | J  |
| 2. 556-67-2 | CYCLOTETRASILOXANE, OCTAMETH | 14.11 | 11         | JN |
| 3.          | UNKNOWN                      | 16.75 | 30         | J  |
| 4.          | UNKNOWN                      | 20.01 | 6          | J  |
| 5.          |                              |       |            |    |
| 6.          |                              |       |            |    |
| 7.          |                              |       |            |    |
| 8.          |                              |       |            |    |
| 9.          |                              |       |            |    |
| 10.         |                              |       |            |    |
| 11.         |                              |       |            |    |
| 12.         |                              |       |            |    |
| 13.         |                              |       |            |    |
| 14.         |                              |       |            |    |
| 15.         |                              |       |            |    |
| 16.         |                              |       |            |    |
| 17.         |                              |       |            |    |
| 18.         |                              |       |            |    |
| 19.         |                              |       |            |    |
| 20.         |                              |       |            |    |
| 21.         |                              |       |            |    |
| 22.         |                              |       |            |    |
| 23.         |                              |       |            |    |
| 24.         |                              |       |            |    |
| 25.         |                              |       |            |    |
| 26.         |                              |       |            |    |
| 27.         |                              |       |            |    |
| 28.         |                              |       |            |    |
| 29.         |                              |       |            |    |
| 30.         |                              |       |            |    |

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1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNQ89

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) SOIL Lab Sample ID: 45798.15  
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: 315A22.D  
 Level: (low/med) LOW Date Received: 11/08/00  
 % Moisture: not dec. 22 Date Analyzed: 11/10/00  
 GC Column: DB-624 ID: 0.53 (mm) Dilution Factor: 1.0  
 Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

|           |                                       |    |   |
|-----------|---------------------------------------|----|---|
| 75-71-8   | Dichlorodifluoromethane               | 13 | U |
| 74-87-3   | Chloromethane                         | 13 | U |
| 75-01-4   | Vinyl Chloride                        | 13 | U |
| 74-83-9   | Bromomethane                          | 13 | U |
| 75-00-3   | Chloroethane                          | 13 | U |
| 75-69-4   | Trichlorofluoromethane                | 13 | U |
| 75-35-4   | 1,1-Dichloroethene                    | 13 | U |
| 76-13-1   | 1,1,2-Trichloro-1,2,2-trifluoroethane | 13 | U |
| 67-64-1   | Acetone                               | 13 | U |
| 75-15-0   | Carbon Disulfide                      | 13 | U |
| 79-20-9   | Methyl Acetate                        | 13 | U |
| 75-09-2   | Methylene Chloride                    | 13 | U |
| 156-60-5  | trans-1,2-Dichloroethene              | 13 | U |
| 1634-04-4 | Methyl tert-Butyl Ether               | 13 | U |
| 75-34-3   | 1,1-Dichloroethane                    | 13 | U |
| 156-59-2  | cis-1,2-Dichloroethene                | 13 | U |
| 78-93-3   | 2-Butanone                            | 13 | U |
| 67-66-3   | Chloroform                            | 13 | U |
| 71-55-6   | 1,1,1-Trichloroethane                 | 13 | U |
| 110-82-7  | Cyclohexane                           | 13 | U |
| 56-23-5   | Carbon Tetrachloride                  | 13 | U |
| 71-43-2   | Benzene                               | 13 | U |
| 107-06-2  | 1,2-Dichloroethane                    | 13 | U |

000134

1B  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNQ89

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) SOIL Lab Sample ID: 45798.15  
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: 315A22.D  
 Level: (low/med) LOW Date Received: 11/08/00  
 % Moisture: not dec. 22 Date Analyzed: 11/10/00  
 GC Column: DB-624 ID: 0.53 (mm) Dilution Factor: 1.0  
 Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

| CAS NO.    | COMPOUND                    | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) UG/KG | Q |
|------------|-----------------------------|---|---|
| 79-01-6    | Trichloroethene             | 13  | U |
| 108-87-2   | Methylcyclohexane           | 13  | U |
| 78-87-5    | 1,2-Dichloropropane         | 13  | U |
| 75-27-4    | Bromodichloromethane        | 13  | U |
| 10061-01-5 | cis-1,3-Dichloropropene     | 13  | U |
| 108-10-1   | 4-Methyl-2-Pentanone        | 13  | U |
| 108-88-3   | Toluene                     | 13  | U |
| 10061-02-6 | trans-1,3-Dichloropropene   | 13  | U |
| 79-00-5    | 1,1,2-Trichloroethane       | 13  | U |
| 127-18-4   | Tetrachloroethene           | 13  | U |
| 591-78-6   | 2-Hexanone                  | 13  | U |
| 124-48-1   | Dibromochloromethane        | 13  | U |
| 106-93-4   | 1,2-Dibromoethane           | 13  | U |
| 108-90-7   | Chlorobenzene               | 13  | U |
| 100-41-4   | Ethylbenzene                | 13  | U |
| 1330-20-7  | Xylene (total)              | 13  | U |
| 100-42-5   | Styrene                     | 13  | U |
| 75-25-2    | Bromoform                   | 13  | U |
| 98-82-8    | Isopropylbenzene            | 13  | U |
| 79-34-5    | 1,1,2,2-Tetrachloroethane   | 13  | U |
| 541-73-1   | 1,3-Dichlorobenzene         | 13  | U |
| 106-46-7   | 1,4-Dichlorobenzene         | 13  | U |
| 95-50-1    | 1,2-Dichlorobenzene         | 13  | U |
| 96-12-8    | 1,2-Dibromo-3-chloropropane | 13  | U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 13  | U |

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000135

1F  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

BNQ89

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.15

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: 315A22.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: not dec. 22

Date Analyzed: 11/10/00

GC Column: DB-624 ID: 0.53 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs found: 1

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

| CAS NUMBER | COMPOUND NAME | RT    | EST. CONC. | Q |
|------------|---------------|-------|------------|---|
| 1.         | UNKNOWN       | 14.11 | 8          | J |
| 2.         |               |       |            |   |
| 3.         |               |       |            |   |
| 4.         |               |       |            |   |
| 5.         |               |       |            |   |
| 6.         |               |       |            |   |
| 7.         |               |       |            |   |
| 8.         |               |       |            |   |
| 9.         |               |       |            |   |
| 10.        |               |       |            |   |
| 11.        |               |       |            |   |
| 12.        |               |       |            |   |
| 13.        |               |       |            |   |
| 14.        |               |       |            |   |
| 15.        |               |       |            |   |
| 16.        |               |       |            |   |
| 17.        |               |       |            |   |
| 18.        |               |       |            |   |
| 19.        |               |       |            |   |
| 20.        |               |       |            |   |
| 21.        |               |       |            |   |
| 22.        |               |       |            |   |
| 23.        |               |       |            |   |
| 24.        |               |       |            |   |
| 25.        |               |       |            |   |
| 26.        |               |       |            |   |
| 27.        |               |       |            |   |
| 28.        |               |       |            |   |
| 29.        |               |       |            |   |
| 30.        |               |       |            |   |

2.121



000141

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNQ91

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) SOIL Lab Sample ID: 45798.16  
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: 315A23.D  
 Level: (low/med) LOW Date Received: 11/08/00  
 % Moisture: not dec. 22 Date Analyzed: 11/10/00  
 GC Column: DB-624 ID: 0.53 (mm) Dilution Factor: 1.0  
 Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NO. COMPOUND

Q

| CAS NO.   | COMPOUND                              | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) UG/KG | Q    |
|-----------|---------------------------------------|---|------|
| 75-71-8   | Dichlorodifluoromethane               | 13  | U    |
| 74-87-3   | Chloromethane                         | 13  | U    |
| 75-01-4   | Vinyl Chloride                        | 13  | U    |
| 74-83-9   | Bromomethane                          | 13  | U    |
| 75-00-3   | Chloroethane                          | 13  | U    |
| 75-69-4   | Trichlorofluoromethane                | 13  | U    |
| 75-35-4   | 1,1-Dichloroethene                    | 13  | U    |
| 76-13-1   | 1,1,2-Trichloro-1,2,2-trifluoroethane | 13  | U    |
| 67-64-1   | Acetone                               | 13 4  | U B+ |
| 75-15-0   | Carbon Disulfide                      | 13  | U    |
| 79-20-9   | Methyl Acetate                        | 13  | U    |
| 75-09-2   | Methylene Chloride                    | 13 2  | U B+ |
| 156-60-5  | trans-1,2-Dichloroethene              | 13  | U    |
| 1634-04-4 | Methyl tert-Butyl Ether               | 13  | U    |
| 75-34-3   | 1,1-Dichloroethane                    | 13  | U    |
| 156-59-2  | cis-1,2-Dichloroethene                | 13  | U    |
| 78-93-3   | 2-Butanone                            | 1   | J    |
| 67-66-3   | Chloroform                            | 13  | U    |
| 71-55-6   | 1,1,1-Trichloroethane                 | 13  | U    |
| 110-82-7  | Cyclohexane                           | 13  | U    |
| 56-23-5   | Carbon Tetrachloride                  | 13  | U    |
| 71-43-2   | Benzene                               | 13  | U    |
| 107-06-2  | 1,2-Dichloroethane                    | 13  | U    |

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1B  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNQ91

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) SOIL Lab Sample ID: 45798.16  
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: 315A23.D  
 Level: (low/med) LOW Date Received: 11/08/00  
 % Moisture: not dec. 22 Date Analyzed: 11/10/00  
 GC Column: DB-624 ID: 0.53 (mm) Dilution Factor: 1.0  
 Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

CAS NO. COMPOUND

| CAS NO.    | COMPOUND                    | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) UG/KG | Q |
|------------|-----------------------------|---|---|
| 79-01-6    | Trichloroethene             | 13  | U |
| 108-87-2   | Methylcyclohexane           | 13  | U |
| 78-87-5    | 1,2-Dichloropropane         | 13  | U |
| 75-27-4    | Bromodichloromethane        | 13  | U |
| 10061-01-5 | cis-1,3-Dichloropropene     | 13  | U |
| 108-10-1   | 4-Methyl-2-Pentanone        | 13  | U |
| 108-88-3   | Toluene                     | 13  | U |
| 10061-02-6 | trans-1,3-Dichloropropene   | 13  | U |
| 79-00-5    | 1,1,2-Trichloroethane       | 13  | U |
| 127-18-4   | Tetrachloroethene           | 13  | U |
| 591-78-6   | 2-Hexanone                  | 13  | U |
| 124-48-1   | Dibromochloromethane        | 13  | U |
| 106-93-4   | 1,2-Dibromoethane           | 13  | U |
| 108-90-7   | Chlorobenzene               | 13  | U |
| 100-41-4   | Ethylbenzene                | 13  | U |
| 1330-20-7  | Xylene (total)              | 13  | U |
| 100-42-5   | Styrene                     | 13  | U |
| 75-25-2    | Bromoform                   | 13  | U |
| 98-82-8    | Isopropylbenzene            | 13  | U |
| 79-34-5    | 1,1,2,2-Tetrachloroethane   | 13  | U |
| 541-73-1   | 1,3-Dichlorobenzene         | 13  | U |
| 106-46-7   | 1,4-Dichlorobenzene         | 13  | U |
| 95-50-1    | 1,2-Dichlorobenzene         | 13  | U |
| 96-12-8    | 1,2-Dibromo-3-chloropropane | 13  | U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 13  | U |

0.123

000143

EPA SAMPLE NO.

1F  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

BNQ91

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.16

Sample wt/vol: 5.0 (g/mL) G

Lab File ID: 315A23.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: not dec. 22

Date Analyzed: 11/10/00

GC Column: DB-624 ID: 0.53 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs found: 0

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

| CAS NUMBER | COMPOUND NAME | RT | EST. CONC. | Q |
|------------|---------------|----|------------|---|
| 1.         |               |    |            |   |
| 2.         |               |    |            |   |
| 3.         |               |    |            |   |
| 4.         |               |    |            |   |
| 5.         |               |    |            |   |
| 6.         |               |    |            |   |
| 7.         |               |    |            |   |
| 8.         |               |    |            |   |
| 9.         |               |    |            |   |
| 10.        |               |    |            |   |
| 11.        |               |    |            |   |
| 12.        |               |    |            |   |
| 13.        |               |    |            |   |
| 14.        |               |    |            |   |
| 15.        |               |    |            |   |
| 16.        |               |    |            |   |
| 17.        |               |    |            |   |
| 18.        |               |    |            |   |
| 19.        |               |    |            |   |
| 20.        |               |    |            |   |
| 21.        |               |    |            |   |
| 22.        |               |    |            |   |
| 23.        |               |    |            |   |
| 24.        |               |    |            |   |
| 25.        |               |    |            |   |
| 26.        |               |    |            |   |
| 27.        |               |    |            |   |
| 28.        |               |    |            |   |
| 29.        |               |    |            |   |
| 30.        |               |    |            |   |

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1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNQ92 FB

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) WATER Lab Sample ID: 45798.17  
 Sample wt/vol: 5 (g/mL) ML Lab File ID: 318B23.D  
 Level: (low/med) LOW Date Received: 11/08/00  
 % Moisture: not dec. \_\_\_\_\_ Date Analyzed: 11/13/00  
 GC Column: DB-624 ID: 0.53 (mm) Dilution Factor: 1.0  
 Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L Q

| CAS NO.   | COMPOUND                              | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) UG/L | Q |
|-----------|---------------------------------------|--|---|
| 75-71-8   | Dichlorodifluoromethane               | 10   | U |
| 74-87-3   | Chloromethane                         | 10   | U |
| 75-01-4   | Vinyl Chloride                        | 10   | U |
| 74-83-9   | Bromomethane                          | 10   | U |
| 75-00-3   | Chloroethane                          | 10   | U |
| 75-69-4   | Trichlorofluoromethane                | 10   | U |
| 75-35-4   | 1,1-Dichloroethene                    | 10   | U |
| 76-13-1   | 1,1,2-Trichloro-1,2,2-trifluoroethane | 10   | U |
| 67-64-1   | Acetone                               | 10   | U |
| 75-15-0   | Carbon Disulfide                      | 10   | U |
| 79-20-9   | Methyl Acetate                        | 10   | U |
| 75-09-2   | Methylene Chloride                    | 10   | U |
| 156-60-5  | trans-1,2-Dichloroethene              | 10   | U |
| 1634-04-4 | Methyl tert-Butyl Ether               | 10   | U |
| 75-34-3   | 1,1-Dichloroethane                    | 10   | U |
| 156-59-2  | cis-1,2-Dichloroethene                | 10   | U |
| 78-93-3   | 2-Butanone                            | 10   | U |
| 67-66-3   | Chloroform                            | 10   | U |
| 71-55-6   | 1,1,1-Trichloroethane                 | 10   | U |
| 110-82-7  | Cyclohexane                           | 10   | U |
| 56-23-5   | Carbon Tetrachloride                  | 10   | U |
| 71-43-2   | Benzene                               | 10   | U |
| 107-06-2  | 1,2-Dichloroethane                    | 10   | U |

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1B  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNQ92

FB

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) WATER Lab Sample ID: 45798.17  
 Sample wt/vol: 5 (g/mL) ML Lab File ID: 318B23.D  
 Level: (low/med) LOW Date Received: 11/08/00  
 % Moisture: not dec. \_\_\_\_\_ Date Analyzed: 11/13/00  
 GC Column: DB-624 ID: 0.53 (mm) Dilution Factor: 1.0  
 Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L

CAS NO. COMPOUND

Q

| CAS NO.    | COMPOUND                    | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) UG/L | Q |
|------------|-----------------------------|--|---|
| 79-01-6    | Trichloroethene             | 10   | U |
| 108-87-2   | Methylcyclohexane           | 10   | U |
| 78-87-5    | 1,2-Dichloropropane         | 10   | U |
| 75-27-4    | Bromodichloromethane        | 10   | U |
| 10061-01-5 | cis-1,3-Dichloropropene     | 10   | U |
| 108-10-1   | 4-Methyl-2-Pentanone        | 10   | U |
| 108-88-3   | Toluene                     | 10   | U |
| 10061-02-6 | trans-1,3-Dichloropropene   | 10   | U |
| 79-00-5    | 1,1,2-Trichloroethane       | 10   | U |
| 127-18-4   | Tetrachloroethene           | 10   | U |
| 591-78-6   | 2-Hexanone                  | 10   | U |
| 124-48-1   | Dibromochloromethane        | 10   | U |
| 106-93-4   | 1,2-Dibromoethane           | 10   | U |
| 108-90-7   | Chlorobenzene               | 10   | U |
| 100-41-4   | Ethylbenzene                | 10   | U |
| 1330-20-7  | Xylene (total)              | 10   | U |
| 100-42-5   | Styrene                     | 10   | U |
| 75-25-2    | Bromoform                   | 10   | U |
| 98-82-8    | Isopropylbenzene            | 10   | U |
| 79-34-5    | 1,1,2,2-Tetrachloroethane   | 10   | U |
| 541-73-1   | 1,3-Dichlorobenzene         | 10   | U |
| 106-46-7   | 1,4-Dichlorobenzene         | 10   | U |
| 95-50-1    | 1,2-Dichlorobenzene         | 10   | U |
| 96-12-8    | 1,2-Dibromo-3-chloropropane | 10   | U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 10   | U |

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1F  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

BNQ92

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) WATER Lab Sample ID: 45798.17  
 Sample wt/vol: 5 (g/mL) ML Lab File ID: 318B23.D  
 Level: (low/med) LOW Date Received: 11/08/00  
 % Moisture: not dec. \_\_\_\_\_ Date Analyzed: 11/13/00  
 GC Column: DB-624 ID: 0.53 (mm) Dilution Factor: 1.0  
 Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 Number TICs found: 0 CONCENTRATION UNITS:  
 (ug/L or ug/Kg) UG/L

| CAS NUMBER | COMPOUND NAME | RT | EST. CONC. | Q |
|------------|---------------|----|------------|---|
| 1.         |               |    |            |   |
| 2.         |               |    |            |   |
| 3.         |               |    |            |   |
| 4.         |               |    |            |   |
| 5.         |               |    |            |   |
| 6.         |               |    |            |   |
| 7.         |               |    |            |   |
| 8.         |               |    |            |   |
| 9.         |               |    |            |   |
| 10.        |               |    |            |   |
| 11.        |               |    |            |   |
| 12.        |               |    |            |   |
| 13.        |               |    |            |   |
| 14.        |               |    |            |   |
| 15.        |               |    |            |   |
| 16.        |               |    |            |   |
| 17.        |               |    |            |   |
| 18.        |               |    |            |   |
| 19.        |               |    |            |   |
| 20.        |               |    |            |   |
| 21.        |               |    |            |   |
| 22.        |               |    |            |   |
| 23.        |               |    |            |   |
| 24.        |               |    |            |   |
| 25.        |               |    |            |   |
| 26.        |               |    |            |   |
| 27.        |               |    |            |   |
| 28.        |               |    |            |   |
| 29.        |               |    |            |   |
| 30.        |               |    |            |   |

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000278

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: AATSLA

Contract: 68-W0-0081

BNP05

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: <sup>SS-01</sup>BNP05

Matrix: (soil/water) SOIL ✓

Lab Sample ID: 45798.01

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 319F08.D

Level: (low/med) LOW

Date Received: 11/08/00 ✓

% Moisture: 27 ✓ decanted: (Y/N) N

Date Extracted: 11/08/00 ✓

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/14/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y ✓

pH: 4.2 ✓

Extraction: (Type) SONC

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

Q

| CAS NO.  | COMPOUND                     |      |   |
|----------|------------------------------|------|---|
| 100-52-7 | Benzaldehyde                 | 450  | U |
| 108-95-2 | Phenol                       | 450  | U |
| 111-44-4 | bis(2-Chloroethyl) ether     | 450  | U |
| 95-57-8  | 2-Chlorophenol               | 450  | U |
| 95-48-7  | 2-Methylphenol               | 450  | U |
| 108-60-1 | 2,2'-oxybis(1-Chloropropane) | 450  | U |
| 98-86-2  | Acetophenone                 | 450  | U |
| 106-44-5 | 4-Methylphenol               | 450  | U |
| 621-64-7 | N-Nitroso-di-n-propylamine   | 450  | U |
| 67-72-1  | Hexachloroethane             | 450  | U |
| 98-95-3  | Nitrobenzene                 | 450  | U |
| 78-59-1  | Isophorone                   | 450  | U |
| 88-75-5  | 2-Nitrophenol                | 450  | U |
| 105-67-9 | 2,4-Dimethylphenol           | 450  | U |
| 111-91-1 | bis(2-Chloroethoxy)methane   | 450  | U |
| 120-83-2 | 2,4-Dichlorophenol           | 450  | U |
| 91-20-3  | Naphthalene                  | 450  | U |
| 106-47-8 | 4-Chloroaniline              | 450  | U |
| 87-68-3  | Hexachlorobutadiene          | 450  | U |
| 105-60-2 | Caprolactam                  | 450  | U |
| 59-50-7  | 4-Chloro-3-methylphenol      | 450  | U |
| 91-57-6  | 2-Methylnaphthalene          | 450  | U |
| 77-47-4  | Hexachlorocyclopentadiene    | 450  | U |
| 88-06-2  | 2,4,6-Trichlorophenol        | 450  | U |
| 95-95-4  | 2,4,5-Trichlorophenol        | 1100 | U |
| 92-52-4  | 1,1'-Biphenyl                | 450  | U |
| 91-58-7  | 2-Chloronaphthalene          | 450  | U |
| 88-74-4  | 2-Nitroaniline               | 1100 | U |
| 131-11-3 | Dimethylphthalate            | 450  | U |
| 606-20-2 | 2,6-Dinitrotoluene           | 450  | U |
| 208-96-8 | Acenaphthylene               | 450  | U |
| 99-09-2  | 3-Nitroaniline               | 1100 | U |
| 83-32-9  | Acenaphthene                 | 450  | U |

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000279

1D  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP05

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) SOIL Lab Sample ID: 45798.01  
 Sample wt/vol: 30.0 (g/mL) G Lab File ID: 319F08.D  
 Level: (low/med) LOW Date Received: 11/08/00  
 % Moisture: 27 decanted: (Y/N) N Date Extracted: 11/08/00  
 Concentrated Extract Volume: 500 (uL) Date Analyzed: 11/14/00  
 Injection Volume: 2.0 (uL) Dilution Factor: 1.0  
 GPC Cleanup: (Y/N) Y pH: 4.2 Extraction: (Type) SONC

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

| CAS NO.   | COMPOUND                   | UG/KG | Q |
|-----------|----------------------------|-------|---|
| 51-28-5   | 2,4-Dinitrophenol          | 1100  | U |
| 100-02-7  | 4-Nitrophenol              | 1100  | U |
| 132-64-9  | Dibenzofuran               | 450   | U |
| 121-14-2  | 2,4-Dinitrotoluene         | 450   | U |
| 84-66-2   | Diethylphthalate           | 450   | U |
| 86-73-7   | Fluorene                   | 450   | U |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 450   | U |
| 100-01-6  | 4-Nitroaniline             | 1100  | U |
| 534-52-1  | 4,6-Dinitro-2-methylphenol | 1100  | U |
| 86-30-6   | N-Nitrosodiphenylamine (1) | 450   | U |
| 101-55-3  | 4-Bromophenyl-phenylether  | 450   | U |
| 118-74-1  | Hexachlorobenzene          | 450   | U |
| 1912-24-9 | Atrazine                   | 450   | U |
| 87-86-5   | Pentachlorophenol          | 1100  | U |
| 85-01-8   | Phenanthrene               | 450   | U |
| 120-12-7  | Anthracene                 | 450   | U |
| 86-74-8   | Carbazole                  | 450   | U |
| 84-74-2   | Di-n-butylphthalate        | 450   | U |
| 206-44-0  | Fluoranthene               | 450   | U |
| 129-00-0  | Pyrene                     | 450   | U |
| 85-68-7   | Butylbenzylphthalate       | 450   | U |
| 91-94-1   | 3,3'-Dichlorobenzidine     | 450   | U |
| 56-55-3   | Benzo(a)anthracene         | 450   | U |
| 218-01-9  | Chrysene                   | 450   | U |
| 117-81-7  | bis(2-Ethylhexyl)phthalate | 450   | U |
| 117-84-0  | Di-n-octylphthalate        | 450   | U |
| 205-99-2  | Benzo(b)fluoranthene       | 450   | U |
| 207-08-9  | Benzo(k)fluoranthene       | 450   | U |
| 50-32-8   | Benzo(a)pyrene             | 170   | J |
| 193-39-5  | Indeno(1,2,3-cd)pyrene     | 450   | U |
| 53-70-3   | Dibenz(a,h)anthracene      | 450   | U |
| 191-24-2  | Benzo(g,h,i)perylene       | 450   | U |

(1) Cannot be seperated from Diphenylamine

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000230

1G  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

BNP05

SS-01

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.01

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 319F08.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: 27 decanted: (Y/N) N

Date Extracted: 11/08/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/14/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y

pH: 4.2

Extraction: (Type) SONC

Number TICs found: 19

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

| CAS NUMBER | COMPOUND NAME                 | RT    | EST. CONC. | Q             |
|------------|-------------------------------|-------|------------|---------------|
| 1.         | UNKNOWN                       | 3.19  | 150        | <del>BJ</del> |
| 2.         | UNKNOWN                       | 8.01  | 96         | J             |
| 3.         | UNKNOWN                       | 10.06 | 640        | <del>BJ</del> |
| 4.         | 6624-79-9<br>1-DOTRIACONTANOL | 10.38 | 200        | JN            |
| 5.         | UNKNOWN                       | 10.50 | 170        | J             |
| 6.         | 1599-67-3<br>1-DOCOSENE       | 11.26 | 150        | JN            |
| 7.         | 629-96-9<br>1-EICOSANOL       | 12.06 | 100        | JN            |
| 8.         | UNKNOWN                       | 12.87 | 95         | J             |
| 9.         | UNKNOWN                       | 13.71 | 200        | J             |
| 10.        | UNKNOWN                       | 13.91 | 370        | J             |
| 11.        | 127-22-0<br>TARAXEROL         | 14.07 | 200        | JN            |
| 12.        | UNKNOWN                       | 14.15 | 210        | J             |
| 13.        | UNKNOWN                       | 14.28 | 370        | J             |
| 14.        | UNKNOWN                       | 14.43 | 690        | J             |
| 15.        | UNKNOWN                       | 14.49 | 110        | J             |
| 16.        | UNKNOWN                       | 14.58 | 740        | J             |
| 17.        | UNKNOWN                       | 14.76 | 96         | J             |
| 18.        | UNKNOWN                       | 14.99 | 200        | J             |
| 19.        | UNKNOWN                       | 15.15 | 1600       | J             |
| 20.        |                               |       |            |               |
| 21.        |                               |       |            |               |
| 22.        |                               |       |            |               |
| 23.        |                               |       |            |               |
| 24.        |                               |       |            |               |
| 25.        |                               |       |            |               |
| 26.        |                               |       |            |               |
| 27.        |                               |       |            |               |
| 28.        |                               |       |            |               |
| 29.        |                               |       |            |               |
| 30.        |                               |       |            |               |

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP06

55-00

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.04

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 319F11.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: 28 decanted: (Y/N) N

Date Extracted: 11/08/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/14/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 5.4

Extraction: (Type) SONC  
CONCENTRATION UNITS:

CAS NO. COMPOUND

(ug/L or ug/Kg) UG/KG

Q

|          |                              |      |   |
|----------|------------------------------|------|---|
| 100-52-7 | Benzaldehyde                 | 460  | U |
| 108-95-2 | Phenol                       | 460  | U |
| 111-44-4 | bis(2-Chloroethyl) ether     | 460  | U |
| 95-57-8  | 2-Chlorophenol               | 460  | U |
| 95-48-7  | 2-Methylphenol               | 460  | U |
| 108-60-1 | 2,2'-oxybis(1-Chloropropane) | 460  | U |
| 98-86-2  | Acetophenone                 | 460  | U |
| 106-44-5 | 4-Methylphenol               | 460  | U |
| 621-64-7 | N-Nitroso-di-n-propylamine   | 460  | U |
| 67-72-1  | Hexachloroethane             | 460  | U |
| 98-95-3  | Nitrobenzene                 | 460  | U |
| 78-59-1  | Isophorone                   | 460  | U |
| 88-75-5  | 2-Nitrophenol                | 460  | U |
| 105-67-9 | 2,4-Dimethylphenol           | 460  | U |
| 111-91-1 | bis(2-Chloroethoxy)methane   | 460  | U |
| 120-83-2 | 2,4-Dichlorophenol           | 460  | U |
| 91-20-3  | Naphthalene                  | 460  | U |
| 106-47-8 | 4-Chloroaniline              | 460  | U |
| 87-68-3  | Hexachlorobutadiene          | 460  | U |
| 105-60-2 | Caprolactam                  | 460  | U |
| 59-50-7  | 4-Chloro-3-methylphenol      | 460  | U |
| 91-57-6  | 2-Methylnaphthalene          | 460  | U |
| 77-47-4  | Hexachlorocyclopentadiene    | 460  | U |
| 88-06-2  | 2,4,6-Trichlorophenol        | 460  | U |
| 95-95-4  | 2,4,5-Trichlorophenol        | 1200 | U |
| 92-52-4  | 1,1'-Biphenyl                | 460  | U |
| 91-58-7  | 2-Chloronaphthalene          | 460  | U |
| 88-74-4  | 2-Nitroaniline               | 1200 | U |
| 131-11-3 | Dimethylphthalate            | 460  | U |
| 606-20-2 | 2,6-Dinitrotoluene           | 460  | U |
| 208-96-8 | Acenaphthylene               | 460  | U |
| 99-09-2  | 3-Nitroaniline               | 1200 | U |
| 83-32-9  | Acenaphthene                 | 460  | U |

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1D  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP06

SS-02

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.04

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 319F11.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: 28 decanted: (Y/N) N

Date Extracted: 11/08/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/14/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 5.4

Extraction: (Type) SONC  
CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

| CAS NO.   | COMPOUND                   |      |   |
|-----------|----------------------------|------|---|
| 51-28-5   | 2,4-Dinitrophenol          | 1200 | U |
| 100-02-7  | 4-Nitrophenol              | 1200 | U |
| 132-64-9  | Dibenzofuran               | 460  | U |
| 121-14-2  | 2,4-Dinitrotoluene         | 460  | U |
| 84-66-2   | Diethylphthalate           | 460  | U |
| 86-73-7   | Fluorene                   | 460  | U |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 460  | U |
| 100-01-6  | 4-Nitroaniline             | 1200 | U |
| 534-52-1  | 4,6-Dinitro-2-methylphenol | 1200 | U |
| 86-30-6   | N-Nitrosodiphenylamine (1) | 460  | U |
| 101-55-3  | 4-Bromophenyl-phenylether  | 460  | U |
| 118-74-1  | Hexachlorobenzene          | 460  | U |
| 1912-24-9 | Atrazine                   | 460  | U |
| 87-86-5   | Pentachlorophenol          | 1200 | U |
| 85-01-8   | Phenanthrene               | 460  | U |
| 120-12-7  | Anthracene                 | 460  | U |
| 86-74-8   | Carbazole                  | 460  | U |
| 84-74-2   | Di-n-butylphthalate        | 460  | U |
| 206-44-0  | Fluoranthene               | 460  | U |
| 129-00-0  | Pyrene                     | 460  | U |
| 85-68-7   | Butylbenzylphthalate       | 460  | U |
| 91-94-1   | 3,3'-Dichlorobenzidine     | 460  | U |
| 56-55-3   | Benzo(a)anthracene         | 460  | U |
| 218-01-9  | Chrysene                   | 460  | U |
| 117-81-7  | bis(2-Ethylhexyl)phthalate | 60   | J |
| 117-84-0  | Di-n-octylphthalate        | 460  | U |
| 205-99-2  | Benzo(b)fluoranthene       | 460  | U |
| 207-08-9  | Benzo(k)fluoranthene       | 460  | U |
| 50-32-8   | Benzo(a)pyrene             | 230  | J |
| 193-39-5  | Indeno(1,2,3-cd)pyrene     | 460  | U |
| 53-70-3   | Dibenz(a,h)anthracene      | 460  | U |
| 191-24-2  | Benzo(g,h,i)perylene       | 460  | U |

(1) Cannot be seperated from Diphenylamine

1G  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

Lab Name: AATSLA

Contract: 68-W0-0081

BNP06  
SS-02

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.04

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 319F11.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: 28 decanted: (Y/N) N

Date Extracted: 11/08/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/14/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y

pH: 5.4

Extraction: (Type) SONC

Number TICs found: 17

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

| CAS NUMBER     | COMPOUND NAME                | RT    | EST. CONC. | Q    |
|----------------|------------------------------|-------|------------|------|
| 1. 541-02-6    | CYCLOPENTASILOXANE, DECAMETH | 3.19  | 98         | JN R |
| 2.             | UNKNOWN                      | 9.95  | 190        | J    |
| 3.             | UNKNOWN                      | 10.06 | 110        | BJ R |
| 4. 6624-79-9   | 1-DOTRIACONTANOL             | 10.39 | 290        | JN   |
| 5.             | UNKNOWN                      | 11.26 | 230        | J    |
| 6. 629-96-9    | 1-EICOSANOL                  | 12.07 | 96         | JN   |
| 7.             | UNKNOWN                      | 13.91 | 200        | J    |
| 8. 514-07-8    | D-FRIEDOOLEAN-14-EN-3-ONE    | 13.94 | 300        | JN   |
| 9. 127-22-0    | TARAXEROL                    | 14.07 | 200        | JN   |
| 10.            | UNKNOWN                      | 14.16 | 130        | J    |
| 11. 22611-26-3 | D:C-FRIEDOOLEAN-8-EN-3-ONE   | 14.25 | 180        | JN   |
| 12. 58801-23-3 | HOP-22(29)-EN-3.BETA.-OL     | 14.37 | 260        | JN   |
| 13.            | UNKNOWN                      | 14.42 | 180        | J    |
| 14.            | UNKNOWN                      | 14.49 | 100        | J    |
| 15.            | UNKNOWN                      | 14.58 | 400        | J    |
| 16.            | UNKNOWN                      | 14.99 | 220        | J    |
| 17.            | UNKNOWN                      | 15.16 | 1900       | J    |
| 18.            |                              |       |            |      |
| 19.            |                              |       |            |      |
| 20.            |                              |       |            |      |
| 21.            |                              |       |            |      |
| 22.            |                              |       |            |      |
| 23.            |                              |       |            |      |
| 24.            |                              |       |            |      |
| 25.            |                              |       |            |      |
| 26.            |                              |       |            |      |
| 27.            |                              |       |            |      |
| 28.            |                              |       |            |      |
| 29.            |                              |       |            |      |
| 30.            |                              |       |            |      |

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000330

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP07

SS-03

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.05

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 319F12.D

Level: (low/med) LOW

Date Received: 11/08/00 ✓

% Moisture: 21 ✓ decanted: (Y/N) N ✓

Date Extracted: 11/08/00 ✓

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/14/00

Injection Volume: 2.0 (uL) ✓

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y ✓

pH: 4.7 ✓

Extraction: (Type) SONC  
CONCENTRATION UNITS:

CAS NO. COMPOUND

(ug/L or ug/Kg) UG/KG

Q

| CAS NO.  | COMPOUND                     | UG/KG | Q   |
|----------|------------------------------|-------|-----|
| 100-52-7 | Benzaldehyde                 | 420   | U J |
| 108-95-2 | Phenol                       | 420   | U   |
| 111-44-4 | bis(2-Chloroethyl) ether     | 420   | U   |
| 95-57-8  | 2-Chlorophenol               | 420   | U   |
| 95-48-7  | 2-Methylphenol               | 420   | U   |
| 108-60-1 | 2,2'-oxybis(1-Chloropropane) | 420   | U   |
| 98-86-2  | Acetophenone                 | 420   | U   |
| 106-44-5 | 4-Methylphenol               | 420   | U   |
| 621-64-7 | N-Nitroso-di-n-propylamine   | 420   | U   |
| 67-72-1  | Hexachloroethane             | 420   | U   |
| 98-95-3  | Nitrobenzene                 | 420   | U   |
| 78-59-1  | Isophorone                   | 420   | U   |
| 88-75-5  | 2-Nitrophenol                | 420   | U   |
| 105-67-9 | 2,4-Dimethylphenol           | 420   | U   |
| 111-91-1 | bis(2-Chloroethoxy)methane   | 420   | U   |
| 120-83-2 | 2,4-Dichlorophenol           | 420   | U   |
| 91-20-3  | Naphthalene                  | 420   | U   |
| 106-47-8 | 4-Chloroaniline              | 420   | U   |
| 87-68-3  | Hexachlorobutadiene          | 420   | U   |
| 105-60-2 | Caprolactam                  | 420   | U   |
| 59-50-7  | 4-Chloro-3-methylphenol      | 420   | U   |
| 91-57-6  | 2-Methylnaphthalene          | 420   | U   |
| 77-47-4  | Hexachlorocyclopentadiene    | 420   | U   |
| 88-06-2  | 2,4,6-Trichlorophenol        | 420   | U   |
| 95-95-4  | 2,4,5-Trichlorophenol        | 1100  | U   |
| 92-52-4  | 1,1'-Biphenyl                | 420   | U   |
| 91-58-7  | 2-Chloronaphthalene          | 420   | U   |
| 88-74-4  | 2-Nitroaniline               | 1100  | U   |
| 131-11-3 | Dimethylphthalate            | 420   | U   |
| 606-20-2 | 2,6-Dinitrotoluene           | 420   | U   |
| 208-96-8 | Acenaphthylene               | 420   | U   |
| 99-09-2  | 3-Nitroaniline               | 1100  | U   |
| 83-32-9  | Acenaphthene                 | 420   | U   |

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1D  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

BNP07

SS-07

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.05

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 319F12.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: 21 decanted: (Y/N) N

Date Extracted: 11/08/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/14/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y

pH: 4.7

Extraction: (Type) SONC  
CONCENTRATION UNITS:

CAS NO. COMPOUND

(ug/L or ug/Kg) UG/KG

Q

|           |                            |      |   |
|-----------|----------------------------|------|---|
| 51-28-5   | 2,4-Dinitrophenol          | 1100 | U |
| 100-02-7  | 4-Nitrophenol              | 1100 | U |
| 132-64-9  | Dibenzofuran               | 420  | U |
| 121-14-2  | 2,4-Dinitrotoluene         | 420  | U |
| 84-66-2   | Diethylphthalate           | 420  | U |
| 86-73-7   | Fluorene                   | 420  | U |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 420  | U |
| 100-01-6  | 4-Nitroaniline             | 1100 | U |
| 534-52-1  | 4,6-Dinitro-2-methylphenol | 1100 | U |
| 86-30-6   | N-Nitrosodiphenylamine (1) | 420  | U |
| 101-55-3  | 4-Bromophenyl-phenylether  | 420  | U |
| 118-74-1  | Hexachlorobenzene          | 420  | U |
| 1912-24-9 | Atrazine                   | 420  | U |
| 87-86-5   | Pentachlorophenol          | 1100 | U |
| 85-01-8   | Phenanthrene               | 420  | U |
| 120-12-7  | Anthracene                 | 420  | U |
| 86-74-8   | Carbazole                  | 420  | U |
| 84-74-2   | Di-n-butylphthalate        | 420  | U |
| 206-44-0  | Fluoranthene               | 420  | U |
| 129-00-0  | Pyrene                     | 420  | U |
| 85-68-7   | Butylbenzylphthalate       | 420  | U |
| 91-94-1   | 3,3'-Dichlorobenzidine     | 420  | U |
| 56-55-3   | Benzo (a) anthracene       | 420  | U |
| 218-01-9  | Chrysene                   | 420  | U |
| 117-81-7  | bis(2-Ethylhexyl)phthalate | 380  | J |
| 117-84-0  | Di-n-octylphthalate        | 420  | U |
| 205-99-2  | Benzo (b) fluoranthene     | 420  | U |
| 207-08-9  | Benzo (k) fluoranthene     | 420  | U |
| 50-32-8   | Benzo (a) pyrene           | 420  | U |
| 193-39-5  | Indeno (1,2,3-cd) pyrene   | 420  | U |
| 53-70-3   | Dibenz (a,h) anthracene    | 420  | U |
| 191-24-2  | Benzo (g,h,i) perylene     | 420  | U |

(1) Cannot be seperated from Diphenylamine

000332

1G  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

BNP07

SS-03

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.05

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 319F12.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: 21 decanted: (Y/N) N

Date Extracted: 11/08/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/14/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y

pH: 4.7

Extraction: (Type) SONC

Number TICs found: 6

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

| CAS NUMBER  | COMPOUND NAME                | RT    | EST. CONC.    | Q  |
|-------------|------------------------------|-------|---------------|----|
| 1.          | UNKNOWN                      | 2.37  | 120           | J  |
| 2. 541-02-6 | CYCLOPENTASILOXANE, DECAMETH | 3.19  | <del>89</del> | JN |
| 3.          | UNKNOWN                      | 9.84  | 99            | J  |
| 4.          | UNKNOWN                      | 10.06 | 85            | BU |
| 5.          | UNKNOWN                      | 11.10 | 200           | J  |
| 6.          | UNKNOWN                      | 12.23 | 310           | J  |
| 7.          |                              |       |               |    |
| 8.          |                              |       |               |    |
| 9.          |                              |       |               |    |
| 10.         |                              |       |               |    |
| 11.         |                              |       |               |    |
| 12.         |                              |       |               |    |
| 13.         |                              |       |               |    |
| 14.         |                              |       |               |    |
| 15.         |                              |       |               |    |
| 16.         |                              |       |               |    |
| 17.         |                              |       |               |    |
| 18.         |                              |       |               |    |
| 19.         |                              |       |               |    |
| 20.         |                              |       |               |    |
| 21.         |                              |       |               |    |
| 22.         |                              |       |               |    |
| 23.         |                              |       |               |    |
| 24.         |                              |       |               |    |
| 25.         |                              |       |               |    |
| 26.         |                              |       |               |    |
| 27.         |                              |       |               |    |
| 28.         |                              |       |               |    |
| 29.         |                              |       |               |    |
| 30.         |                              |       |               |    |

R  
R

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1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: AATSLA

Contract: 68-W0-0081

BNP08

Lab Code: AATSLA Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL ✓

Lab Sample ID: 45798.06

Sample wt/vol: 30.0 (g/mL) G ✓

Lab File ID: 319F13.D

Level: (low/med) LOW

Date Received: 11/08/00 ✓

% Moisture: 3 ✓ decanted: (Y/N) N

Date Extracted: 11/08/00 ✓

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/14/00

Injection Volume: 2.0 (uL) ✓

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y ✓ pH: 6.0 ✓

Extraction: (Type) SONC  
CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

| CAS NO.  | COMPOUND                     |                         |   |
|----------|------------------------------|-------------------------|---|
| 100-52-7 | Benzaldehyde                 | 340                     | U |
| 108-95-2 | Phenol                       | 7900 * <del>16000</del> | U |
| 111-44-4 | bis(2-Chloroethyl) ether     | 340                     | U |
| 95-57-8  | 2-Chlorophenol               | 340                     | U |
| 95-48-7  | 2-Methylphenol               | 340                     | U |
| 108-60-1 | 2,2'-oxybis(1-Chloropropane) | 340                     | U |
| 98-86-2  | Acetophenone                 | 830                     | U |
| 106-44-5 | 4-Methylphenol               | 340                     | U |
| 621-64-7 | N-Nitroso-di-n-propylamine   | 340                     | U |
| 67-72-1  | Hexachloroethane             | 340                     | U |
| 98-95-3  | Nitrobenzene                 | 340                     | U |
| 78-59-1  | Isophorone                   | 340                     | U |
| 88-75-5  | 2-Nitrophenol                | 340                     | U |
| 105-67-9 | 2,4-Dimethylphenol           | 340                     | U |
| 111-91-1 | bis(2-Chloroethoxy) methane  | 340                     | U |
| 120-83-2 | 2,4-Dichlorophenol           | 340                     | U |
| 91-20-3  | Naphthalene                  | 1100                    | U |
| 106-47-8 | 4-Chloroaniline              | 340                     | U |
| 87-68-3  | Hexachlorobutadiene          | 340                     | U |
| 105-60-2 | Caprolactam                  | 340                     | U |
| 59-50-7  | 4-Chloro-3-methylphenol      | 340                     | U |
| 91-57-6  | 2-Methylnaphthalene          | 130                     | J |
| 77-47-4  | Hexachlorocyclopentadiene    | 340                     | U |
| 88-06-2  | 2,4,6-Trichlorophenol        | 340                     | U |
| 95-95-4  | 2,4,5-Trichlorophenol        | 860                     | U |
| 92-52-4  | 1,1'-Biphenyl                | 340                     | U |
| 91-58-7  | 2-Chloronaphthalene          | 340                     | U |
| 88-74-4  | 2-Nitroaniline               | 860                     | U |
| 131-11-3 | Dimethylphthalate            | 340                     | U |
| 606-20-2 | 2,6-Dinitrotoluene           | 340                     | U |
| 208-96-8 | Acenaphthylene               | 340                     | U |
| 99-09-2  | 3-Nitroaniline               | 860                     | U |
| 83-32-9  | Acenaphthene                 | 340                     | U |

\* Value transferred from dilution run.



000343

EPA SAMPLE NO.

1D  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

BNP08

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.06

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 319F13.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: 3 decanted: (Y/N) N

Date Extracted: 11/08/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/14/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.0

Extraction: (Type) SONC  
CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

CAS NO. COMPOUND

| CAS NO.   | COMPOUND                   | UG/KG | Q |
|-----------|----------------------------|-------|---|
| 51-28-5   | 2,4-Dinitrophenol          | 860   | U |
| 100-02-7  | 4-Nitrophenol              | 860   | U |
| 132-64-9  | Dibenzofuran               | 340   | U |
| 121-14-2  | 2,4-Dinitrotoluene         | 340   | U |
| 84-66-2   | Diethylphthalate           | 340   | U |
| 86-73-7   | Fluorene                   | 340   | U |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 340   | U |
| 100-01-6  | 4-Nitroaniline             | 860   | U |
| 534-52-1  | 4,6-Dinitro-2-methylphenol | 860   | U |
| 86-30-6   | N-Nitrosodiphenylamine (1) | 340   | U |
| 101-55-3  | 4-Bromophenyl-phenylether  | 340   | U |
| 118-74-1  | Hexachlorobenzene          | 340   | U |
| 1912-24-9 | Atrazine                   | 340   | U |
| 87-86-5   | Pentachlorophenol          | 860   | U |
| 85-01-8   | Phenanthrene               | 340   | U |
| 120-12-7  | Anthracene                 | 340   | U |
| 86-74-8   | Carbazole                  | 340   | U |
| 84-74-2   | Di-n-butylphthalate        | 58    | J |
| 206-44-0  | Fluoranthene               | 67    | J |
| 129-00-0  | Pyrene                     | 59    | J |
| 85-68-7   | Butylbenzylphthalate       | 340   | U |
| 91-94-1   | 3,3'-Dichlorobenzidine     | 340   | U |
| 56-55-3   | Benzo(a)anthracene         | 340   | U |
| 218-01-9  | Chrysene                   | 41    | J |
| 117-81-7  | bis(2-Ethylhexyl)phthalate | 520   |   |
| 117-84-0  | Di-n-octylphthalate        | 340   | U |
| 205-99-2  | Benzo(b)fluoranthene       | 340   | U |
| 207-08-9  | Benzo(k)fluoranthene       | 340   | U |
| 50-32-8   | Benzo(a)pyrene             | 340   | U |
| 193-39-5  | Indeno(1,2,3-cd)pyrene     | 340   | U |
| 53-70-3   | Dibenz(a,h)anthracene      | 340   | U |
| 191-24-2  | Benzo(g,h,i)perylene       | 340   | U |

(1) Cannot be seperated from Diphenylamine

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000344

1G  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

BNP08

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

SS-04A

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.06

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 319F13.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: 3 decanted: (Y/N) N

Date Extracted: 11/08/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/14/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y

pH: 6.0

Extraction: (Type) SONC

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

Number TICs found: 24

| CAS NUMBER    | COMPOUND NAME                | RT    | EST. CONC. | Q  |
|---------------|------------------------------|-------|------------|----|
| 1. 103-65-1   | BENZENE, PROPYL-             | 2.41  | 240        | JN |
| 2. 622-96-8   | BENZENE, 1-ETHYL-4-METHYL-   | 2.45  | 1600       | JN |
| 3. 95-36-3    | 1,2,4-TRIMETHYLBENZENE       | 2.48  | 3500       | JN |
| 4.            | UNKNOWN                      | 3.00  | 12000      | J  |
| 5. 2870-04-4  | BENZENE, 2-ETHYL-1,3-DIMETHY | 3.12  | 100        | JN |
| 6. 934-80-5   | BENZENE, 4-ETHYL-1,2-DIMETHY | 3.16  | 280        | JN |
| 7. 95-93-2    | BENZENE, 1,2,4,5-TETRAMETHYL | 3.18  | 350        | JN |
| 8.            | UNKNOWN                      | 3.20  | 170        | BJ |
| 9.            | UNKNOWN                      | 3.24  | 85         | BJ |
| 10. 933-98-2  | BENZENE, 1-ETHYL-2,3-DIMETHY | 3.26  | 450        | JN |
| 11.           | UNKNOWN                      | 3.30  | 73         | J  |
| 12.           | UNKNOWN                      | 3.35  | 660        | J  |
| 13. 5161-04-6 | BENZENE, 1-METHYL-4-(2-METHY | 3.39  | 180        | JN |
| 14. 2049-95-8 | BENZENE, (1,1-DIMETHYLPROPYL | 3.42  | 120        | JN |
| 15. 627-93-0  | HEXANEDIOIC ACID, DIMETHYL E | 3.71  | 270        | JN |
| 16.           | UNKNOWN                      | 3.96  | 690        | J  |
| 17.           | UNKNOWN                      | 4.64  | 220        | J  |
| 18.           | UNKNOWN                      | 6.81  | 950        | J  |
| 19. 84-69-5   | 1,2-BENZENEDICARBOXYLIC ACID | 7.37  | 4900       | JN |
| 20.           | UNKNOWN                      | 7.68  | 150        | J  |
| 21.           | UNKNOWN                      | 8.28  | 260        | J  |
| 22.           | UNKNOWN                      | 8.42  | 280        | J  |
| 23. 101-68-8  | BENZENE, 1,1'-METHYLENEBIS[4 | 8.68  | 1600       | JN |
| 24. 1058-61-3 | STIGMAST-4-EN-3-ONE          | 14.57 | 130        | JN |
| 25.           |                              |       |            |    |
| 26.           |                              |       |            |    |
| 27.           |                              |       |            |    |
| 28.           |                              |       |            |    |
| 29.           |                              |       |            |    |
| 30.           |                              |       |            |    |

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000396

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: AATSLA

Contract: 68-W0-0081

BNP09

Lab Code: AATSLA Case No.: 28706

SAS No.:

SDG No.: BNP05

SS 04B

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.07

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 319F14.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: 10 decanted: (Y/N) N

Date Extracted: 11/08/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/14/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 9.2

Extraction: (Type) SONC  
CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

CAS NO. COMPOUND

| CAS NO.  | COMPOUND                     | UG/KG                  | Q            |
|----------|------------------------------|------------------------|--------------|
| 100-52-7 | Benzaldehyde                 | 370                    | U J          |
| 108-95-2 | Phenol                       | <del>3500 # 5900</del> | <del>E</del> |
| 111-44-4 | bis(2-Chloroethyl) ether     | 370                    | U            |
| 95-57-8  | 2-Chlorophenol               | 370                    | U            |
| 95-48-7  | 2-Methylphenol               | 370                    | U            |
| 108-60-1 | 2,2'-oxybis(1-Chloropropane) | 370                    | U            |
| 98-86-2  | Acetophenone                 | 370                    | U            |
| 106-44-5 | 4-Methylphenol               | 370                    | U            |
| 621-64-7 | N-Nitroso-di-n-propylamine   | 370                    | U            |
| 67-72-1  | Hexachloroethane             | 370                    | U            |
| 98-95-3  | Nitrobenzene                 | 370                    | U            |
| 78-59-1  | Isophorone                   | 370                    | U            |
| 88-75-5  | 2-Nitrophenol                | 370                    | U            |
| 105-67-9 | 2,4-Dimethylphenol           | 370                    | U            |
| 111-91-1 | bis(2-Chloroethoxy)methane   | 370                    | U            |
| 120-83-2 | 2,4-Dichlorophenol           | 370                    | U            |
| 91-20-3  | Naphthalene                  | 380                    | U            |
| 106-47-8 | 4-Chloroaniline              | 370                    | U            |
| 87-68-3  | Hexachlorobutadiene          | 370                    | U            |
| 105-60-2 | Caprolactam                  | 370                    | U            |
| 59-50-7  | 4-Chloro-3-methylphenol      | 370                    | U            |
| 91-57-6  | 2-Methylnaphthalene          | 150                    | J            |
| 77-47-4  | Hexachlorocyclopentadiene    | 370                    | U            |
| 88-06-2  | 2,4,6-Trichlorophenol        | 370                    | U            |
| 95-95-4  | 2,4,5-Trichlorophenol        | 920                    | U            |
| 92-52-4  | 1,1'-Biphenyl                | 370                    | U            |
| 91-58-7  | 2-Chloronaphthalene          | 370                    | U            |
| 88-74-4  | 2-Nitroaniline               | 920                    | U            |
| 131-11-3 | Dimethylphthalate            | 370                    | U            |
| 606-20-2 | 2,6-Dinitrotoluene           | 370                    | U            |
| 208-96-8 | Acenaphthylene               | 370                    | U            |
| 99-09-2  | 3-Nitroaniline               | 920                    | U            |
| 83-32-9  | Acenaphthene                 | 300                    | J            |

\* Values transferred from dilution runs.

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000397

1D  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP09

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.07

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 319F14.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: 10 decanted: (Y/N) N

Date Extracted: 11/08/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/14/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 9.2

Extraction: (Type) SONC  
CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

| CAS NO.   | COMPOUND                   |                        |   |
|-----------|----------------------------|------------------------|---|
| 51-28-5   | 2,4-Dinitrophenol          | 920                    | U |
| 100-02-7  | 4-Nitrophenol              | 920                    | U |
| 132-64-9  | Dibenzofuran               | 170                    | J |
| 121-14-2  | 2,4-Dinitrotoluene         | 370                    | U |
| 84-66-2   | Diethylphthalate           | 370                    | U |
| 86-73-7   | Fluorene                   | 260                    | J |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 370                    | U |
| 100-01-6  | 4-Nitroaniline             | 920                    | U |
| 534-52-1  | 4,6-Dinitro-2-methylphenol | 920                    | U |
| 86-30-6   | N-Nitrosodiphenylamine (1) | 370                    | U |
| 101-55-3  | 4-Bromophenyl-phenylether  | 370                    | U |
| 118-74-1  | Hexachlorobenzene          | 370                    | U |
| 1912-24-9 | Atrazine                   | 370                    | U |
| 87-86-5   | Pentachlorophenol          | 920                    | U |
| 85-01-8   | Phenanthrene               | <del>3700</del> * 5300 | H |
| 120-12-7  | Anthracene                 | 600                    |   |
| 86-74-8   | Carbazole                  | 370                    | U |
| 84-74-2   | Di-n-butylphthalate        | 61                     | J |
| 206-44-0  | Fluoranthene               | <del>4600</del> * 6900 | H |
| 129-00-0  | Pyrene                     | <del>4000</del> * 4800 | H |
| 85-68-7   | Butylbenzylphthalate       | 370                    | U |
| 91-94-1   | 3,3'-Dichlorobenzidine     | 370                    | U |
| 56-55-3   | Benzo(a)anthracene         | 2400                   |   |
| 218-01-9  | Chrysene                   | 2800                   |   |
| 117-81-7  | bis(2-Ethylhexyl)phthalate | 870                    |   |
| 117-84-0  | Di-n-octylphthalate        | 370                    | U |
| 205-99-2  | Benzo(b)fluoranthene       | <del>2100</del> * 3400 | H |
| 207-08-9  | Benzo(k)fluoranthene       | 760                    |   |
| 50-32-8   | Benzo(a)pyrene             | 2100                   |   |
| 193-39-5  | Indeno(1,2,3-cd)pyrene     | 1300                   |   |
| 53-70-3   | Dibenz(a,h)anthracene      | 530                    |   |
| 191-24-2  | Benzo(g,h,i)perylene       | 1300                   |   |

(1) Cannot be separated from Diphenylamine

\* Values transferred from dilution num.

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1G  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

BNP09

Lab Name: AATSLA

Contract: 68-W0-0081

SS-040

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.07

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 319F14.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: 10 decanted: (Y/N) N

Date Extracted: 11/08/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/14/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y

pH: 9.2

Extraction: (Type) SONC

Number TICs found: 23

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

| CAS NUMBER    | COMPOUND NAME                    | RT    | EST. CONC. | Q    |
|---------------|----------------------------------|-------|------------|------|
| 1. 103-65-1   | BENZENE, PROPYL-                 | 2.41  | 160        | JN   |
| 2. 620-14-4   | BENZENE, 1-ETHYL-3-METHYL-       | 2.45  | 790        | JN   |
| 3. 526-73-8   | BENZENE, 1,2,3-TRIMETHYL-        | 2.47  | 1400       | JN   |
| 4. 933-98-2   | BENZENE, 1-ETHYL-2,3-DIMETHY     | 2.88  | 950        | JN   |
| 5.            | UNKNOWN                          | 2.98  | 4000       | J    |
| 6. 933-98-2   | BENZENE, 1-ETHYL-2,3-DIMETHY     | 3.11  | 80         | JN   |
| 7. 95-93-2    | BENZENE, 1,2,4,5-TETRAMETHYL     | 3.15  | 170        | JN   |
| 8. 934-74-7   | BENZENE, 1-ETHYL-3,5-DIMETHY     | 3.18  | 250        | JN   |
| 9.            | UNKNOWN                          | 3.20  | 200        | BJ R |
| 10. 2050-24-0 | BENZENE, 1,3-DIETHYL-5-METHY     | 3.26  | 270        | JN   |
| 11. 1005-64-7 | (E)-1-PHENYL-1-BUTENE            | 3.34  | 370        | JN   |
| 12.           | UNKNOWN                          | 3.39  | 100        | J    |
| 13.           | UNKNOWN                          | 3.41  | 87         | J    |
| 14. 627-93-0  | HEXANEDIOIC ACID, DIMETHYL E     | 3.70  | 150        | JN   |
| 15.           | UNKNOWN                          | 3.94  | 330        | J    |
| 16.           | UNKNOWN                          | 6.82  | 77         | J    |
| 17.           | UNKNOWN                          | 7.82  | 82         | J    |
| 18. 101-68-8  | BENZENE, 1,1'-METHYLENEBIS [4    | 8.70  | 170        | JN   |
| 19.           | UNKNOWN                          | 10.10 | 100        | BJ   |
| 20. 239-35-0  | BENZO [B] NAPHTHO [2,1-D] THIOPH | 10.24 | 85         | JN   |
| 21.           | UNKNOWN                          | 10.39 | 240        | J    |
| 22. 205-99-2  | BENZ [E] ACEPHENANTHRYLENE       | 12.09 | 140        | JN   |
| 23. 207-08-9  | BENZO [K] FLUORANTHENE           | 12.30 | 340        | JN R |
| 24.           |                                  |       |            |      |
| 25.           |                                  |       |            |      |
| 26.           |                                  |       |            |      |
| 27.           |                                  |       |            |      |
| 28.           |                                  |       |            |      |
| 29.           |                                  |       |            |      |
| 30.           |                                  |       |            |      |

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1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP10

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.08

Sample wt/vol: 30.0 (g/mL) G ✓

Lab File ID: 320F04.D

Level: (low/med) LOW ✓

Date Received: 11/08/00 ✓

% Moisture: 19 ✓ decanted: (Y/N) N

Date Extracted: 11/08/00 ✓

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/15/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y ✓ pH: 8.3 ✓

Extraction: (Type) SONC  
CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

CAS NO. COMPOUND

| CAS NO.  | COMPOUND                     | UG/KG | Q |
|----------|------------------------------|-------|---|
| 100-52-7 | Benzaldehyde                 | 410   | U |
| 108-95-2 | Phenol                       | 410   | U |
| 111-44-4 | bis(2-Chloroethyl) ether     | 410   | U |
| 95-57-8  | 2-Chlorophenol               | 410   | U |
| 95-48-7  | 2-Methylphenol               | 410   | U |
| 108-60-1 | 2,2'-oxybis(1-Chloropropane) | 410   | U |
| 98-86-2  | Acetophenone                 | 410   | U |
| 106-44-5 | 4-Methylphenol               | 410   | U |
| 621-64-7 | N-Nitroso-di-n-propylamine   | 410   | U |
| 67-72-1  | Hexachloroethane             | 410   | U |
| 98-95-3  | Nitrobenzene                 | 410   | U |
| 78-59-1  | Isophorone                   | 410   | U |
| 88-75-5  | 2-Nitrophenol                | 410   | U |
| 105-67-9 | 2,4-Dimethylphenol           | 410   | U |
| 111-91-1 | bis(2-Chloroethoxy) methane  | 410   | U |
| 120-83-2 | 2,4-Dichlorophenol           | 410   | U |
| 91-20-3  | Naphthalene                  | 410   | U |
| 106-47-8 | 4-Chloroaniline              | 410   | U |
| 87-68-3  | Hexachlorobutadiene          | 410   | U |
| 105-60-2 | Caprolactam                  | 410   | U |
| 59-50-7  | 4-Chloro-3-methylphenol      | 410   | U |
| 91-57-6  | 2-Methylnaphthalene          | 410   | U |
| 77-47-4  | Hexachlorocyclopentadiene    | 410   | U |
| 88-06-2  | 2,4,6-Trichlorophenol        | 410   | U |
| 95-95-4  | 2,4,5-Trichlorophenol        | 1000  | U |
| 92-52-4  | 1,1'-Biphenyl                | 410   | U |
| 91-58-7  | 2-Chloronaphthalene          | 410   | U |
| 88-74-4  | 2-Nitroaniline               | 1000  | U |
| 131-11-3 | Dimethylphthalate            | 410   | U |
| 606-20-2 | 2,6-Dinitrotoluene           | 410   | U |
| 208-96-8 | Acenaphthylene               | 410   | U |
| 99-09-2  | 3-Nitroaniline               | 1000  | U |
| 83-32-9  | Acenaphthene                 | 410   | U |

1D  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: AATSLA

Contract: 68-W0-0081

BNP10

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.08

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 320F04.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: 19 decanted: (Y/N) N

Date Extracted: 11/08/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/15/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 8.3

Extraction: (Type) SONC  
CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

CAS NO. COMPOUND

| CAS NO.   | COMPOUND                   | UG/KG | Q |
|-----------|----------------------------|-------|---|
| 51-28-5   | 2,4-Dinitrophenol          | 1000  | U |
| 100-02-7  | 4-Nitrophenol              | 1000  | U |
| 132-64-9  | Dibenzofuran               | 410   | U |
| 121-14-2  | 2,4-Dinitrotoluene         | 410   | U |
| 84-66-2   | Diethylphthalate           | 410   | U |
| 86-73-7   | Fluorene                   | 410   | U |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 410   | U |
| 100-01-6  | 4-Nitroaniline             | 1000  | U |
| 534-52-1  | 4,6-Dinitro-2-methylphenol | 1000  | U |
| 86-30-6   | N-Nitrosodiphenylamine (1) | 410   | U |
| 101-55-3  | 4-Bromophenyl-phenylether  | 410   | U |
| 118-74-1  | Hexachlorobenzene          | 410   | U |
| 1912-24-9 | Atrazine                   | 410   | U |
| 87-86-5   | Pentachlorophenol          | 1000  | U |
| 85-01-8   | Phenanthrene               | 410   | U |
| 120-12-7  | Anthracene                 | 410   | U |
| 86-74-8   | Carbazole                  | 410   | U |
| 84-74-2   | Di-n-butylphthalate        | 410   | U |
| 206-44-0  | Fluoranthene               | 410   | U |
| 129-00-0  | Pyrene                     | 410   | U |
| 85-68-7   | Butylbenzylphthalate       | 410   | U |
| 91-94-1   | 3,3'-Dichlorobenzidine     | 410   | U |
| 56-55-3   | Benzo(a)anthracene         | 410   | U |
| 218-01-9  | Chrysene                   | 410   | U |
| 117-81-7  | bis(2-Ethylhexyl)phthalate | 94    | J |
| 117-84-0  | Di-n-octylphthalate        | 410   | U |
| 205-99-2  | Benzo(b)fluoranthene       | 410   | U |
| 207-08-9  | Benzo(k)fluoranthene       | 410   | U |
| 50-32-8   | Benzo(a)pyrene             | 410   | U |
| 193-39-5  | Indeno(1,2,3-cd)pyrene     | 410   | U |
| 53-70-3   | Dibenz(a,h)anthracene      | 410   | U |
| 191-24-2  | Benzo(g,h,i)perylene       | 410   | U |

(1) Cannot be separated from Diphenylamine

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1G  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

BNP10

Lab Name: AATSLA

Contract: 68-W0-0081

SS-04C

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.08

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 320F04.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: 19 decanted: (Y/N) N

Date Extracted: 11/08/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/15/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 8.3

Extraction: (Type) SONC

Number TICs found: 3

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

| CAS NUMBER | COMPOUND NAME | RT    | EST. CONC. | Q  |
|------------|---------------|-------|------------|----|
| 1.         | UNKNOWN       | 2.37  | 130        | J  |
| 2.         | UNKNOWN       | 3.19  | 170        | BJ |
| 3.         | UNKNOWN       | 10.08 | 95         | BJ |
| 4.         |               |       |            |    |
| 5.         |               |       |            |    |
| 6.         |               |       |            |    |
| 7.         |               |       |            |    |
| 8.         |               |       |            |    |
| 9.         |               |       |            |    |
| 10.        |               |       |            |    |
| 11.        |               |       |            |    |
| 12.        |               |       |            |    |
| 13.        |               |       |            |    |
| 14.        |               |       |            |    |
| 15.        |               |       |            |    |
| 16.        |               |       |            |    |
| 17.        |               |       |            |    |
| 18.        |               |       |            |    |
| 19.        |               |       |            |    |
| 20.        |               |       |            |    |
| 21.        |               |       |            |    |
| 22.        |               |       |            |    |
| 23.        |               |       |            |    |
| 24.        |               |       |            |    |
| 25.        |               |       |            |    |
| 26.        |               |       |            |    |
| 27.        |               |       |            |    |
| 28.        |               |       |            |    |
| 29.        |               |       |            |    |
| 30.        |               |       |            |    |

R  
R



000495

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP11

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.09

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 320F05.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: 20 decanted: (Y/N) N

Date Extracted: 11/08/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/15/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

Extraction: (Type) SONC  
CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

| CAS NO.  | COMPOUND                     | UG/KG | Q |
|----------|------------------------------|-------|---|
| 100-52-7 | Benzaldehyde                 | 410   | U |
| 108-95-2 | Phenol                       | 410   | U |
| 111-44-4 | bis(2-Chloroethyl) ether     | 410   | U |
| 95-57-8  | 2-Chlorophenol               | 410   | U |
| 95-48-7  | 2-Methylphenol               | 410   | U |
| 108-60-1 | 2,2'-oxybis(1-Chloropropane) | 410   | U |
| 98-86-2  | Acetophenone                 | 410   | U |
| 106-44-5 | 4-Methylphenol               | 410   | U |
| 621-64-7 | N-Nitroso-di-n-propylamine   | 410   | U |
| 67-72-1  | Hexachloroethane             | 410   | U |
| 98-95-3  | Nitrobenzene                 | 410   | U |
| 78-59-1  | Isophorone                   | 410   | U |
| 88-75-5  | 2-Nitrophenol                | 410   | U |
| 105-67-9 | 2,4-Dimethylphenol           | 410   | U |
| 111-91-1 | bis(2-Chloroethoxy) methane  | 410   | U |
| 120-83-2 | 2,4-Dichlorophenol           | 410   | U |
| 91-20-3  | Naphthalene                  | 410   | U |
| 106-47-8 | 4-Chloroaniline              | 410   | U |
| 87-68-3  | Hexachlorobutadiene          | 410   | U |
| 105-60-2 | Caprolactam                  | 410   | U |
| 59-50-7  | 4-Chloro-3-methylphenol      | 410   | U |
| 91-57-6  | 2-Methylnaphthalene          | 410   | U |
| 77-47-4  | Hexachlorocyclopentadiene    | 410   | U |
| 88-06-2  | 2,4,6-Trichlorophenol        | 410   | U |
| 95-95-4  | 2,4,5-Trichlorophenol        | 1000  | U |
| 92-52-4  | 1,1'-Biphenyl                | 410   | U |
| 91-58-7  | 2-Chloronaphthalene          | 410   | U |
| 88-74-4  | 2-Nitroaniline               | 1000  | U |
| 131-11-3 | Dimethylphthalate            | 410   | U |
| 606-20-2 | 2,6-Dinitrotoluene           | 410   | U |
| 208-96-8 | Acenaphthylene               | 410   | U |
| 99-09-2  | 3-Nitroaniline               | 1000  | U |
| 83-32-9  | Acenaphthene                 | 410   | U |

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1D  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP11

SS-04D

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) SOIL Lab Sample ID: 45798.09  
 Sample wt/vol: 30.0 (g/mL) G Lab File ID: 320F05.D  
 Level: (low/med) LOW Date Received: 11/08/00  
 % Moisture: 20 decanted: (Y/N) N Date Extracted: 11/08/00  
 Concentrated Extract Volume: 500 (uL) Date Analyzed: 11/15/00  
 Injection Volume: 2.0 (uL) Dilution Factor: 1.0  
 GPC Cleanup: (Y/N) Y pH: 7.0 Extraction: (Type) SONC

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

Q

| CAS NO.   | COMPOUND                   | UG/KG | Q |
|-----------|----------------------------|-------|---|
| 51-28-5   | 2,4-Dinitrophenol          | 1000  | U |
| 100-02-7  | 4-Nitrophenol              | 1000  | U |
| 132-64-9  | Dibenzofuran               | 410   | U |
| 121-14-2  | 2,4-Dinitrotoluene         | 410   | U |
| 84-66-2   | Diethylphthalate           | 410   | U |
| 86-73-7   | Fluorene                   | 410   | U |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 410   | U |
| 100-01-6  | 4-Nitroaniline             | 1000  | U |
| 534-52-1  | 4,6-Dinitro-2-methylphenol | 1000  | U |
| 86-30-6   | N-Nitrosodiphenylamine (1) | 410   | U |
| 101-55-3  | 4-Bromophenyl-phenylether  | 410   | U |
| 118-74-1  | Hexachlorobenzene          | 410   | U |
| 1912-24-9 | Atrazine                   | 410   | U |
| 87-86-5   | Pentachlorophenol          | 1000  | U |
| 85-01-8   | Phenanthrene               | 410   | U |
| 120-12-7  | Anthracene                 | 410   | U |
| 86-74-8   | Carbazole                  | 410   | U |
| 84-74-2   | Di-n-butylphthalate        | 410   | U |
| 206-44-0  | Fluoranthene               | 410   | U |
| 129-00-0  | Pyrene                     | 410   | U |
| 85-68-7   | Butylbenzylphthalate       | 410   | U |
| 91-94-1   | 3,3'-Dichlorobenzidine     | 410   | U |
| 56-55-3   | Benzo(a)anthracene         | 410   | U |
| 218-01-9  | Chrysene                   | 410   | U |
| 117-81-7  | bis(2-Ethylhexyl)phthalate | 240   | J |
| 117-84-0  | Di-n-octylphthalate        | 410   | U |
| 205-99-2  | Benzo(b)fluoranthene       | 410   | U |
| 207-08-9  | Benzo(k)fluoranthene       | 410   | U |
| 50-32-8   | Benzo(a)pyrene             | 410   | U |
| 193-39-5  | Indeno(1,2,3-cd)pyrene     | 410   | U |
| 53-70-3   | Dibenz(a,h)anthracene      | 410   | U |
| 191-24-2  | Benzo(g,h,i)perylene       | 410   | U |

(1) Cannot be separated from Diphenylamine

000497

1G  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

BNP11

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

SS-04 D

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.09

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 320F05.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: 20 decanted: (Y/N) N

Date Extracted: 11/08/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/15/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

Extraction: (Type) SONC

Number TICs found: 3

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

| CAS NUMBER | COMPOUND NAME | RT    | EST. CONC. | Q  |
|------------|---------------|-------|------------|----|
| 1.         | UNKNOWN       | 2.37  | 120        | J  |
| 2.         | UNKNOWN       | 3.19  | 160        | BJ |
| 3.         | UNKNOWN       | 10.08 | 120        | BJ |
| 4.         |               |       |            |    |
| 5.         |               |       |            |    |
| 6.         |               |       |            |    |
| 7.         |               |       |            |    |
| 8.         |               |       |            |    |
| 9.         |               |       |            |    |
| 10.        |               |       |            |    |
| 11.        |               |       |            |    |
| 12.        |               |       |            |    |
| 13.        |               |       |            |    |
| 14.        |               |       |            |    |
| 15.        |               |       |            |    |
| 16.        |               |       |            |    |
| 17.        |               |       |            |    |
| 18.        |               |       |            |    |
| 19.        |               |       |            |    |
| 20.        |               |       |            |    |
| 21.        |               |       |            |    |
| 22.        |               |       |            |    |
| 23.        |               |       |            |    |
| 24.        |               |       |            |    |
| 25.        |               |       |            |    |
| 26.        |               |       |            |    |
| 27.        |               |       |            |    |
| 28.        |               |       |            |    |
| 29.        |               |       |            |    |
| 30.        |               |       |            |    |

R  
R

P. 148

000504

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP12

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

SS-05A

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.10

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 320F06.D

Level: (low/med) LOW

Date Received: 11/08/00 ✓

% Moisture: 21 ✓ decanted: (Y/N) N

Date Extracted: 11/08/00 ✓

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/15/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y ✓ pH: 7.3 ✓

Extraction: (Type) SONC  
CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

CAS NO. COMPOUND

| CAS NO.  | COMPOUND                     | UG/KG | Q |
|----------|------------------------------|-------|---|
| 100-52-7 | Benzaldehyde                 | 420   | U |
| 108-95-2 | Phenol                       | 420   | U |
| 111-44-4 | bis(2-Chloroethyl) ether     | 420   | U |
| 95-57-8  | 2-Chlorophenol               | 420   | U |
| 95-48-7  | 2-Methylphenol               | 420   | U |
| 108-60-1 | 2,2'-oxybis(1-Chloropropane) | 420   | U |
| 98-86-2  | Acetophenone                 | 420   | U |
| 106-44-5 | 4-Methylphenol               | 420   | U |
| 621-64-7 | N-Nitroso-di-n-propylamine   | 420   | U |
| 67-72-1  | Hexachloroethane             | 420   | U |
| 98-95-3  | Nitrobenzene                 | 420   | U |
| 78-59-1  | Isophorone                   | 420   | U |
| 88-75-5  | 2-Nitrophenol                | 420   | U |
| 105-67-9 | 2,4-Dimethylphenol           | 420   | U |
| 111-91-1 | bis(2-Chloroethoxy)methane   | 420   | U |
| 120-83-2 | 2,4-Dichlorophenol           | 420   | U |
| 91-20-3  | Naphthalene                  | 420   | U |
| 106-47-8 | 4-Chloroaniline              | 420   | U |
| 87-68-3  | Hexachlorobutadiene          | 420   | U |
| 105-60-2 | Caprolactam                  | 420   | U |
| 59-50-7  | 4-Chloro-3-methylphenol      | 420   | U |
| 91-57-6  | 2-Methylnaphthalene          | 420   | U |
| 77-47-4  | Hexachlorocyclopentadiene    | 420   | U |
| 88-06-2  | 2,4,6-Trichlorophenol        | 420   | U |
| 95-95-4  | 2,4,5-Trichlorophenol        | 1100  | U |
| 92-52-4  | 1,1'-Biphenyl                | 420   | U |
| 91-58-7  | 2-Chloronaphthalene          | 420   | U |
| 88-74-4  | 2-Nitroaniline               | 1100  | U |
| 131-11-3 | Dimethylphthalate            | 420   | U |
| 606-20-2 | 2,6-Dinitrotoluene           | 420   | U |
| 208-96-8 | Acenaphthylene               | 420   | U |
| 99-09-2  | 3-Nitroaniline               | 1100  | U |
| 83-32-9  | Acenaphthene                 | 420   | U |

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1D  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP12

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.10

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 320F06.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: 21 decanted: (Y/N) N

Date Extracted: 11/08/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/15/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.3

Extraction: (Type) SONC  
CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

| CAS NO.   | COMPOUND                   |      |   |
|-----------|----------------------------|------|---|
| 51-28-5   | 2,4-Dinitrophenol          | 1100 | U |
| 100-02-7  | 4-Nitrophenol              | 1100 | U |
| 132-64-9  | Dibenzofuran               | 420  | U |
| 121-14-2  | 2,4-Dinitrotoluene         | 420  | U |
| 84-66-2   | Diethylphthalate           | 420  | U |
| 86-73-7   | Fluorene                   | 420  | U |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 420  | U |
| 100-01-6  | 4-Nitroaniline             | 1100 | U |
| 534-52-1  | 4,6-Dinitro-2-methylphenol | 1100 | U |
| 86-30-6   | N-Nitrosodiphenylamine (1) | 420  | U |
| 101-55-3  | 4-Bromophenyl-phenylether  | 420  | U |
| 118-74-1  | Hexachlorobenzene          | 420  | U |
| 1912-24-9 | Atrazine                   | 420  | U |
| 87-86-5   | Pentachlorophenol          | 1100 | U |
| 85-01-8   | Phenanthrene               | 420  | U |
| 120-12-7  | Anthracene                 | 420  | U |
| 86-74-8   | Carbazole                  | 420  | U |
| 84-74-2   | Di-n-butylphthalate        | 420  | U |
| 206-44-0  | Fluoranthene               | 420  | U |
| 129-00-0  | Pyrene                     | 130  | J |
| 85-68-7   | Butylbenzylphthalate       | 420  | U |
| 91-94-1   | 3,3'-Dichlorobenzidine     | 420  | U |
| 56-55-3   | Benzo (a) anthracene       | 420  | U |
| 218-01-9  | Chrysene                   | 420  | U |
| 117-81-7  | bis(2-Ethylhexyl)phthalate | 59   | J |
| 117-84-0  | Di-n-octylphthalate        | 120  | J |
| 205-99-2  | Benzo (b) fluoranthene     | 420  | U |
| 207-08-9  | Benzo (k) fluoranthene     | 420  | U |
| 50-32-8   | Benzo (a) pyrene           | 420  | U |
| 193-39-5  | Indeno (1,2,3-cd) pyrene   | 420  | U |
| 53-70-3   | Dibenz (a,h) anthracene    | 420  | U |
| 191-24-2  | Benzo (g,h,i) perylene     | 420  | U |

(1) Cannot be separated from Diphenylamine

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1G  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

BNP12

Lab Name: AATSLA

Contract: 68-W0-0081

SS-05A

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.10

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 320F06.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: 21 decanted: (Y/N) N

Date Extracted: 11/08/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/15/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y

pH: 7.3

Extraction: (Type) SONC

Number TICs found: 8

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

| CAS NUMBER | COMPOUND NAME | RT   | EST. CONC. | Q    |
|------------|---------------|------|------------|------|
| 1.         | UNKNOWN       | 2.35 | 140        | J    |
| 2.         | UNKNOWN       | 3.19 | 190        | BJ R |
| 3.         | UNKNOWN       | 4.82 | 130        | J    |
| 4.         | UNKNOWN       | 5.04 | 290        | J    |
| 5.         | UNKNOWN       | 5.31 | 86         | J    |
| 6.         | UNKNOWN       | 5.54 | 200        | J    |
| 7.         | UNKNOWN       | 5.58 | 110        | J    |
| 8.         | UNKNOWN       | 5.62 | 340        | J    |
| 9.         |               |      |            |      |
| 10.        |               |      |            |      |
| 11.        |               |      |            |      |
| 12.        |               |      |            |      |
| 13.        |               |      |            |      |
| 14.        |               |      |            |      |
| 15.        |               |      |            |      |
| 16.        |               |      |            |      |
| 17.        |               |      |            |      |
| 18.        |               |      |            |      |
| 19.        |               |      |            |      |
| 20.        |               |      |            |      |
| 21.        |               |      |            |      |
| 22.        |               |      |            |      |
| 23.        |               |      |            |      |
| 24.        |               |      |            |      |
| 25.        |               |      |            |      |
| 26.        |               |      |            |      |
| 27.        |               |      |            |      |
| 28.        |               |      |            |      |
| 29.        |               |      |            |      |
| 30.        |               |      |            |      |

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1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP13

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.11

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 322F05.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: 24 / decanted: (Y/N) N

Date Extracted: 11/08/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/17/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y / pH: 7.1 /

Extraction: (Type) SONC  
CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

| CAS NO.  | COMPOUND                     | UG/KG | Q |
|----------|------------------------------|-------|---|
| 100-52-7 | Benzaldehyde                 | 430   | U |
| 108-95-2 | Phenol                       | 430   | U |
| 111-44-4 | bis(2-Chloroethyl) ether     | 430   | U |
| 95-57-8  | 2-Chlorophenol               | 430   | U |
| 95-48-7  | 2-Methylphenol               | 430   | U |
| 108-60-1 | 2,2'-oxybis(1-Chloropropane) | 430   | U |
| 98-86-2  | Acetophenone                 | 430   | U |
| 106-44-5 | 4-Methylphenol               | 430   | U |
| 621-64-7 | N-Nitroso-di-n-propylamine   | 430   | U |
| 67-72-1  | Hexachloroethane             | 430   | U |
| 98-95-3  | Nitrobenzene                 | 430   | U |
| 78-59-1  | Isophorone                   | 430   | U |
| 88-75-5  | 2-Nitrophenol                | 430   | U |
| 105-67-9 | 2,4-Dimethylphenol           | 430   | U |
| 111-91-1 | bis(2-Chloroethoxy) methane  | 430   | U |
| 120-83-2 | 2,4-Dichlorophenol           | 430   | U |
| 91-20-3  | Naphthalene                  | 430   | U |
| 106-47-8 | 4-Chloroaniline              | 430   | U |
| 87-68-3  | Hexachlorobutadiene          | 430   | U |
| 105-60-2 | Caprolactam                  | 430   | U |
| 59-50-7  | 4-Chloro-3-methylphenol      | 430   | U |
| 91-57-6  | 2-Methylnaphthalene          | 430   | U |
| 77-47-4  | Hexachlorocyclopentadiene    | 430   | U |
| 88-06-2  | 2,4,6-Trichlorophenol        | 430   | U |
| 95-95-4  | 2,4,5-Trichlorophenol        | 1100  | U |
| 92-52-4  | 1,1'-Biphenyl                | 430   | U |
| 91-58-7  | 2-Chloronaphthalene          | 430   | U |
| 88-74-4  | 2-Nitroaniline               | 1100  | U |
| 131-11-3 | Dimethylphthalate            | 430   | U |
| 606-20-2 | 2,6-Dinitrotoluene           | 430   | U |
| 208-96-8 | Acenaphthylene               | 430   | U |
| 99-09-2  | 3-Nitroaniline               | 1100  | U |
| 83-32-9  | Acenaphthene                 | 430   | U |

1D  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP13

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.11

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 322F05.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: 24 decanted: (Y/N) N

Date Extracted: 11/08/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/17/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.1

Extraction: (Type) SONC  
CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

| CAS NO.   | COMPOUND                   |      |   |
|-----------|----------------------------|------|---|
| 51-28-5   | 2,4-Dinitrophenol          | 1100 | U |
| 100-02-7  | 4-Nitrophenol              | 1100 | U |
| 132-64-9  | Dibenzofuran               | 430  | U |
| 121-14-2  | 2,4-Dinitrotoluene         | 430  | U |
| 84-66-2   | Diethylphthalate           | 430  | U |
| 86-73-7   | Fluorene                   | 430  | U |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 430  | U |
| 100-01-6  | 4-Nitroaniline             | 1100 | U |
| 534-52-1  | 4,6-Dinitro-2-methylphenol | 1100 | U |
| 86-30-6   | N-Nitrosodiphenylamine (1) | 430  | U |
| 101-55-3  | 4-Bromophenyl-phenylether  | 430  | U |
| 118-74-1  | Hexachlorobenzene          | 430  | U |
| 1912-24-9 | Atrazine                   | 430  | U |
| 87-86-5   | Pentachlorophenol          | 1100 | U |
| 85-01-8   | Phenanthrene               | 150  | J |
| 120-12-7  | Anthracene                 | 430  | U |
| 86-74-8   | Carbazole                  | 430  | U |
| 84-74-2   | Di-n-butylphthalate        | 430  | U |
| 206-44-0  | Fluoranthene               | 430  | U |
| 129-00-0  | Pyrene                     | 430  | U |
| 85-68-7   | Butylbenzylphthalate       | 430  | U |
| 91-94-1   | 3,3'-Dichlorobenzidine     | 430  | U |
| 56-55-3   | Benzo(a)anthracene         | 430  | U |
| 218-01-9  | Chrysene                   | 430  | U |
| 117-81-7  | bis(2-Ethylhexyl)phthalate | 430  | U |
| 117-84-0  | Di-n-octylphthalate        | 430  | U |
| 205-99-2  | Benzo(b)fluoranthene       | 430  | U |
| 207-08-9  | Benzo(k)fluoranthene       | 430  | U |
| 50-32-8   | Benzo(a)pyrene             | 430  | U |
| 193-39-5  | Indeno(1,2,3-cd)pyrene     | 430  | U |
| 53-70-3   | Dibenz(a,h)anthracene      | 430  | U |
| 191-24-2  | Benzo(g,h,i)perylene       | 430  | U |

(1) Cannot be seperated from Diphenylamine



1G  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

BNP13

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

SS-058

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.11

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 322F05.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: 24 decanted: (Y/N) N

Date Extracted: 11/08/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/17/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.1

Extraction: (Type) SONC

Number TICs found: 7

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

| CAS NUMBER | COMPOUND NAME                | RT   | EST. CONC. | Q    |
|------------|------------------------------|------|------------|------|
| 1.         | UNKNOWN                      | 3.19 | 240        | BJ R |
| 2.         | UNKNOWN                      | 4.82 | 200        | J    |
| 3.         | UNKNOWN                      | 4.99 | 99         | J    |
| 4.         | DECAHYDRO-4,4,8,9,10-PENTAME | 5.03 | 330        | JN   |
| 5.         | UNKNOWN                      | 5.53 | 270        | J    |
| 6.         | UNKNOWN                      | 5.62 | 340        | J    |
| 7.         | UNKNOWN                      | 6.33 | 260        | J    |
| 8.         |                              |      |            |      |
| 9.         |                              |      |            |      |
| 10.        |                              |      |            |      |
| 11.        |                              |      |            |      |
| 12.        |                              |      |            |      |
| 13.        |                              |      |            |      |
| 14.        |                              |      |            |      |
| 15.        |                              |      |            |      |
| 16.        |                              |      |            |      |
| 17.        |                              |      |            |      |
| 18.        |                              |      |            |      |
| 19.        |                              |      |            |      |
| 20.        |                              |      |            |      |
| 21.        |                              |      |            |      |
| 22.        |                              |      |            |      |
| 23.        |                              |      |            |      |
| 24.        |                              |      |            |      |
| 25.        |                              |      |            |      |
| 26.        |                              |      |            |      |
| 27.        |                              |      |            |      |
| 28.        |                              |      |            |      |
| 29.        |                              |      |            |      |
| 30.        |                              |      |            |      |

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNQ85

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.12

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 320F08.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: 25 decanted: (Y/N) N

Date Extracted: 11/08/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/15/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

Extraction: (Type) SONC  
CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

| CAS NO.  | COMPOUND                     |      |   |
|----------|------------------------------|------|---|
| 100-52-7 | Benzaldehyde                 | 440  | U |
| 108-95-2 | Phenol                       | 440  | U |
| 111-44-4 | bis(2-Chloroethyl) ether     | 440  | U |
| 95-57-8  | 2-Chlorophenol               | 440  | U |
| 95-48-7  | 2-Methylphenol               | 440  | U |
| 108-60-1 | 2,2'-oxybis(1-Chloropropane) | 440  | U |
| 98-86-2  | Acetophenone                 | 440  | U |
| 106-44-5 | 4-Methylphenol               | 440  | U |
| 621-64-7 | N-Nitroso-di-n-propylamine   | 440  | U |
| 67-72-1  | Hexachloroethane             | 440  | U |
| 98-95-3  | Nitrobenzene                 | 440  | U |
| 78-59-1  | Isophorone                   | 440  | U |
| 88-75-5  | 2-Nitrophenol                | 440  | U |
| 105-67-9 | 2,4-Dimethylphenol           | 440  | U |
| 111-91-1 | bis(2-Chloroethoxy)methane   | 440  | U |
| 120-83-2 | 2,4-Dichlorophenol           | 440  | U |
| 91-20-3  | Naphthalene                  | 440  | U |
| 106-47-8 | 4-Chloroaniline              | 440  | U |
| 87-68-3  | Hexachlorobutadiene          | 440  | U |
| 105-60-2 | Caprolactam                  | 440  | U |
| 59-50-7  | 4-Chloro-3-methylphenol      | 440  | U |
| 91-57-6  | 2-Methylnaphthalene          | 440  | U |
| 77-47-4  | Hexachlorocyclopentadiene    | 440  | U |
| 88-06-2  | 2,4,6-Trichlorophenol        | 440  | U |
| 95-95-4  | 2,4,5-Trichlorophenol        | 1100 | U |
| 92-52-4  | 1,1'-Biphenyl                | 440  | U |
| 91-58-7  | 2-Chloronaphthalene          | 440  | U |
| 88-74-4  | 2-Nitroaniline               | 1100 | U |
| 131-11-3 | Dimethylphthalate            | 440  | U |
| 606-20-2 | 2,6-Dinitrotoluene           | 440  | U |
| 208-96-8 | Acenaphthylene               | 440  | U |
| 99-09-2  | 3-Nitroaniline               | 1100 | U |
| 83-32-9  | Acenaphthene                 | 440  | U |

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1D  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNQ85

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) SOIL Lab Sample ID: 45798.12  
 Sample wt/vol: 30.0 (g/mL) G Lab File ID: 320F08.D  
 Level: (low/med) LOW Date Received: 11/08/00  
 % Moisture: 25 decanted: (Y/N) N Date Extracted: 11/08/00  
 Concentrated Extract Volume: 500 (uL) Date Analyzed: 11/15/00  
 Injection Volume: 2.0 (uL) Dilution Factor: 1.0  
 GPC Cleanup: (Y/N) Y pH: 7.0 Extraction: (Type) SONC

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

| CAS NO.   | COMPOUND                   |      |   |
|-----------|----------------------------|------|---|
| 51-28-5   | 2,4-Dinitrophenol          | 1100 | U |
| 100-02-7  | 4-Nitrophenol              | 1100 | U |
| 132-64-9  | Dibenzofuran               | 440  | U |
| 121-14-2  | 2,4-Dinitrotoluene         | 440  | U |
| 84-66-2   | Diethylphthalate           | 440  | U |
| 86-73-7   | Fluorene                   | 440  | U |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 440  | U |
| 100-01-6  | 4-Nitroaniline             | 1100 | U |
| 534-52-1  | 4,6-Dinitro-2-methylphenol | 1100 | U |
| 86-30-6   | N-Nitrosodiphenylamine (1) | 440  | U |
| 101-55-3  | 4-Bromophenyl-phenylether  | 440  | U |
| 118-74-1  | Hexachlorobenzene          | 440  | U |
| 1912-24-9 | Atrazine                   | 440  | U |
| 87-86-5   | Pentachlorophenol          | 1100 | U |
| 85-01-8   | Phenanthrene               | 440  | U |
| 120-12-7  | Anthracene                 | 440  | U |
| 86-74-8   | Carbazole                  | 440  | U |
| 84-74-2   | Di-n-butylphthalate        | 440  | U |
| 206-44-0  | Fluoranthene               | 440  | U |
| 129-00-0  | Pyrene                     | 440  | U |
| 85-68-7   | Butylbenzylphthalate       | 440  | U |
| 91-94-1   | 3,3'-Dichlorobenzidine     | 440  | U |
| 56-55-3   | Benzo(a)anthracene         | 440  | U |
| 218-01-9  | Chrysene                   | 440  | U |
| 117-81-7  | bis(2-Ethylhexyl)phthalate | 78   | J |
| 117-84-0  | Di-n-octylphthalate        | 440  | U |
| 205-99-2  | Benzo(b)fluoranthene       | 440  | U |
| 207-08-9  | Benzo(k)fluoranthene       | 440  | U |
| 50-32-8   | Benzo(a)pyrene             | 440  | U |
| 193-39-5  | Indeno(1,2,3-cd)pyrene     | 440  | U |
| 53-70-3   | Dibenz(a,h)anthracene      | 440  | U |
| 191-24-2  | Benzo(g,h,i)perylene       | 440  | U |

(1) Cannot be seperated from Diphenylamine

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1G  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

BNQ85

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.12

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 320F08.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: 25 decanted: (Y/N) N

Date Extracted: 11/08/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/15/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

Extraction: (Type) SONC

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

Number TICs found: 3

| CAS NUMBER | COMPOUND NAME | RT   | EST. CONC. | Q  |
|------------|---------------|------|------------|----|
| 1.         | UNKNOWN       | 2.36 | 120        | J  |
| 2.         | UNKNOWN       | 3.20 | 200        | BJ |
| 3.         | UNKNOWN       | 6.87 | 260        | J  |
| 4.         |               |      |            |    |
| 5.         |               |      |            |    |
| 6.         |               |      |            |    |
| 7.         |               |      |            |    |
| 8.         |               |      |            |    |
| 9.         |               |      |            |    |
| 10.        |               |      |            |    |
| 11.        |               |      |            |    |
| 12.        |               |      |            |    |
| 13.        |               |      |            |    |
| 14.        |               |      |            |    |
| 15.        |               |      |            |    |
| 16.        |               |      |            |    |
| 17.        |               |      |            |    |
| 18.        |               |      |            |    |
| 19.        |               |      |            |    |
| 20.        |               |      |            |    |
| 21.        |               |      |            |    |
| 22.        |               |      |            |    |
| 23.        |               |      |            |    |
| 24.        |               |      |            |    |
| 25.        |               |      |            |    |
| 26.        |               |      |            |    |
| 27.        |               |      |            |    |
| 28.        |               |      |            |    |
| 29.        |               |      |            |    |
| 30.        |               |      |            |    |

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000545

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNQ87

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

SS-06 A

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.13

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 320F09.D

Level: (low/med) LOW

Date Received: 11/08/00 ✓

% Moisture: 27 ✓ decanted: (Y/N) N

Date Extracted: 11/08/00 ✓

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/15/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y ✓

pH: 7.3 ✓

Extraction: (Type) SONC  
CONCENTRATION UNITS:

CAS NO. COMPOUND

(ug/L or ug/Kg) UG/KG

Q

| CAS NO.  | COMPOUND                     | UG/KG | Q |
|----------|------------------------------|-------|---|
| 100-52-7 | Benzaldehyde                 | 450   | U |
| 108-95-2 | Phenol                       | 450   | U |
| 111-44-4 | bis(2-Chloroethyl) ether     | 450   | U |
| 95-57-8  | 2-Chlorophenol               | 450   | U |
| 95-48-7  | 2-Methylphenol               | 450   | U |
| 108-60-1 | 2,2'-oxybis(1-Chloropropane) | 450   | U |
| 98-86-2  | Acetophenone                 | 450   | U |
| 106-44-5 | 4-Methylphenol               | 450   | U |
| 621-64-7 | N-Nitroso-di-n-propylamine   | 450   | U |
| 67-72-1  | Hexachloroethane             | 450   | U |
| 98-95-3  | Nitrobenzene                 | 450   | U |
| 78-59-1  | Isophorone                   | 450   | U |
| 88-75-5  | 2-Nitrophenol                | 450   | U |
| 105-67-9 | 2,4-Dimethylphenol           | 450   | U |
| 111-91-1 | bis(2-Chloroethoxy)methane   | 450   | U |
| 120-83-2 | 2,4-Dichlorophenol           | 450   | U |
| 91-20-3  | Naphthalene                  | 450   | U |
| 106-47-8 | 4-Chloroaniline              | 450   | U |
| 87-68-3  | Hexachlorobutadiene          | 450   | U |
| 105-60-2 | Caprolactam                  | 450   | U |
| 59-50-7  | 4-Chloro-3-methylphenol      | 450   | U |
| 91-57-6  | 2-Methylnaphthalene          | 450   | U |
| 77-47-4  | Hexachlorocyclopentadiene    | 450   | U |
| 88-06-2  | 2,4,6-Trichlorophenol        | 450   | U |
| 95-95-4  | 2,4,5-Trichlorophenol        | 1100  | U |
| 92-52-4  | 1,1'-Biphenyl                | 450   | U |
| 91-58-7  | 2-Chloronaphthalene          | 450   | U |
| 88-74-4  | 2-Nitroaniline               | 1100  | U |
| 131-11-3 | Dimethylphthalate            | 450   | U |
| 606-20-2 | 2,6-Dinitrotoluene           | 450   | U |
| 208-96-8 | Acenaphthylene               | 450   | U |
| 99-09-2  | 3-Nitroaniline               | 1100  | U |
| 83-32-9  | Acenaphthene                 | 450   | U |

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1D  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNQ87

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

SS-06A

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.13

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 320F09.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: 27 decanted: (Y/N) N

Date Extracted: 11/08/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/15/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y

pH: 7.3

Extraction: (Type) SONC  
CONCENTRATION UNITS:

CAS NO. COMPOUND

(ug/L or ug/Kg) UG/KG

Q

|           |                            |      |   |
|-----------|----------------------------|------|---|
| 51-28-5   | 2,4-Dinitrophenol          | 1100 | U |
| 100-02-7  | 4-Nitrophenol              | 1100 | U |
| 132-64-9  | Dibenzofuran               | 450  | U |
| 121-14-2  | 2,4-Dinitrotoluene         | 450  | U |
| 84-66-2   | Diethylphthalate           | 450  | U |
| 86-73-7   | Fluorene                   | 450  | U |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 450  | U |
| 100-01-6  | 4-Nitroaniline             | 1100 | U |
| 534-52-1  | 4,6-Dinitro-2-methylphenol | 1100 | U |
| 86-30-6   | N-Nitrosodiphenylamine (1) | 450  | U |
| 101-55-3  | 4-Bromophenyl-phenylether  | 450  | U |
| 118-74-1  | Hexachlorobenzene          | 450  | U |
| 1912-24-9 | Atrazine                   | 450  | U |
| 87-86-5   | Pentachlorophenol          | 1100 | U |
| 85-01-8   | Phenanthrene               | 450  | U |
| 120-12-7  | Anthracene                 | 450  | U |
| 86-74-8   | Carbazole                  | 450  | U |
| 84-74-2   | Di-n-butylphthalate        | 450  | U |
| 206-44-0  | Fluoranthene               | 450  | U |
| 129-00-0  | Pyrene                     | 450  | U |
| 85-68-7   | Butylbenzylphthalate       | 450  | U |
| 91-94-1   | 3,3'-Dichlorobenzidine     | 450  | U |
| 56-55-3   | Benzo(a)anthracene         | 450  | U |
| 218-01-9  | Chrysene                   | 450  | U |
| 117-81-7  | bis(2-Ethylhexyl)phthalate | 450  | U |
| 117-84-0  | Di-n-octylphthalate        | 450  | U |
| 205-99-2  | Benzo(b)fluoranthene       | 450  | U |
| 207-08-9  | Benzo(k)fluoranthene       | 450  | U |
| 50-32-8   | Benzo(a)pyrene             | 190  | J |
| 193-39-5  | Indeno(1,2,3-cd)pyrene     | 450  | U |
| 53-70-3   | Dibenz(a,h)anthracene      | 450  | U |
| 191-24-2  | Benzo(g,h,i)perylene       | 450  | U |

(1) Cannot be separated from Diphenylamine

000547

1G  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

BNQ87

Lab Name: AATSLA

Contract: 68-W0-0081

SS-06A

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.13

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 320F09.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: 27 decanted: (Y/N) N

Date Extracted: 11/08/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/15/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y

pH: 7.3

Extraction: (Type) SONC

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

Number TICs found: 6

| CAS NUMBER | COMPOUND NAME                            | RT    | EST. CONC. | Q  |
|------------|--|-------|------------|----|
| 1.         | UNKNOWN                                  | 2.37  | 180        | J  |
| 2.         | UNKNOWN                                  | 3.20  | 220        | BJ |
| 3.         | 540-97-6<br>CYCLOHEXASILOXANE, DODECAMET | 4.04  | 93         | JN |
| 4.         | UNKNOWN                                  | 6.86  | 420        | J  |
| 5.         | UNKNOWN                                  | 10.09 | 160        | BJ |
| 6.         | UNKNOWN                                  | 15.18 | 95         | J  |
| 7.         |  |       |            |    |
| 8.         |  |       |            |    |
| 9.         |  |       |            |    |
| 10.        |  |       |            |    |
| 11.        |  |       |            |    |
| 12.        |  |       |            |    |
| 13.        |  |       |            |    |
| 14.        |  |       |            |    |
| 15.        |  |       |            |    |
| 16.        |  |       |            |    |
| 17.        |  |       |            |    |
| 18.        |  |       |            |    |
| 19.        |  |       |            |    |
| 20.        |  |       |            |    |
| 21.        |  |       |            |    |
| 22.        |  |       |            |    |
| 23.        |  |       |            |    |
| 24.        |  |       |            |    |
| 25.        |  |       |            |    |
| 26.        |  |       |            |    |
| 27.        |  |       |            |    |
| 28.        |  |       |            |    |
| 29.        |  |       |            |    |
| 30.        |  |       |            |    |

R  
R  
R

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1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNQ88

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.14

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 320F10.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: 20 decanted: (Y/N) N

Date Extracted: 11/08/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/15/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.5

 Extraction: (Type) SONC  
 CONCENTRATION UNITS:  
 (ug/L or ug/Kg) UG/KG Q

CAS NO. COMPOUND

| CAS NO.  | COMPOUND                     | UG/KG | Q |
|----------|------------------------------|-------|---|
| 100-52-7 | Benzaldehyde                 | 410   | U |
| 108-95-2 | Phenol                       | 410   | U |
| 111-44-4 | bis(2-Chloroethyl) ether     | 410   | U |
| 95-57-8  | 2-Chlorophenol               | 410   | U |
| 95-48-7  | 2-Methylphenol               | 410   | U |
| 108-60-1 | 2,2'-oxybis(1-Chloropropane) | 410   | U |
| 98-86-2  | Acetophenone                 | 410   | U |
| 106-44-5 | 4-Methylphenol               | 410   | U |
| 621-64-7 | N-Nitroso-di-n-propylamine   | 410   | U |
| 67-72-1  | Hexachloroethane             | 410   | U |
| 98-95-3  | Nitrobenzene                 | 410   | U |
| 78-59-1  | Isophorone                   | 410   | U |
| 88-75-5  | 2-Nitrophenol                | 410   | U |
| 105-67-9 | 2,4-Dimethylphenol           | 410   | U |
| 111-91-1 | bis(2-Chloroethoxy)methane   | 410   | U |
| 120-83-2 | 2,4-Dichlorophenol           | 410   | U |
| 91-20-3  | Naphthalene                  | 410   | U |
| 106-47-8 | 4-Chloroaniline              | 410   | U |
| 87-68-3  | Hexachlorobutadiene          | 410   | U |
| 105-60-2 | Caprolactam                  | 410   | U |
| 59-50-7  | 4-Chloro-3-methylphenol      | 410   | U |
| 91-57-6  | 2-Methylnaphthalene          | 410   | U |
| 77-47-4  | Hexachlorocyclopentadiene    | 410   | U |
| 88-06-2  | 2,4,6-Trichlorophenol        | 410   | U |
| 95-95-4  | 2,4,5-Trichlorophenol        | 1000  | U |
| 92-52-4  | 1,1'-Biphenyl                | 410   | U |
| 91-58-7  | 2-Chloronaphthalene          | 410   | U |
| 88-74-4  | 2-Nitroaniline               | 1000  | U |
| 131-11-3 | Dimethylphthalate            | 410   | U |
| 606-20-2 | 2,6-Dinitrotoluene           | 410   | U |
| 208-96-8 | Acenaphthylene               | 410   | U |
| 99-09-2  | 3-Nitroaniline               | 1000  | U |
| 83-32-9  | Acenaphthene                 | 410   | U |



1D  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNQ88

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.14

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 320F10.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: 20 decanted: (Y/N) N

Date Extracted: 11/08/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/15/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.5

Extraction: (Type) SONC

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

Q

| CAS NO.   | COMPOUND                   |      |   |
|-----------|----------------------------|------|---|
| 51-28-5   | 2,4-Dinitrophenol          | 1000 | U |
| 100-02-7  | 4-Nitrophenol              | 1000 | U |
| 132-64-9  | Dibenzofuran               | 410  | U |
| 121-14-2  | 2,4-Dinitrotoluene         | 410  | U |
| 84-66-2   | Diethylphthalate           | 410  | U |
| 86-73-7   | Fluorene                   | 410  | U |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 410  | U |
| 100-01-6  | 4-Nitroaniline             | 1000 | U |
| 534-52-1  | 4,6-Dinitro-2-methylphenol | 1000 | U |
| 86-30-6   | N-Nitrosodiphenylamine (1) | 410  | U |
| 101-55-3  | 4-Bromophenyl-phenylether  | 410  | U |
| 118-74-1  | Hexachlorobenzene          | 410  | U |
| 1912-24-9 | Atrazine                   | 410  | U |
| 87-86-5   | Pentachlorophenol          | 1000 | U |
| 85-01-8   | Phenanthrene               | 410  | U |
| 120-12-7  | Anthracene                 | 410  | U |
| 86-74-8   | Carbazole                  | 410  | U |
| 84-74-2   | Di-n-butylphthalate        | 410  | U |
| 206-44-0  | Fluoranthene               | 410  | U |
| 129-00-0  | Pyrene                     | 410  | U |
| 85-68-7   | Butylbenzylphthalate       | 410  | U |
| 91-94-1   | 3,3'-Dichlorobenzidine     | 410  | U |
| 56-55-3   | Benzo (a) anthracene       | 410  | U |
| 218-01-9  | Chrysene                   | 410  | U |
| 117-81-7  | bis(2-Ethylhexyl)phthalate | 100  | J |
| 117-84-0  | Di-n-octylphthalate        | 410  | U |
| 205-99-2  | Benzo (b) fluoranthene     | 410  | U |
| 207-08-9  | Benzo (k) fluoranthene     | 410  | U |
| 50-32-8   | Benzo (a) pyrene           | 410  | U |
| 193-39-5  | Indeno (1,2,3-cd) pyrene   | 410  | U |
| 53-70-3   | Dibenz (a,h) anthracene    | 410  | U |
| 191-24-2  | Benzo (g,h,i) perylene     | 410  | U |

(1) Cannot be separated from Diphenylamine

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000559

1G  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

Lab Name: AATSLA

Contract: 68-W0-0081

BNQ88

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

55-063

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.14

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 320F10.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: 20 decanted: (Y/N) N

Date Extracted: 11/08/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/15/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y

pH: 7.5

Extraction: (Type) SONC

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

Number TICs found: 4

| CAS NUMBER   | COMPOUND NAME                | RT    | EST. CONC. | Q    |
|--------------|------------------------------|-------|------------|------|
| 1.           | UNKNOWN                      | 2.36  | 95         | J    |
| 2. 541-02-6  | CYCLOPENTASILOXANE, DECAMETH | 3.20  | 170        | JN R |
| 3. 2057-49-0 | PYRIDINE, 4-(3-PHENYLPROPYL) | 6.86  | 170        | JN   |
| 4.           | UNKNOWN                      | 10.08 | 90         | BJ   |
| 5.           |                              |       |            |      |
| 6.           |                              |       |            |      |
| 7.           |                              |       |            |      |
| 8.           |                              |       |            |      |
| 9.           |                              |       |            |      |
| 10.          |                              |       |            |      |
| 11.          |                              |       |            |      |
| 12.          |                              |       |            |      |
| 13.          |                              |       |            |      |
| 14.          |                              |       |            |      |
| 15.          |                              |       |            |      |
| 16.          |                              |       |            |      |
| 17.          |                              |       |            |      |
| 18.          |                              |       |            |      |
| 19.          |                              |       |            |      |
| 20.          |                              |       |            |      |
| 21.          |                              |       |            |      |
| 22.          |                              |       |            |      |
| 23.          |                              |       |            |      |
| 24.          |                              |       |            |      |
| 25.          |                              |       |            |      |
| 26.          |                              |       |            |      |
| 27.          |                              |       |            |      |
| 28.          |                              |       |            |      |
| 29.          |                              |       |            |      |
| 30.          |                              |       |            |      |

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000567

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNQ89

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.15

Sample wt/vol: / 30.0 (g/mL) G

Lab File ID: 320F11.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: 22 decanted: (Y/N) N

Date Extracted: 11/08/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/15/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y / pH: 7.1 /

Extraction: (Type) SONC  
CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

| CAS NO.  | COMPOUND                     |      |   |
|----------|------------------------------|------|---|
| 100-52-7 | Benzaldehyde                 | 420  | U |
| 108-95-2 | Phenol                       | 420  | U |
| 111-44-4 | bis(2-Chloroethyl) ether     | 420  | U |
| 95-57-8  | 2-Chlorophenol               | 420  | U |
| 95-48-7  | 2-Methylphenol               | 420  | U |
| 108-60-1 | 2,2'-oxybis(1-Chloropropane) | 420  | U |
| 98-86-2  | Acetophenone                 | 420  | U |
| 106-44-5 | 4-Methylphenol               | 420  | U |
| 621-64-7 | N-Nitroso-di-n-propylamine   | 420  | U |
| 67-72-1  | Hexachloroethane             | 420  | U |
| 98-95-3  | Nitrobenzene                 | 420  | U |
| 78-59-1  | Isophorone                   | 420  | U |
| 88-75-5  | 2-Nitrophenol                | 420  | U |
| 105-67-9 | 2,4-Dimethylphenol           | 420  | U |
| 111-91-1 | bis(2-Chloroethoxy) methane  | 420  | U |
| 120-83-2 | 2,4-Dichlorophenol           | 420  | U |
| 91-20-3  | Naphthalene                  | 420  | U |
| 106-47-8 | 4-Chloroaniline              | 420  | U |
| 87-68-3  | Hexachlorobutadiene          | 420  | U |
| 105-60-2 | Caprolactam                  | 420  | U |
| 59-50-7  | 4-Chloro-3-methylphenol      | 420  | U |
| 91-57-6  | 2-Methylnaphthalene          | 420  | U |
| 77-47-4  | Hexachlorocyclopentadiene    | 420  | U |
| 88-06-2  | 2,4,6-Trichlorophenol        | 420  | U |
| 95-95-4  | 2,4,5-Trichlorophenol        | 1100 | U |
| 92-52-4  | 1,1'-Biphenyl                | 420  | U |
| 91-58-7  | 2-Chloronaphthalene          | 420  | U |
| 88-74-4  | 2-Nitroaniline               | 1100 | U |
| 131-11-3 | Dimethylphthalate            | 420  | U |
| 606-20-2 | 2,6-Dinitrotoluene           | 420  | U |
| 208-96-8 | Acenaphthylene               | 420  | U |
| 99-09-2  | 3-Nitroaniline               | 1100 | U |
| 83-32-9  | Acenaphthene                 | 420  | U |

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1D  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNQ89

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.15

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 320F11.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: 22 decanted: (Y/N) N

Date Extracted: 11/08/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/15/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.1

Extraction: (Type) SONC  
CONCENTRATION UNITS:

CAS NO. COMPOUND

(ug/L or ug/Kg) UG/KG

Q

|           |                            |      |   |
|-----------|----------------------------|------|---|
| 51-28-5   | 2,4-Dinitrophenol          | 1100 | U |
| 100-02-7  | 4-Nitrophenol              | 1100 | U |
| 132-64-9  | Dibenzofuran               | 420  | U |
| 121-14-2  | 2,4-Dinitrotoluene         | 420  | U |
| 84-66-2   | Diethylphthalate           | 420  | U |
| 86-73-7   | Fluorene                   | 420  | U |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 420  | U |
| 100-01-6  | 4-Nitroaniline             | 1100 | U |
| 534-52-1  | 4,6-Dinitro-2-methylphenol | 1100 | U |
| 86-30-6   | N-Nitrosodiphenylamine (1) | 420  | U |
| 101-55-3  | 4-Bromophenyl-phenylether  | 420  | U |
| 118-74-1  | Hexachlorobenzene          | 420  | U |
| 1912-24-9 | Atrazine                   | 420  | U |
| 87-86-5   | Pentachlorophenol          | 1100 | U |
| 85-01-8   | Phenanthrene               | 420  | U |
| 120-12-7  | Anthracene                 | 420  | U |
| 86-74-8   | Carbazole                  | 420  | U |
| 84-74-2   | Di-n-butylphthalate        | 420  | U |
| 206-44-0  | Fluoranthene               | 420  | U |
| 129-00-0  | Pyrene                     | 420  | U |
| 85-68-7   | Butylbenzylphthalate       | 420  | U |
| 91-94-1   | 3,3'-Dichlorobenzidine     | 420  | U |
| 56-55-3   | Benzo(a)anthracene         | 420  | U |
| 218-01-9  | Chrysene                   | 420  | U |
| 117-81-7  | bis(2-Ethylhexyl)phthalate | 81   | J |
| 117-84-0  | Di-n-octylphthalate        | 420  | U |
| 205-99-2  | Benzo(b)fluoranthene       | 420  | U |
| 207-08-9  | Benzo(k)fluoranthene       | 420  | U |
| 50-32-8   | Benzo(a)pyrene             | 420  | U |
| 193-39-5  | Indeno(1,2,3-cd)pyrene     | 420  | U |
| 53-70-3   | Dibenz(a,h)anthracene      | 420  | U |
| 191-24-2  | Benzo(g,h,i)perylene       | 420  | U |

(1) Cannot be seperated from Diphenylamine

1G  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

BNQ89

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: <sup>SS-06 C</sup>BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.15

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 320F11.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: 22 decanted: (Y/N) N

Date Extracted: 11/08/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/15/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.1

Extraction: (Type) SONC

Number TICs found: 4

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

| CAS NUMBER | COMPOUND NAME | RT    | EST. CONC. | Q    |
|------------|---------------|-------|------------|------|
| 1.         | UNKNOWN       | 2.36  | 120        | J    |
| 2.         | UNKNOWN       | 3.20  | 170        | BJ R |
| 3.         | UNKNOWN       | 6.86  | 130        | J    |
| 4.         | UNKNOWN       | 10.08 | 97         | BJ R |
| 5.         |               |       |            |      |
| 6.         |               |       |            |      |
| 7.         |               |       |            |      |
| 8.         |               |       |            |      |
| 9.         |               |       |            |      |
| 10.        |               |       |            |      |
| 11.        |               |       |            |      |
| 12.        |               |       |            |      |
| 13.        |               |       |            |      |
| 14.        |               |       |            |      |
| 15.        |               |       |            |      |
| 16.        |               |       |            |      |
| 17.        |               |       |            |      |
| 18.        |               |       |            |      |
| 19.        |               |       |            |      |
| 20.        |               |       |            |      |
| 21.        |               |       |            |      |
| 22.        |               |       |            |      |
| 23.        |               |       |            |      |
| 24.        |               |       |            |      |
| 25.        |               |       |            |      |
| 26.        |               |       |            |      |
| 27.        |               |       |            |      |
| 28.        |               |       |            |      |
| 29.        |               |       |            |      |
| 30.        |               |       |            |      |

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1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNQ91

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

SS-03 (Dup)

Matrix: (soil/water) SOIL ✓

Lab Sample ID: 45798.16

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 320F12.D

Level: (low/med) LOW

Date Received: 11/08/00 ✓

% Moisture: 22 ✓ decanted: (Y/N) N

Date Extracted: 11/08/00 ✓

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/15/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y ✓ pH: 7.3 ✓

 Extraction: (Type) SONC  
 CONCENTRATION UNITS:  
 (ug/L or ug/Kg) UG/KG Q

CAS NO. COMPOUND

| CAS NO.  | COMPOUND                     | UG/KG | Q |
|----------|------------------------------|-------|---|
| 100-52-7 | Benzaldehyde                 | 420   | U |
| 108-95-2 | Phenol                       | 420   | U |
| 111-44-4 | bis(2-Chloroethyl) ether     | 420   | U |
| 95-57-8  | 2-Chlorophenol               | 420   | U |
| 95-48-7  | 2-Methylphenol               | 420   | U |
| 108-60-1 | 2,2'-oxybis(1-Chloropropane) | 420   | U |
| 98-86-2  | Acetophenone                 | 420   | U |
| 106-44-5 | 4-Methylphenol               | 420   | U |
| 621-64-7 | N-Nitroso-di-n-propylamine   | 420   | U |
| 67-72-1  | Hexachloroethane             | 420   | U |
| 98-95-3  | Nitrobenzene                 | 420   | U |
| 78-59-1  | Isophorone                   | 420   | U |
| 88-75-5  | 2-Nitrophenol                | 420   | U |
| 105-67-9 | 2,4-Dimethylphenol           | 420   | U |
| 111-91-1 | bis(2-Chloroethoxy)methane   | 420   | U |
| 120-83-2 | 2,4-Dichlorophenol           | 420   | U |
| 91-20-3  | Naphthalene                  | 420   | U |
| 106-47-8 | 4-Chloroaniline              | 420   | U |
| 87-68-3  | Hexachlorobutadiene          | 420   | U |
| 105-60-2 | Caprolactam                  | 420   | U |
| 59-50-7  | 4-Chloro-3-methylphenol      | 420   | U |
| 91-57-6  | 2-Methylnaphthalene          | 420   | U |
| 77-47-4  | Hexachlorocyclopentadiene    | 420   | U |
| 88-06-2  | 2,4,6-Trichlorophenol        | 420   | U |
| 95-95-4  | 2,4,5-Trichlorophenol        | 1100  | U |
| 92-52-4  | 1,1'-Biphenyl                | 420   | U |
| 91-58-7  | 2-Chloronaphthalene          | 420   | U |
| 88-74-4  | 2-Nitroaniline               | 1100  | U |
| 131-11-3 | Dimethylphthalate            | 420   | U |
| 606-20-2 | 2,6-Dinitrotoluene           | 420   | U |
| 208-96-8 | Acenaphthylene               | 420   | U |
| 99-09-2  | 3-Nitroaniline               | 1100  | U |
| 83-32-9  | Acenaphthene                 | 420   | U |

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1D  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNQ91

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.16

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 320F12.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: 22 decanted: (Y/N) N

Date Extracted: 11/08/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/15/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.3

Extraction: (Type) SONC  
CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

| CAS NO.   | COMPOUND                   |      |   |
|-----------|----------------------------|------|---|
| 51-28-5   | 2,4-Dinitrophenol          | 1100 | U |
| 100-02-7  | 4-Nitrophenol              | 1100 | U |
| 132-64-9  | Dibenzofuran               | 420  | U |
| 121-14-2  | 2,4-Dinitrotoluene         | 420  | U |
| 84-66-2   | Diethylphthalate           | 420  | U |
| 86-73-7   | Fluorene                   | 420  | U |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 420  | U |
| 100-01-6  | 4-Nitroaniline             | 1100 | U |
| 534-52-1  | 4,6-Dinitro-2-methylphenol | 1100 | U |
| 86-30-6   | N-Nitrosodiphenylamine (1) | 420  | U |
| 101-55-3  | 4-Bromophenyl-phenylether  | 420  | U |
| 118-74-1  | Hexachlorobenzene          | 420  | U |
| 1912-24-9 | Atrazine                   | 420  | U |
| 87-86-5   | Pentachlorophenol          | 1100 | U |
| 85-01-8   | Phenanthrene               | 420  | U |
| 120-12-7  | Anthracene                 | 420  | U |
| 86-74-8   | Carbazole                  | 420  | U |
| 84-74-2   | Di-n-butylphthalate        | 420  | U |
| 206-44-0  | Fluoranthene               | 420  | U |
| 129-00-0  | Pyrene                     | 420  | U |
| 85-68-7   | Butylbenzylphthalate       | 420  | U |
| 91-94-1   | 3,3'-Dichlorobenzidine     | 420  | U |
| 56-55-3   | Benzo(a)anthracene         | 420  | U |
| 218-01-9  | Chrysene                   | 420  | U |
| 117-81-7  | bis(2-Ethylhexyl)phthalate | 83   | J |
| 117-84-0  | Di-n-octylphthalate        | 420  | U |
| 205-99-2  | Benzo(b)fluoranthene       | 420  | U |
| 207-08-9  | Benzo(k)fluoranthene       | 420  | U |
| 50-32-8   | Benzo(a)pyrene             | 420  | U |
| 193-39-5  | Indeno(1,2,3-cd)pyrene     | 420  | U |
| 53-70-3   | Dibenz(a,h)anthracene      | 420  | U |
| 191-24-2  | Benzo(g,h,i)perylene       | 420  | U |

(1) Cannot be seperated from Diphenylamine

1G  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

Lab Name: AATSLA

Contract: 68-W0-0081

BNQ91

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: <sup>SS-03 (dup)</sup> BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.16

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: 320F12.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: 22 decanted: (Y/N) N

Date Extracted: 11/08/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 11/15/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.3

Extraction: (Type) SONC

Number TICs found: 3

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

| CAS NUMBER | COMPOUND NAME | RT    | EST. CONC.     | Q             |
|------------|---------------|-------|----------------|---------------|
| 1.         | UNKNOWN       | 2.36  | 98             | J             |
| 2.         | UNKNOWN       | 3.20  | <del>190</del> | <del>BJ</del> |
| 3.         | UNKNOWN       | 10.08 | <del>230</del> | <del>BJ</del> |
| 4.         |               |       |                |               |
| 5.         |               |       |                |               |
| 6.         |               |       |                |               |
| 7.         |               |       |                |               |
| 8.         |               |       |                |               |
| 9.         |               |       |                |               |
| 10.        |               |       |                |               |
| 11.        |               |       |                |               |
| 12.        |               |       |                |               |
| 13.        |               |       |                |               |
| 14.        |               |       |                |               |
| 15.        |               |       |                |               |
| 16.        |               |       |                |               |
| 17.        |               |       |                |               |
| 18.        |               |       |                |               |
| 19.        |               |       |                |               |
| 20.        |               |       |                |               |
| 21.        |               |       |                |               |
| 22.        |               |       |                |               |
| 23.        |               |       |                |               |
| 24.        |               |       |                |               |
| 25.        |               |       |                |               |
| 26.        |               |       |                |               |
| 27.        |               |       |                |               |
| 28.        |               |       |                |               |
| 29.        |               |       |                |               |
| 30.        |               |       |                |               |

RP

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1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNQ92 **FB**

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: <sup>RB-01</sup>BNP05

Matrix: (soil/water) WATER

Lab Sample ID: 45798.17

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: 320F15.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: \_\_\_\_\_ decanted: (Y/N) N

Date Extracted: 11/13/00

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 11/15/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N /

pH: 7.0 /

Extraction: (Type) CONT  
CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L

Q

| CAS NO.  | COMPOUND                     |    |   |
|----------|------------------------------|----|---|
| 100-52-7 | Benzaldehyde                 | 10 | U |
| 108-95-2 | Phenol                       | 10 | U |
| 111-44-4 | bis(2-Chloroethyl) ether     | 10 | U |
| 95-57-8  | 2-Chlorophenol               | 10 | U |
| 95-48-7  | 2-Methylphenol               | 10 | U |
| 108-60-1 | 2,2'-oxybis(1-Chloropropane) | 10 | U |
| 98-86-2  | Acetophenone                 | 10 | U |
| 106-44-5 | 4-Methylphenol               | 10 | U |
| 621-64-7 | N-Nitroso-di-n-propylamine   | 10 | U |
| 67-72-1  | Hexachloroethane             | 10 | U |
| 98-95-3  | Nitrobenzene                 | 10 | U |
| 78-59-1  | Isophorone                   | 10 | U |
| 88-75-5  | 2-Nitrophenol                | 10 | U |
| 105-67-9 | 2,4-Dimethylphenol           | 10 | U |
| 111-91-1 | bis(2-Chloroethoxy) methane  | 10 | U |
| 120-83-2 | 2,4-Dichlorophenol           | 10 | U |
| 91-20-3  | Naphthalene                  | 10 | U |
| 106-47-8 | 4-Chloroaniline              | 10 | U |
| 87-68-3  | Hexachlorobutadiene          | 10 | U |
| 105-60-2 | Caprolactam                  | 10 | U |
| 59-50-7  | 4-Chloro-3-methylphenol      | 10 | U |
| 91-57-6  | 2-Methylnaphthalene          | 10 | U |
| 77-47-4  | Hexachlorocyclopentadiene    | 10 | U |
| 88-06-2  | 2,4,6-Trichlorophenol        | 10 | U |
| 95-95-4  | 2,4,5-Trichlorophenol        | 25 | U |
| 92-52-4  | 1,1'-Biphenyl                | 10 | U |
| 91-58-7  | 2-Chloronaphthalene          | 10 | U |
| 88-74-4  | 2-Nitroaniline               | 25 | U |
| 131-11-3 | Dimethylphthalate            | 10 | U |
| 606-20-2 | 2,6-Dinitrotoluene           | 10 | U |
| 208-96-8 | Acenaphthylene               | 10 | U |
| 99-09-2  | 3-Nitroaniline               | 25 | U |
| 83-32-9  | Acenaphthene                 | 10 | U |

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1D  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNQ92

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

RB-01

Matrix: (soil/water) WATER

Lab Sample ID: 45798.17

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: 320F15.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: \_\_\_\_\_ decanted: (Y/N) N

Date Extracted: 11/13/00

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 11/15/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N

pH: 7.0

Extraction: (Type) CONT  
CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L

CAS NO. COMPOUND

Q

|           |                            |    |   |
|-----------|----------------------------|----|---|
| 51-28-5   | 2,4-Dinitrophenol          | 25 | U |
| 100-02-7  | 4-Nitrophenol              | 25 | U |
| 132-64-9  | Dibenzofuran               | 10 | U |
| 121-14-2  | 2,4-Dinitrotoluene         | 10 | U |
| 84-66-2   | Diethylphthalate           | 10 | U |
| 86-73-7   | Fluorene                   | 10 | U |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 10 | U |
| 100-01-6  | 4-Nitroaniline             | 25 | U |
| 534-52-1  | 4,6-Dinitro-2-methylphenol | 25 | U |
| 86-30-6   | N-Nitrosodiphenylamine (1) | 10 | U |
| 101-55-3  | 4-Bromophenyl-phenylether  | 10 | U |
| 118-74-1  | Hexachlorobenzene          | 10 | U |
| 1912-24-9 | Atrazine                   | 10 | U |
| 87-86-5   | Pentachlorophenol          | 25 | U |
| 85-01-8   | Phenanthrene               | 10 | U |
| 120-12-7  | Anthracene                 | 10 | U |
| 86-74-8   | Carbazole                  | 10 | U |
| 84-74-2   | Di-n-butylphthalate        | 10 | U |
| 206-44-0  | Fluoranthene               | 10 | U |
| 129-00-0  | Pyrene                     | 10 | U |
| 85-68-7   | Butylbenzylphthalate       | 10 | U |
| 91-94-1   | 3,3'-Dichlorobenzidine     | 10 | U |
| 56-55-3   | Benzo(a)anthracene         | 10 | U |
| 218-01-9  | Chrysene                   | 10 | U |
| 117-81-7  | bis(2-Ethylhexyl)phthalate | 2  | J |
| 117-84-0  | Di-n-octylphthalate        | 10 | U |
| 205-99-2  | Benzo(b)fluoranthene       | 10 | U |
| 207-08-9  | Benzo(k)fluoranthene       | 10 | U |
| 50-32-8   | Benzo(a)pyrene             | 10 | U |
| 193-39-5  | Indeno(1,2,3-cd)pyrene     | 10 | U |
| 53-70-3   | Dibenz(a,h)anthracene      | 10 | U |
| 191-24-2  | Benzo(g,h,i)perylene       | 10 | U |

(1) Cannot be seperated from Diphenylamine

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1G  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

BNQ92

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) WATER

Lab Sample ID: 45798.17

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: 320F15.D

Level: (low/med) LOW

Date Received: 11/08/00

% Moisture: \_\_\_\_\_ decanted: (Y/N) N

Date Extracted: 11/13/00

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 11/15/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N

pH: 7.0

Extraction: (Type) CONT

Number TICs found: 2

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L

| CAS NUMBER    | COMPOUND NAME               | RT   | EST. CONC. | Q              |
|---------------|-----------------------------|------|------------|----------------|
| 1.            | UNKNOWN                     | 2.32 | 9          | <del>BJ</del>  |
| 2. 21400-25-9 | 1-PROPENE, 1,1,2-TRICHLORO- | 2.37 | 9          | <del>BJN</del> |
| 3.            |                             |      |            |                |
| 4.            |                             |      |            |                |
| 5.            |                             |      |            |                |
| 6.            |                             |      |            |                |
| 7.            |                             |      |            |                |
| 8.            |                             |      |            |                |
| 9.            |                             |      |            |                |
| 10.           |                             |      |            |                |
| 11.           |                             |      |            |                |
| 12.           |                             |      |            |                |
| 13.           |                             |      |            |                |
| 14.           |                             |      |            |                |
| 15.           |                             |      |            |                |
| 16.           |                             |      |            |                |
| 17.           |                             |      |            |                |
| 18.           |                             |      |            |                |
| 19.           |                             |      |            |                |
| 20.           |                             |      |            |                |
| 21.           |                             |      |            |                |
| 22.           |                             |      |            |                |
| 23.           |                             |      |            |                |
| 24.           |                             |      |            |                |
| 25.           |                             |      |            |                |
| 26.           |                             |      |            |                |
| 27.           |                             |      |            |                |
| 28.           |                             |      |            |                |
| 29.           |                             |      |            |                |
| 30.           |                             |      |            |                |

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1E  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP05

Lab Name: AATSLA Contract: 68-W0-0081

Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05

Matrix: (soil/water) SOIL Lab Sample ID: 45798.01

Sample wt/vol: 30.0 (g/mL) G Lab File ID: \_\_\_\_\_

% Moisture: 27 Decanted: (Y/N) N Date Received: 11/08/00

Extraction: (Type) SONC Date Extracted: 11/08/00

Concentrated Extract Volume: 5000 (uL) Date Analyzed: 11/15/00

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 4.2 Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

| CAS NO.    | COMPOUND            |     |      |       |
|------------|---------------------|-----|------|-------|
| 319-84-6   | alpha-BHC           |     | 2.3  | U     |
| 319-85-7   | beta-BHC            |     | 2.3  | U     |
| 319-86-8   | delta-BHC           |     | 2.3  | U     |
| 58-89-9    | gamma-BHC (Lindane) | 2.3 | 0.15 | PJB U |
| 76-44-8    | Heptachlor          | 2.3 | 0.27 | PJ U  |
| 309-00-2   | Aldrin              |     | 2.3  | U     |
| 1024-57-3  | Heptachlor epoxide  | 2.3 | 0.17 | PJ U  |
| 959-98-8   | Endosulfan I        |     | 2.3  | U     |
| 60-57-1    | Dieldrin            |     | 0.68 | J     |
| 72-55-9    | 4,4'-DDE            |     | 4.5  | U     |
| 72-20-8    | Endrin              | 4.5 | 0.42 | U JB  |
| 33213-65-9 | Endosulfan II       |     | 4.5  | U     |
| 72-54-8    | 4,4'-DDD            |     | 4.5  | U     |
| 1031-07-8  | Endosulfan sulfate  | 4.5 | 0.26 | U PJ  |
| 50-29-3    | 4,4'-DDT            |     | 4.5  | U     |
| 72-43-5    | Methoxychlor        | 23  | 1.1  | PJB U |
| 53494-70-5 | Endrin ketone       |     | 4.5  | U     |
| 7421-93-4  | Endrin aldehyde     |     | 4.5  | U     |
| 5103-71-9  | alpha-Chlordane     |     | 2.3  | U     |
| 5103-74-2  | gamma-Chlordane     |     | 0.27 | J PJ  |
| 8001-35-2  | Toxaphene           |     | 230  | U     |
| 12674-11-2 | Aroclor-1016        |     | 45   | U     |
| 11104-28-2 | Aroclor-1221        |     | 92   | U     |
| 11141-16-5 | Aroclor-1232        |     | 45   | U     |
| 53469-21-9 | Aroclor-1242        |     | 45   | U     |
| 12672-29-6 | Aroclor-1248        |     | 45   | U     |
| 11097-69-1 | Aroclor-1254        |     | 45   | U     |
| 11096-82-5 | Aroclor-1260        |     | 45   | U     |

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000730

1E  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP06

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.04

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: \_\_\_\_\_

% Moisture: 28 Decanted: (Y/N) N

Date Received: 11/08/00

Extraction: (Type) SONC

Date Extracted: 11/08/00

Concentrated Extract Volume: 5000 (uL)

Date Analyzed: 11/17/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 5.4

Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

Q

| CAS NO.    | COMPOUND            |     |         |       |
|------------|---------------------|-----|---------|-------|
| 319-84-6   | alpha-BHC           | 2.4 | 0.42    | PJ U  |
| 319-85-7   | beta-BHC            |     | 4.0     | PR R  |
| 319-86-8   | delta-BHC           |     | 2.4     | U     |
| 58-89-9    | gamma-BHC (Lindane) |     | 2.4     | U     |
| 76-44-8    | Heptachlor          |     | 2.4     | U     |
| 309-00-2   | Aldrin              |     | 2.4     | U     |
| 1024-57-3  | Heptachlor epoxide  |     | 4.6     | U     |
| 959-98-8   | Endosulfan I        |     | 2.4     | U     |
| 60-57-1    | Dieldrin            |     | 4.6     | U     |
| 72-55-9    | 4,4'-DDE            |     | 4.6     | U     |
| 72-20-8    | Endrin              |     | 4.6     | U     |
| 33213-65-9 | Endosulfan II       |     | 2.5     | J     |
| 72-54-8    | 4,4'-DDD            | 4.6 | 1.5     | PJ U  |
| 1031-07-8  | Endosulfan sulfate  |     | 4.6     | U     |
| 50-29-3    | 4,4'-DDT            | 4.6 | 0.79    | PJB U |
| 72-43-5    | Methoxychlor        |     | 24 0.91 | JB U  |
| 53494-70-5 | Endrin ketone       |     | 4.6     | U     |
| 7421-93-4  | Endrin aldehyde     | 4.6 | 1.0     | PJB U |
| 5103-71-9  | alpha-Chlordane     |     | 2.4     | U     |
| 5103-74-2  | gamma-Chlordane     | 2.4 | 0.52    | PJ U  |
| 8001-35-2  | Toxaphene           |     | 240     | U     |
| 12674-11-2 | Aroclor-1016        |     | 46      | U     |
| 11104-28-2 | Aroclor-1221        |     | 93      | U     |
| 11141-16-5 | Aroclor-1232        |     | 46      | U     |
| 53469-21-9 | Aroclor-1242        |     | 46      | U     |
| 12672-29-6 | Aroclor-1248        |     | 46      | U     |
| 11097-69-1 | Aroclor-1254        |     | 46      | U     |
| 11096-82-5 | Aroclor-1260        |     | 46      | U     |

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1E  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP07

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) SOIL Lab Sample ID: 45798.05  
 Sample wt/vol: 30.0 (g/mL) G Lab File ID: \_\_\_\_\_  
 % Moisture: 21 Decanted: (Y/N) N Date Received: 11/08/00  
 Extraction: (Type) SONC Date Extracted: 11/08/00  
 Concentrated Extract Volume: 5000 (uL) Date Analyzed: 11/15/00  
 Injection Volume: 2.0 (uL) Dilution Factor: 1.0  
 GPC Cleanup: (Y/N) Y pH: 4.7 Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

| CAS NO.    | COMPOUND            | UG/KG    | Q     |
|------------|---------------------|----------|-------|
| 319-84-6   | alpha-BHC           | 0.38     | J     |
| 319-85-7   | beta-BHC            | 2.2      | U     |
| 319-86-8   | delta-BHC           | 2.2      | U     |
| 58-89-9    | gamma-BHC (Lindane) | 2.2      | U     |
| 76-44-8    | Heptachlor          | 2.2      | U     |
| 309-00-2   | Aldrin              | 2.2      | U     |
| 1024-57-3  | Heptachlor epoxide  | 0.77     | J     |
| 959-98-8   | Endosulfan I        | 2.2      | U     |
| 60-57-1    | Dieldrin            | 4.2      | U     |
| 72-55-9    | 4,4'-DDE            | 4.2      | U     |
| 72-20-8    | Endrin              | 4.2      | U     |
| 33213-65-9 | Endosulfan II       | 4.2      | U     |
| 72-54-8    | 4,4'-DDD            | 4.2      | U     |
| 1031-07-8  | Endosulfan sulfate  | 4.2      | U     |
| 50-29-3    | 4,4'-DDT            | 4.2      | U     |
| 72-43-5    | Methoxychlor        | 22 0.43  | PJB U |
| 53494-70-5 | Endrin ketone       | 4.2      | U     |
| 7421-93-4  | Endrin aldehyde     | 4.2 0.13 | PJB U |
| 5103-71-9  | alpha-Chlordane     | 2.2      | U     |
| 5103-74-2  | gamma-Chlordane     | 2.2 0.14 | PJB U |
| 8001-35-2  | Toxaphene           | 220      | U     |
| 12674-11-2 | Aroclor-1016        | 42       | U     |
| 11104-28-2 | Aroclor-1221        | 85       | U     |
| 11141-16-5 | Aroclor-1232        | 42       | U     |
| 53469-21-9 | Aroclor-1242        | 42       | U     |
| 12672-29-6 | Aroclor-1248        | 42       | U     |
| 11097-69-1 | Aroclor-1254        | 42       | U     |
| 11096-82-5 | Aroclor-1260        | 42       | U     |

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1E  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP08

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.06

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: \_\_\_\_\_

% Moisture: 3 Decanted: (Y/N) N

Date Received: 11/08/00

Extraction: (Type) SONC

Date Extracted: 11/08/00

Concentrated Extract Volume: 5000 (uL)

Date Analyzed: 11/15/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.0

Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

Q

| CAS NO.    | COMPOUND            |     |      |       |
|------------|---------------------|-----|------|-------|
| 319-84-6   | alpha-BHC           | 1.8 | 0.38 | PJ U  |
| 319-85-7   | beta-BHC            |     | 1.8  | U     |
| 319-86-8   | delta-BHC           |     | 1.8  | U     |
| 58-89-9    | gamma-BHC (Lindane) |     | 1.8  | U     |
| 76-44-8    | Heptachlor          |     | 1.8  | U     |
| 309-00-2   | Aldrin              |     | 1.8  | U     |
| 1024-57-3  | Heptachlor epoxide  | 1.8 | 1.2  | PJ U  |
| 959-98-8   | Endosulfan I        |     | 1.8  | U     |
| 60-57-1    | Dieldrin            | 3.4 | 1.5  | PJ U  |
| 72-55-9    | 4,4'-DDE            | 3.4 | 1.1  | PJ U  |
| 72-20-8    | Endrin              | 3.4 | 0.37 | PJB U |
| 33213-65-9 | Endosulfan II       | 3.4 | 0.86 | PJ U  |
| 72-54-8    | 4,4'-DDD            | 3.4 | 0.46 | PJ U  |
| 1031-07-8  | Endosulfan sulfate  | 3.4 | 0.68 | PJ U  |
| 50-29-3    | 4,4'-DDT            | 3.4 | 0.44 | PJB U |
| 72-43-5    | Methoxychlor        | 18  | 1.5  | PJB U |
| 53494-70-5 | Endrin ketone       | 3.4 | 0.38 | PJB U |
| 7421-93-4  | Endrin aldehyde     | 3.4 | 0.87 | PJB U |
| 5103-71-9  | alpha-Chlordane     | 1.8 | 1.4  | PJ U  |
| 5103-74-2  | gamma-Chlordane     | 1.8 | 0.71 | PJ U  |
| 8001-35-2  | Toxaphene           |     | 180  | U     |
| 12674-11-2 | Aroclor-1016        |     | 34   | U     |
| 11104-28-2 | Aroclor-1221        |     | 69   | U     |
| 11141-16-5 | Aroclor-1232        |     | 34   | U     |
| 53469-21-9 | Aroclor-1242        |     | 34   | U     |
| 12672-29-6 | Aroclor-1248        |     | 170  | U     |
| 11097-69-1 | Aroclor-1254        |     | 34   | U     |
| 11096-82-5 | Aroclor-1260        |     | 34   | U     |

0.176

1E  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: AATSLA

Contract: 68-W0-0081

BNP09

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.07

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: \_\_\_\_\_

% Moisture: 10 Decanted: (Y/N) N

Date Received: 11/08/00

Extraction: (Type) SONC

Date Extracted: 11/08/00

Concentrated Extract Volume: 5000 (uL)

Date Analyzed: 11/16/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 9.2

Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

| CAS NO.    | COMPOUND            |     |      | Q     |
|------------|---------------------|-----|------|-------|
| 319-84-6   | alpha-BHC           |     | 1.9  | U     |
| 319-85-7   | beta-BHC            |     | 1.9  | U     |
| 319-86-8   | delta-BHC           |     | 1.9  | U     |
| 58-89-9    | gamma-BHC (Lindane) |     | 1.9  | U     |
| 76-44-8    | Heptachlor          |     | 1.9  | U     |
| 309-00-2   | Aldrin              |     | 1.9  | U     |
| 1024-57-3  | Heptachlor epoxide  | 1.9 | 0.75 | PJ U  |
| 959-98-8   | Endosulfan I        |     | 1.9  | U     |
| 60-57-1    | Dieldrin            | 3.7 | 0.61 | PJ U  |
| 72-55-9    | 4,4'-DDE            | 3.7 | 1.7  | PJ U  |
| 72-20-8    | Endrin              |     | 3.7  | U     |
| 33213-65-9 | Endosulfan II       | 3.7 | 0.32 | PJ U  |
| 72-54-8    | 4,4'-DDD            | 3.7 | 0.34 | PJ U  |
| 1031-07-8  | Endosulfan sulfate  | 3.7 | 1.3  | PJ U  |
| 50-29-3    | 4,4'-DDT            |     | 3.7  | U     |
| 72-43-5    | Methoxychlor        |     | 19   | U     |
| 53494-70-5 | Endrin ketone       |     | 3.8  | PBJN  |
| 7421-93-4  | Endrin aldehyde     | 3.7 | 2.3  | PJB U |
| 5103-71-9  | alpha-Chlordane     | 1.9 | 0.51 | PJ U  |
| 5103-74-2  | gamma-Chlordane     | 1.9 | 0.43 | PJ U  |
| 8001-35-2  | Toxaphene           |     | 190  | U     |
| 12674-11-2 | Aroclor-1016        |     | 37   | U     |
| 11104-28-2 | Aroclor-1221        |     | 74   | U     |
| 11141-16-5 | Aroclor-1232        |     | 37   | U     |
| 53469-21-9 | Aroclor-1242        |     | 37   | U     |
| 12672-29-6 | Aroclor-1248        |     | 210  | U     |
| 11097-69-1 | Aroclor-1254        |     | 37   | U     |
| 11096-82-5 | Aroclor-1260        |     | 37   | U     |

0.177



1E  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP10

Lab Name: AATSLA Contract: 68-W0-0081

Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05

Matrix: (soil/water) SOIL Lab Sample ID: 45798.08

Sample wt/vol: 30.0 (g/mL) G Lab File ID: \_\_\_\_\_

% Moisture: 19 Decanted: (Y/N) N Date Received: 11/08/00

Extraction: (Type) SONC Date Extracted: 11/08/00

Concentrated Extract Volume: 5000 (uL) Date Analyzed: 11/15/00

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 8.2 Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

| CAS NO.    | COMPOUND            | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) | UG/KG | Q     |
|------------|---------------------|---|-------|-------|
| 319-84-6   | alpha-BHC           |   | 0.34  | J     |
| 319-85-7   | beta-BHC            |   | 2.7   | B     |
| 319-86-8   | delta-BHC           |   | 2.1   | U     |
| 58-89-9    | gamma-BHC (Lindane) |   | 2.1   | U     |
| 76-44-8    | Heptachlor          |   | 2.1   | U     |
| 309-00-2   | Aldrin              |   | 2.1   | U     |
| 1024-57-3  | Heptachlor epoxide  |   | 2.1   | U     |
| 959-98-8   | Endosulfan I        |   | 2.1   | U     |
| 60-57-1    | Dieldrin            | 2.1                                     | 0.12  | PJ U  |
| 72-55-9    | 4,4'-DDE            |   | 4.1   | U     |
| 72-20-8    | Endrin              |   | 4.1   | U     |
| 33213-65-9 | Endosulfan II       |   | 4.1   | U     |
| 72-54-8    | 4,4'-DDD            | 4.1                                     | 0.10  | PJ U  |
| 1031-07-8  | Endosulfan sulfate  |   | 4.1   | U     |
| 50-29-3    | 4,4'-DDT            |   | 4.1   | U     |
| 72-43-5    | Methoxychlor        | 2.1                                     | 0.23  | PJB U |
| 53494-70-5 | Endrin ketone       |   | 4.1   | U     |
| 7421-93-4  | Endrin aldehyde     | 4.1                                     | 0.17  | PJB U |
| 5103-71-9  | alpha-Chlordane     |   | 2.1   | U     |
| 5103-74-2  | gamma-Chlordane     |   | 2.1   | U     |
| 8001-35-2  | Toxaphene           |   | 210   | U     |
| 12674-11-2 | Aroclor-1016        |   | 41    | U     |
| 11104-28-2 | Aroclor-1221        |   | 83    | U     |
| 11141-16-5 | Aroclor-1232        |   | 41    | U     |
| 53469-21-9 | Aroclor-1242        |   | 41    | U     |
| 12672-29-6 | Aroclor-1248        |   | 41    | U     |
| 11097-69-1 | Aroclor-1254        |   | 41    | U     |
| 11096-82-5 | Aroclor-1260        |   | 41    | U     |

0.178

0075

1E  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP11

Lab Name: AATSLA Contract: 68-W0-0081

Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05

Matrix: (soil/water) SOIL Lab Sample ID: 45798.09

Sample wt/vol: 30.0 (g/mL) G Lab File ID: \_\_\_\_\_

% Moisture: 20 Decanted: (Y/N) N Date Received: 11/08/00

Extraction: (Type) SONC Date Extracted: 11/08/00

Concentrated Extract Volume: 5000 (uL) Date Analyzed: 11/15/00

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0 Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

| CAS NO.    | COMPOUND            |     |      | Q     |
|------------|---------------------|-----|------|-------|
| 319-84-6   | alpha-BHC           | 2.1 | 0.25 | PJ U  |
| 319-85-7   | beta-BHC            | 2.1 | 0.90 | PJB U |
| 319-86-8   | delta-BHC           |     | 2.1  | U     |
| 58-89-9    | gamma-BHC (Lindane) |     | 2.1  | U     |
| 76-44-8    | Heptachlor          |     | 2.1  | U     |
| 309-00-2   | Aldrin              |     | 2.1  | U     |
| 1024-57-3  | Heptachlor epoxide  | 2.1 | 0.67 | PJ U  |
| 959-98-8   | Endosulfan I        |     | 2.1  | U     |
| 60-57-1    | Dieldrin            |     | 4.1  | U     |
| 72-55-9    | 4,4'-DDE            |     | 4.1  | U     |
| 72-20-8    | Endrin              |     | 4.1  | U     |
| 33213-65-9 | Endosulfan II       |     | 4.1  | U     |
| 72-54-8    | 4,4'-DDD            |     | 4.1  | U     |
| 1031-07-8  | Endosulfan sulfate  |     | 4.1  | U     |
| 50-29-3    | 4,4'-DDT            |     | 4.1  | U     |
| 72-43-5    | Methoxychlor        | 2.1 | 0.35 | PJB U |
| 53494-70-5 | Endrin ketone       |     | 4.1  | U     |
| 7421-93-4  | Endrin aldehyde     |     | 4.1  | U     |
| 5103-71-9  | alpha-Chlordane     |     | 2.1  | U     |
| 5103-74-2  | gamma-Chlordane     | 2.1 | 0.13 | PJ U  |
| 8001-35-2  | Toxaphene           |     | 210  | U     |
| 12674-11-2 | Aroclor-1016        |     | 41   | U     |
| 11104-28-2 | Aroclor-1221        |     | 84   | U     |
| 11141-16-5 | Aroclor-1232        |     | 41   | U     |
| 53469-21-9 | Aroclor-1242        |     | 41   | U     |
| 12672-29-6 | Aroclor-1248        |     | 41   | U     |
| 11097-69-1 | Aroclor-1254        |     | 41   | U     |
| 11096-82-5 | Aroclor-1260        |     | 41   | U     |

0.179

000754

1E  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP12

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.10

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: \_\_\_\_\_

% Moisture: 21 Decanted: (Y/N) N

Date Received: 11/08/00

Extraction: (Type) SONC

Date Extracted: 11/08/00

Concentrated Extract Volume: 5000 (uL)

Date Analyzed: 11/15/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.3

Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

Q

| CAS NO.    | COMPOUND            |     |      |       |
|------------|---------------------|-----|------|-------|
| 319-84-6   | alpha-BHC           | 2.2 | 0.14 | PJ U  |
| 319-85-7   | beta-BHC            |     | 2.2  | U     |
| 319-86-8   | delta-BHC           |     | 2.2  | U     |
| 58-89-9    | gamma-BHC (Lindane) |     | 2.2  | U     |
| 76-44-8    | Heptachlor          | 2.2 | 0.11 | PJ U  |
| 309-00-2   | Aldrin              |     | 2.2  | U     |
| 1024-57-3  | Heptachlor epoxide  |     | 0.79 | J     |
| 959-98-8   | Endosulfan I        |     | 0.64 | J     |
| 60-57-1    | Dieldrin            | 4.4 | 0.41 | PJ U  |
| 72-55-9    | 4,4'-DDE            | 4.4 | 0.66 | PJ U  |
| 72-20-8    | Endrin              |     | 2.3  | PJB   |
| 33213-65-9 | Endosulfan II       | 4.4 | 1.1  | PJ U  |
| 72-54-8    | 4,4'-DDD            |     | 1.8  | J     |
| 1031-07-8  | Endosulfan sulfate  |     | 6.1  |       |
| 50-29-3    | 4,4'-DDT            | 4.4 | 1.2  | PJB U |
| 72-43-5    | Methoxychlor        |     | 9.1  | J     |
| 53494-70-5 | Endrin ketone       |     | 4.1  | PJB   |
| 7421-93-4  | Endrin aldehyde     |     | 7.7  | PB    |
| 5103-71-9  | alpha-Chlordane     |     | 2.2  | U     |
| 5103-74-2  | gamma-Chlordane     |     | 0.36 | PJ    |
| 8001-35-2  | Toxaphene           |     | 220  | U     |
| 12674-11-2 | Aroclor-1016        |     | 42   | U     |
| 11104-28-2 | Aroclor-1221        |     | 85   | U     |
| 11141-16-5 | Aroclor-1232        |     | 42   | U     |
| 53469-21-9 | Aroclor-1242        |     | 42   | U     |
| 12672-29-6 | Aroclor-1248        |     | 42   | U     |
| 11097-69-1 | Aroclor-1254        |     | 42   | U     |
| 11096-82-5 | Aroclor-1260        |     | 42   | U     |

180

00757

1E  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNP13

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) SOIL Lab Sample ID: 45798.11  
 Sample wt/vol: 30.0 (g/mL) G Lab File ID: \_\_\_\_\_  
 % Moisture: 24 Decanted: (Y/N) N Date Received: 11/08/00  
 Extraction: (Type) SONC Date Extracted: 11/08/00  
 Concentrated Extract Volume: 5000 (uL) Date Analyzed: 11/15/00  
 Injection Volume: 2.0 (uL) Dilution Factor: 1.0  
 GPC Cleanup: (Y/N) Y pH: 7.1 Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

| CAS NO.    | COMPOUND            |     |      | Q     |
|------------|---------------------|-----|------|-------|
| 319-84-6   | alpha-BHC           | 2.2 | 0.15 | PJ U  |
| 319-85-7   | beta-BHC            | 2.2 | 0.17 | PJB U |
| 319-86-8   | delta-BHC           |     | 2.2  | U     |
| 58-89-9    | gamma-BHC (Lindane) |     | 2.2  | U     |
| 76-44-8    | Heptachlor          |     | 0.32 | PJ J  |
| 309-00-2   | Aldrin              |     | 2.2  | U     |
| 1024-57-3  | Heptachlor epoxide  | 4.3 | 0.36 | PJ U  |
| 959-98-8   | Endosulfan I        |     | 0.26 | PJ J  |
| 60-57-1    | Dieldrin            | 4.3 | 0.23 | PJ U  |
| 72-55-9    | 4,4'-DDE            | 4.3 | 0.39 | PJ U  |
| 72-20-8    | Endrin              | 4.3 | 0.97 | PJB U |
| 33213-65-9 | Endosulfan II       |     | 4.3  | U     |
| 72-54-8    | 4,4'-DDD            |     | 0.97 | J     |
| 1031-07-8  | Endosulfan sulfate  |     | 4.3  | U     |
| 50-29-3    | 4,4'-DDT            | 4.3 | 1.2  | PJB U |
| 72-43-5    | Methoxychlor        |     | 7.0  | PJB J |
| 53494-70-5 | Endrin ketone       | 4.3 | 1.9  | PJB U |
| 7421-93-4  | Endrin aldehyde     |     | 5.0  | PB J  |
| 5103-71-9  | alpha-Chlordane     |     | 2.2  | U     |
| 5103-74-2  | gamma-Chlordane     |     | 0.35 | J     |
| 8001-35-2  | Toxaphene           |     | 220  | U     |
| 12674-11-2 | Aroclor-1016        |     | 43   | U     |
| 11104-28-2 | Aroclor-1221        |     | 88   | U     |
| 11141-16-5 | Aroclor-1232        |     | 43   | U     |
| 53469-21-9 | Aroclor-1242        |     | 43   | U     |
| 12672-29-6 | Aroclor-1248        |     | 43   | U     |
| 11097-69-1 | Aroclor-1254        |     | 43   | U     |
| 11096-82-5 | Aroclor-1260        |     | 43   | U     |

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000760

1E  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNQ85

Lab Name: AATSLA Contract: 68-W0-0081

Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05

Matrix: (soil/water) SOIL Lab Sample ID: 45798.12

Sample wt/vol: 30.0 (g/mL) G Lab File ID: \_\_\_\_\_

% Moisture: 25 Decanted: (Y/N) N Date Received: 11/08/00

Extraction: (Type) SONC Date Extracted: 11/08/00

Concentrated Extract Volume: 5000 (uL) Date Analyzed: 11/15/00

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0 Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

| CAS NO.    | COMPOUND            | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) UG/KG | Q     |
|------------|---------------------|---|-------|
| 319-84-6   | alpha-BHC           | 2.3   | U     |
| 319-85-7   | beta-BHC            | 2.3 <del>0.89</del>                           | PJB U |
| 319-86-8   | delta-BHC           | 2.3   | U     |
| 58-89-9    | gamma-BHC (Lindane) | 2.3   | U     |
| 76-44-8    | Heptachlor          | 2.3   | U     |
| 309-00-2   | Aldrin              | 2.3   | U     |
| 1024-57-3  | Heptachlor epoxide  | 2.3 <del>0.34</del>                           | PJB U |
| 959-98-8   | Endosulfan I        | 2.3   | U     |
| 60-57-1    | Dieldrin            | 4.4   | U     |
| 72-55-9    | 4,4'-DDE            | 4.4   | U     |
| 72-20-8    | Endrin              | 4.4 <del>0.18</del>                           | PJB U |
| 33213-65-9 | Endosulfan II       | 0.14  | PJB U |
| 72-54-8    | 4,4'-DDD            | 4.4   | U     |
| 1031-07-8  | Endosulfan sulfate  | 4.4 <del>0.22</del>                           | PJB U |
| 50-29-3    | 4,4'-DDT            | 4.4 <del>0.20</del>                           | PJB U |
| 72-43-5    | Methoxychlor        | 23  | U     |
| 53494-70-5 | Endrin ketone       | 4.4 <del>0.28</del>                           | PJB U |
| 7421-93-4  | Endrin aldehyde     | 4.4 <del>0.47</del>                           | PJB U |
| 5103-71-9  | alpha-Chlordane     | 2.3   | U     |
| 5103-74-2  | gamma-Chlordane     | 2.3   | U     |
| 8001-35-2  | Toxaphene           | 230   | U     |
| 12674-11-2 | Aroclor-1016        | 44  | U     |
| 11104-28-2 | Aroclor-1221        | 89  | U     |
| 11141-16-5 | Aroclor-1232        | 44  | U     |
| 53469-21-9 | Aroclor-1242        | 44  | U     |
| 12672-29-6 | Aroclor-1248        | 44  | U     |
| 11097-69-1 | Aroclor-1254        | 44  | U     |
| 11096-82-5 | Aroclor-1260        | 44  | U     |

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000763

1E  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNQ87

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.13

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: \_\_\_\_\_

% Moisture: 27 Decanted: (Y/N) N

Date Received: 11/08/00

Extraction: (Type) SONC

Date Extracted: 11/08/00

Concentrated Extract Volume: 5000 (uL)

Date Analyzed: 11/15/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.3

Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

| CAS NO.    | COMPOUND            | UG/KG | Q               |
|------------|---------------------|-------|-----------------|
| 319-84-6   | alpha-BHC           | 2.3   | U               |
| 319-85-7   | beta-BHC            | 2.3   | <del>PJ</del> U |
| 319-86-8   | delta-BHC           | 2.3   | U               |
| 58-89-9    | gamma-BHC (Lindane) | 2.3   | U               |
| 76-44-8    | Heptachlor          | 2.3   | <del>PJ</del> U |
| 309-00-2   | Aldrin              | 2.3   | U               |
| 1024-57-3  | Heptachlor epoxide  | 2.3   | <del>PJ</del> U |
| 959-98-8   | Endosulfan I        | 2.3   | U               |
| 60-57-1    | Dieldrin            | 4.5   | U               |
| 72-55-9    | 4,4'-DDE            | 4.5   | U               |
| 72-20-8    | Endrin              | 4.5   | U               |
| 33213-65-9 | Endosulfan II       | 4.5   | U               |
| 72-54-8    | 4,4'-DDD            | 4.5   | U               |
| 1031-07-8  | Endosulfan sulfate  | 4.5   | U               |
| 50-29-3    | 4,4'-DDT            | 4.5   | U               |
| 72-43-5    | Methoxychlor        | 2.3   | <del>PJ</del> U |
| 53494-70-5 | Endrin ketone       | 4.5   | U               |
| 7421-93-4  | Endrin aldehyde     | 4.5   | <del>PJ</del> U |
| 5103-71-9  | alpha-Chlordane     | 2.3   | U               |
| 5103-74-2  | gamma-Chlordane     | 2.3   | U               |
| 8001-35-2  | Toxaphene           | 230   | U               |
| 12674-11-2 | Aroclor-1016        | 45    | U               |
| 11104-28-2 | Aroclor-1221        | 92    | U               |
| 11141-16-5 | Aroclor-1232        | 45    | U               |
| 53469-21-9 | Aroclor-1242        | 45    | U               |
| 12672-29-6 | Aroclor-1248        | 45    | U               |
| 11097-69-1 | Aroclor-1254        | 45    | U               |
| 11096-82-5 | Aroclor-1260        | 45    | U               |

0.183

000766

1E  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNQ88

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.14

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: \_\_\_\_\_

% Moisture: 20 Decanted: (Y/N) N

Date Received: 11/08/00

Extraction: (Type) SONC

Date Extracted: 11/08/00

Concentrated Extract Volume: 5000 (uL)

Date Analyzed: 11/15/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.5

Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

Q

| CAS NO.    | COMPOUND            |     |      |       |
|------------|---------------------|-----|------|-------|
| 319-84-6   | alpha-BHC           | 2.1 | 0.20 | PJ U  |
| 319-85-7   | beta-BHC            | 2.1 | 1.9  | PJB U |
| 319-86-8   | delta-BHC           |     | 2.1  | U     |
| 58-89-9    | gamma-BHC (Lindane) |     | 2.1  | U     |
| 76-44-8    | Heptachlor          |     | 2.1  | U     |
| 309-00-2   | Aldrin              |     | 2.1  | U     |
| 1024-57-3  | Heptachlor epoxide  | 2.1 | 0.36 | PJ U  |
| 959-98-8   | Endosulfan I        |     | 2.1  | U     |
| 60-57-1    | Dieldrin            |     | 4.1  | U     |
| 72-55-9    | 4,4'-DDE            |     | 4.1  | U     |
| 72-20-8    | Endrin              |     | 4.1  | U     |
| 33213-65-9 | Endosulfan II       | 4.1 | 0.16 | PJ U  |
| 72-54-8    | 4,4'-DDD            | 4.1 | 0.20 | PJ U  |
| 1031-07-8  | Endosulfan sulfate  |     | 0.13 | J     |
| 50-29-3    | 4,4'-DDT            |     | 4.1  | U     |
| 72-43-5    | Methoxychlor        | 2.1 | 0.74 | PJB U |
| 53494-70-5 | Endrin ketone       |     | 4.1  | U     |
| 7421-93-4  | Endrin aldehyde     | 4.1 | 0.30 | PJB U |
| 5103-71-9  | alpha-Chlordane     | 2.1 | 0.11 | PJ U  |
| 5103-74-2  | gamma-Chlordane     |     | 2.1  | U     |
| 8001-35-2  | Toxaphene           |     | 210  | U     |
| 12674-11-2 | Aroclor-1016        |     | 41   | U     |
| 11104-28-2 | Aroclor-1221        |     | 84   | U     |
| 11141-16-5 | Aroclor-1232        |     | 41   | U     |
| 53469-21-9 | Aroclor-1242        |     | 41   | U     |
| 12672-29-6 | Aroclor-1248        |     | 41   | U     |
| 11097-69-1 | Aroclor-1254        |     | 41   | U     |
| 11096-82-5 | Aroclor-1260        |     | 41   | U     |

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00769

1E  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNQ89

Lab Name: AATSLA

Contract: 68-W0-0081

Lab Code: AATSLA

Case No.: 28706

SAS No.:

SDG No.: BNP05

Matrix: (soil/water) SOIL

Lab Sample ID: 45798.15

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: \_\_\_\_\_

% Moisture: 22 Decanted: (Y/N) N

Date Received: 11/08/00

Extraction: (Type) SONC

Date Extracted: 11/08/00

Concentrated Extract Volume: 5000 (uL)

Date Analyzed: 11/15/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.1

Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

| CAS NO.    | COMPOUND            | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) UG/KG | Q     |
|------------|---------------------|---|-------|
| 319-84-6   | alpha-BHC           | 2.2   | U     |
| 319-85-7   | beta-BHC            | 2.2   | U     |
| 319-86-8   | delta-BHC           | 2.2   | U     |
| 58-89-9    | gamma-BHC (Lindane) | 2.2   | U     |
| 76-44-8    | Heptachlor          | 2.2   | U     |
| 309-00-2   | Aldrin              | 2.2   | U     |
| 1024-57-3  | Heptachlor epoxide  | 2.2   | U     |
| 959-98-8   | Endosulfan I        | 2.2   | U     |
| 60-57-1    | Dieldrin            | 4.2   | U     |
| 72-55-9    | 4,4'-DDE            | 4.2   | U     |
| 72-20-8    | Endrin              | 4.2   | U     |
| 33213-65-9 | Endosulfan II       | 4.2   | U     |
| 72-54-8    | 4,4'-DDD            | 4.2   | U     |
| 1031-07-8  | Endosulfan sulfate  | 4.2   | U     |
| 50-29-3    | 4,4'-DDT            | 4.2   | U     |
| 72-43-5    | Methoxychlor        | 2.2 0.23                                      | FJB U |
| 53494-70-5 | Endrin ketone       | 4.2 0.034                                     | FJB U |
| 7421-93-4  | Endrin aldehyde     | 4.2   | U     |
| 5103-71-9  | alpha-Chlordane     | 2.2   | U     |
| 5103-74-2  | gamma-Chlordane     | 2.2   | U     |
| 8001-35-2  | Toxaphene           | 220   | U     |
| 12674-11-2 | Aroclor-1016        | 42  | U     |
| 11104-28-2 | Aroclor-1221        | 86  | U     |
| 11141-16-5 | Aroclor-1232        | 42  | U     |
| 53469-21-9 | Aroclor-1242        | 42  | U     |
| 12672-29-6 | Aroclor-1248        | 42  | U     |
| 11097-69-1 | Aroclor-1254        | 42  | U     |
| 11096-82-5 | Aroclor-1260        | 42  | U     |

0.125



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1E  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

BNQ91

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) SOIL Lab Sample ID: 45798.16  
 Sample wt/vol: 30.0 (g/mL) G Lab File ID: \_\_\_\_\_  
 % Moisture: 22 Decanted: (Y/N) N Date Received: 11/08/00  
 Extraction: (Type) SONC Date Extracted: 11/08/00  
 Concentrated Extract Volume: 5000 (uL) Date Analyzed: 11/15/00  
 Injection Volume: 2.0 (uL) Dilution Factor: 1.0  
 GPC Cleanup: (Y/N) Y pH: 7.3 Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

| CAS NO.    | COMPOUND            | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) UG/KG | Q     |
|------------|---------------------|---|-------|
| 319-84-6   | alpha-BHC           | 2.2   | U     |
| 319-85-7   | beta-BHC            | 2.2 <del>0.41</del>                           | PJB U |
| 319-86-8   | delta-BHC           | 2.2   | U     |
| 58-89-9    | gamma-BHC (Lindane) | 2.2   | U     |
| 76-44-8    | Heptachlor          | 2.2   | U     |
| 309-00-2   | Aldrin              | 2.2   | U     |
| 1024-57-3  | Heptachlor epoxide  | 2.2   | U     |
| 959-98-8   | Endosulfan I        | 2.2   | U     |
| 60-57-1    | Dieldrin            | 4.2   | U     |
| 72-55-9    | 4,4'-DDE            | 4.2   | U     |
| 72-20-8    | Endrin              | 4.2   | U     |
| 33213-65-9 | Endosulfan II       | 4.2   | U     |
| 72-54-8    | 4,4'-DDD            | 4.2   | U     |
| 1031-07-8  | Endosulfan sulfate  | 4.2   | U     |
| 50-29-3    | 4,4'-DDT            | 4.2 <del>0.074</del>                          | PJB U |
| 72-43-5    | Methoxychlor        | 2.2 <del>0.47</del>                           | PJB U |
| 53494-70-5 | Endrin ketone       | 4.2   | U     |
| 7421-93-4  | Endrin aldehyde     | 4.2   | U     |
| 5103-71-9  | alpha-Chlordane     | 2.2   | U     |
| 5103-74-2  | gamma-Chlordane     | 2.2   | U     |
| 8001-35-2  | Toxaphene           | 220   | U     |
| 12674-11-2 | Aroclor-1016        | 42  | U     |
| 11104-28-2 | Aroclor-1221        | 86  | U     |
| 11141-16-5 | Aroclor-1232        | 42  | U     |
| 53469-21-9 | Aroclor-1242        | 42  | U     |
| 12672-29-6 | Aroclor-1248        | 42  | U     |
| 11097-69-1 | Aroclor-1254        | 42  | U     |
| 11096-82-5 | Aroclor-1260        | 42  | U     |

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1E  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

000775  
EPA SAMPLE NO.

BNQ92 FE

Lab Name: AATSLA Contract: 68-W0-0081  
 Lab Code: AATSLA Case No.: 28706 SAS No.: SDG No.: BNP05  
 Matrix: (soil/water) WATER Lab Sample ID: 45798.17  
 Sample wt/vol: 1000 (g/mL) ML Lab File ID: \_\_\_\_\_  
 % Moisture: \_\_\_\_\_ Decanted: (Y/N) \_\_\_\_\_ Date Received: 11/08/00  
 Extraction: (Type) CONT Date Extracted: 11/13/00  
 Concentrated Extract Volume: 10000 (uL) Date Analyzed: 11/17/00  
 Injection Volume: 2.0 (uL) Dilution Factor: 1.0  
 GPC Cleanup: (Y/N) N pH: 7.0 Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L Q

| CAS NO.    | COMPOUND            | UG/L   | Q |
|------------|---------------------|--------|---|
| 319-84-6   | alpha-BHC           | 0.050  | U |
| 319-85-7   | beta-BHC            | 0.050  | U |
| 319-86-8   | delta-BHC           | 0.050  | U |
| 58-89-9    | gamma-BHC (Lindane) | 0.050  | U |
| 76-44-8    | Heptachlor          | 0.050  | U |
| 309-00-2   | Aldrin              | 0.050  | U |
| 1024-57-3  | Heptachlor epoxide  | 0.050  | U |
| 959-98-8   | Endosulfan I        | 0.050  | U |
| 60-57-1    | Dieldrin            | 0.10   | U |
| 72-55-9    | 4,4'-DDE            | 0.10   | U |
| 72-20-8    | Endrin              | 0.10   | U |
| 33213-65-9 | Endosulfan II       | 0.10   | U |
| 72-54-8    | 4,4'-DDD            | 0.10   | U |
| 1031-07-8  | Endosulfan sulfate  | 0.10   | U |
| 50-29-3    | 4,4'-DDT            | 0.10   | U |
| 72-43-5    | Methoxychlor        | 0.014  | J |
| 53494-70-5 | Endrin ketone       | 0.10   | U |
| 7421-93-4  | Endrin aldehyde     | 0.032  | J |
| 5103-71-9  | alpha-Chlordane     | 0.0052 | J |
| 5103-74-2  | gamma-Chlordane     | 0.050  | U |
| 8001-35-2  | Toxaphene           | 5.0    | U |
| 12674-11-2 | Aroclor-1016        | 1.0    | U |
| 11104-28-2 | Aroclor-1221        | 2.0    | U |
| 11141-16-5 | Aroclor-1232        | 1.0    | U |
| 53469-21-9 | Aroclor-1242        | 1.0    | U |
| 12672-29-6 | Aroclor-1248        | 1.0    | U |
| 11097-69-1 | Aroclor-1254        | 1.0    | U |
| 11096-82-5 | Aroclor-1260        | 1.0    | U |

0187

REFERENCE NO. 14

RECORD OF COMMUNICATION

TO: YUNRU YANG

FROM: JANET TROTTER  
Region 2, ESAT/RSCC

DATE: Feb. 21, 2001

SUBJECT: QUALITY ASSURED DATA

=====

MESSAGE: PLEASE SIGN BELOW IN ACKNOWLEDGEMENT OF RECEIPT OF THE FOLLOWING AND RETURN ONE COPY OF THIS RECORD OF COMMUNICATION TO THE RSCC - REGION II.

① LACKAWANNA Foundry 28706 Chemical Inorg 14S/1W

\_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

=====

REPLY BY: \_\_\_\_\_

\_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

SIGNATURE: *Janet Trotter* DATE: 2/22/01

DATE RECEIVED BY RSCC: \_\_\_\_\_

# RECORD OF COMMUNICATION

## REGIONAL SAMPLE CONTROL CENTER

DATE: Nov. 24, 2000  
 SUBJECT: CLP Data Package for Quality Assurance Review  
 FROM: RSCC / ESAT  
 TO: Hanif Sheikh, Hazardous Waste Support Section

FEB 02 2001

*Attached is the following INORGANIC Data Package to be reviewed for Quality Assurance*

|                                  |                              |
|----------------------------------|------------------------------|
| SITE <u>LACKAWANNA FOUNDRY</u>   | CASE# <u>28706</u>           |
| SAMPLER <u>SAT</u>               | #SAMPLES <u>14</u>           |
| PHASE <u>BZ</u>                  | MATRIX <u>SOIL</u>           |
| LAB <u>CHEMED</u>                | FRACTION <u>TAL &amp; CN</u> |
| TURN-AROUND-TIME <u>21 DAYS</u>  | SITE SPILL # <u>MY</u>       |
| CERCLIS ID # <u>NYSFNO204209</u> | MATRIX <u>WATER</u>          |

### REGION II RSCC DATA TRANSFER LOG

| Relinquished By     |                 | Received By         |                 |
|---------------------|-----------------|---------------------|-----------------|
| Signature           | Date/Time       | Signature           | Date/Time       |
|                     |                 | <u>John Bulich</u>  | <u>11-22-00</u> |
| <u>John Bulich</u>  | <u>11-24-00</u> | <u>CMAA</u>         | <u>01/29/01</u> |
| <u>CMAA</u>         | <u>1-31-01</u>  | <u>J Bulich</u>     | <u>1-31-01</u>  |
| <u>J Bulich</u>     | <u>1-31-01</u>  | <u>Hanif Sheikh</u> | <u>01-31-01</u> |
| <u>Hanif Sheikh</u> | <u>02-02-01</u> | <u>J Bulich</u>     | <u>2-2-01</u>   |
|                     |                 |                     |                 |
|                     |                 |                     |                 |
|                     |                 |                     |                 |
|                     |                 |                     |                 |
|                     |                 |                     |                 |

USEPA DATA TRANSFER LOG INSTRUCTIONS

The DATA TRANSFER LOG FORM is to be used to document the CUSTODY of original documents transferred to EPA contractors for data validation. Anytime original documents are transferred from one individual to another, the transfer must be documented on the DATA TRANSFER LOG FORM. The completion of this form will assure that a history of the samples is maintained from the time they were collected to the time when the data is presented as evidence in court proceedings.

The documents are considered to be under custody if:

1. They are in your possession, or
2. They are in your view, after being in your possession, or
3. They were in your possession and you locked them up, or
4. They are in a designated secure area.

PLEASE COMPLETE THE FORM IN THE FOLLOWING MANNER:

The DATA TRANSFER LOG is to be completed in the same manner a chain of custody record is completed. Indelible ink must be used for each entry.

The person who initially opens the package containing the data must print their name, sign their name, and enter the date and time received in the "Received By" column. When the data is transferred to another individual, the "Relinquished By" column must also be completed by the individual who originally had possession of the documents. The individual receiving the documents must complete the next line in the "Received By" column. Each individual involved in the transfer of the data must enter their information on the form.

|             |           |           |                   |        |    |
|-------------|-----------|-----------|-------------------|--------|----|
| Case #      | 28706     | Site:     | LACKAWANA FOUNDRY | Soil:  | 14 |
| SDG #       | MBNJ51    | Lab:      | CHEMED            | Water: | 01 |
| Contractor: | SAT/BZ/MY | Reviewer: | C.M. Alaimo/ESAT  | Other: | 0  |

**A.2.1. Validation flags-** The following flags have been applied in red by the data validator and must be considered by the data user.

**J -** This flag indicates the result qualified as estimated.

**Red-Line -** A red-line drawn through a sample result indicates unusable value. The red-lined data are known to contain significant errors based on documented information and must not be used by the data user.

**Fully Usable Data-** The results that do not carry "J" or "red-line" are fully usable.

**Contractual Qualifiers-** The legend of contractual qualifiers applied by the lab on Form I's is found on page B-20 of SOW ILM01.0

**A.2.2. The data assessment is given below and on the attached sheets.**

This package consists of fourteen (14) soil and one (01) aqueous samples collected from the Lackawana Foundry site on 11/07/00 for AL metals plus CN analysis. All holding times were met. QC was performed on sample MBNJ51. No trip report was available, field QC could not be determined.

SDG #MBNJ51

**CRDL**

For both Hg runs, the CRA's % (125.0%) recoveries were between 121-150% requiring estimation of all positive results within the effected range.

J--> Hg in MBNJ54 & MBNJ55.

**ICP SERIAL DILUTION**

For Cu & K, the %D were greater than 10% when the initial sample values were greater than 10xIDL, therefore all associated data greater than or equal to the CRDL (or  $\geq 10xIDL$  when  $10xIDL > CRDL$ ).

J--> Cu in MBNJ51--> 60, MBNJ62, MBNJ63, MBNJ64 & MBNJ66

J--> K in MBNJ59, MBNJ60, MBNJ62, MBNJ63 & MBNJ66.

**A.2.3. Contract-Problems/Non-Compliance**

For field blank MBNJ67, the lead result was reported without a concentration qualifier. According to the raw data, the value was less than the CRDL and required a "B" qualifier. The validator made the necessary corrections.

MMB Reviewer: HS Date: 02/01/01  
Signature

Contractor Reviewer: CM Alaimo Date: 01/30/01  
Signature

Verified by: \_\_\_\_\_ Date: \_\_\_\_\_  
Signature

p.4

**CLP DATA ASSESSMENT SUMMARY FORM**

**TYPE OF REVIEW:** INORGANIC      **DATE:** 01/30/01      **CASE #:** 28706

**SITE:** Lackawana Foundry      **LAB NAME:** CHEMED

**REVIEWER'S INITIALS:** CMA      **NUMBER OF SAMPLES:** WATER 01  
SOIL 14

**ANALYTES REJECTED DUE TO EXCEEDING REVIEW CRITERIA:\***

|            | Holding Time | CRDL | Prep Blank | Field Blank | Interference | Spike Recov. | Tot/Diiso | LCS | Serial Dil. | % TS | Total Analytes | Rejections |
|------------|--------------|------|------------|-------------|--------------|--------------|-----------|-----|-------------|------|----------------|------------|
| ICP        |              |      |            |             |              |              |           |     |             |      | 330            | 0          |
| Flame AA   |              |      |            |             |              |              |           |     |             |      |                |            |
| Furnace AA |              |      |            |             |              |              |           |     |             |      |                |            |
| Mercury    |              |      |            |             |              |              |           |     |             |      | 15             | 0          |
| Cyanide    |              |      |            |             |              |              |           |     |             |      | 15             | 0          |
| Total      |              |      |            |             |              |              |           |     |             |      | 360            | 0          |

**ANALYTES FLAGGED AS ESTIMATED (J) DUE TO EXCEEDING CRITERIA FOR:\***

|            | Holding Time | CRDL | Prep Blank | Tot/Diiso | Preservation | Spike Recov. | Dup Lab | Dup Field | Det. Limit | LCS | Serial Dilution | %TS | Total Analytes | Estimation |
|------------|--------------|------|------------|-----------|--------------|--------------|---------|-----------|------------|-----|-----------------|-----|----------------|------------|
| ICP        |              | 2    |            |           |              |              |         |           |            |     |                 | 19  | 330            | 21         |
| Flame AA   |              |      |            |           |              |              |         |           |            |     |                 |     |                |            |
| Furnace AA |              |      |            |           |              |              |         |           |            |     |                 |     |                |            |
| Mercury    |              |      |            |           |              |              |         |           |            |     |                 |     | 15             | 0          |
| Cyanide    |              |      |            |           |              |              |         |           |            |     |                 |     | 15             | 0          |
| Total      |              | 2    |            |           |              |              |         |           |            |     |                 | 19  | 360            | 21         |

Note: Asterisk (\*) indicates additional exceedances of review criteria.



**Region II  
INORGANIC REGIONAL DATA ASSESSMENT**

**CASE NO.:** 28706

**SITE:** LACKAWANA FOUNDRY

**LABORATORY:** CHEMED

**NO. OF SAMPLES/MATRIX:** 01 WATER, 14 SOIL

**SDG#:** MBNJ51

**REVIEWER (IF NOT ESD):**

**SOW#:** ILM04.1

**REVIEWER'S NAME:** C. ALAIMO

**DPO:** ACTION FYI

**COMPLETION DATE:** 01/30/01

**DATA ASSESSMENT SUMMARY**

|                     | ICP | AA  | MERCURY | CYANIDE |
|---------------------|-----|-----|---------|---------|
| Holding Times       | O   | N/A | O       | O       |
| Calibration         | O   | N/A | O       | O       |
| Blanks              | O   | N/A | O       | O       |
| ICS                 | O   | N/A | N/A     | N/A     |
| LCS                 | O   | N/A | N/A     | N/A     |
| Duplicate Analysis  | O   | N/A | O       | O       |
| Matrix Spike        | O   | N/A | O       | O       |
| MSA                 | N/A | N/A | N/A     | N/A     |
| Serial Dilution     | O   | N/A | N/A     | N/A     |
| Sample Verification | O   | N/A | O       | O       |
| Other QC            | O   | N/A | O       | O       |
| Overall Assessment  | O   | N/A | O       | O       |

O = Data has no problems/or qualified due to minor problems.

M = Data qualified due to major problems.

Z = Data unacceptable.

X = Problems, but do not affect data.

**ACTION ITEMS:**

**AREAS OF CONCERN:**

**NOTABLE PERFORMANCE:**

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Title: Evaluation of Metals Data for the Contract Laboratory Program Appendix A.3: Contract Non-Compliance (SMO Report)

Date: Jan. 1992 Number: HW-2 Revision: 11

CONTRACT NON-COMPLIANCE (SMO REPORT)

Regional Review of Uncontrolled Hazardous Waste Site Contract Laboratory Data Package

CASE NO. 28706

The hardcopied (laboratory name) Inorganic data package received at Region II has been reviewed and the quality assurance and performance data summarized. The data reviewed included: SMO Sample No.:

Conc. & Matrix:

Contract No. ( ) requires that specific analytical work be done and that associated reports be provided by the contractor to the Regions, EMSL-LV, and SMO. The general criteria used to determine the performance were based on an examination of:

- Data Completeness - Duplicate Analysis Results - Matrix Spike Results - Blank Analysis Results - Calibration Standards Results - MSA Results

Items of non-compliance with the above contract are described below.

Comments: N/A

Reviewer's Initial CMA

Date 01/30/07

Title: Evaluation of Metals Data for the  
Contract Laboratory Program  
Appendix A.4: Mailing List for Data Reviewers

Date: Jan. 1992  
Number: HW-2  
Revision: 11

DPO MAILING LIST FOR DATA REVIEWERS

- |  |  |
|--|--|
| <p>1. USEPA Region I (ESD)<br/>60 Westview Street<br/>Lexington, MA 02173<br/>Deb Szaro<br/>(617) 861-4312<br/>CT, ME, MA, NH, RI, VT<br/>CAA, Resource Analysts, York,<br/>E3I, Skinner, TMA</p>  | <p>2. USEPA Region II ESD<br/>Woodbridge Avenue<br/>Edison, NJ 08837<br/>Lisa Gatto Vidulich<br/>(201) 321-6676<br/>NJ, NY, PR, VI<br/>Century, Chemtech, US Test, Nanco<br/>ETC, Gadson, EMS, Galson, ICM</p>   |
| <p>3. USEPA Region III (CRL)<br/>839 Bestgate Road<br/>Annapolis, MD 21401<br/>Chuck Sands<br/>(301) 266-9180<br/>DE, MD, PA, VA, WV, DC<br/>Center, Hitman, JTC, MACK, VERSAR,<br/>ITAS, Weston, MMES, EA Engineering,<br/>Subject Tech., KEYPA</p>   | <p>4. USEPA Region IV (ESD)<br/>Analytical Support Branch<br/>College Station Road<br/>Athens, GA 30613<br/>Tom Bennett, Jr.<br/>(404) 546-3112<br/>AL, FL, GA, KY, MS, NC, SC, TN<br/>CompuChem, EPS, ESE, PBS&amp;J,<br/>Triangle Labs</p>                                     |
| <p>5. USEPA Region V (ESD)<br/>536 South Clark Street<br/>Tenth Floor, CRL<br/>Chicago, IL 60605<br/>Pat Churilla<br/>(312) 353-9087<br/>IL, IN, MI, MN, OH, WI<br/>NLE, TAI/ENG</p>   | <p>6. USEPA Region VI (ESD)<br/>Monterey Park Plaza, Bldg. C<br/>6608 Hornwood Drive<br/>Houston, TX 77074<br/>David Stockton<br/>(713) 953-3425<br/>AR, LA, NM, TX, OK<br/>ANACON, RADIANT, SPECS, EIS, Glochem<br/>Research, Inc., SPL Inc., SWRI,<br/>Allied, KEYTX, EIRA</p> |
| <p>7. USEPA Region VII Laboratory<br/>25 Funston Road<br/>Kansas City, KS 66115<br/>Debra Morey<br/>(913) 236-3881<br/>IO, KS, NB, MO<br/>Wilson, Kansas City Scientific<br/>Enterprises, Eagle Picher</p>   | <p>8. USEPA Region VIII Laboratory<br/>Box 25366<br/>Denver Federal Center<br/>Lakewood, CO 80225<br/>Eva Hoffman<br/>(303) 236-7371<br/>CO, ND, SD, UT, WY, MT<br/>ACCU, CSMRI, RMAL, Data Chem., Centref</p>   |
| <p>9. USEPA Region XI (ESD)<br/>QA Management Section<br/>215 Fremont Street<br/>San Francisco, CA 94105<br/>Kent Kitchingman<br/>(415) 974-0924<br/>AZ, CA, HI, NV, American Samoa,<br/>Guam Trust Territories of Pacific<br/>Islands, Wake Island<br/>ALI, CAL Weston, S-Cubed, IT_CA,<br/>Vegas</p> | <p>10. USEPA Region X Laboratory<br/>P.O. Box 549<br/>Manchester, WA 98353<br/>Gerald Muth<br/>(206) 442-0370<br/>AK, ID, OR, WA<br/>Laucke Testing Labs, Century Testing<br/>Labs (For VOA Only), Weyerhaeuser Co.,<br/>Columbia Testing, Silver Valley</p>                     |
| <p>11. Caria Dempsey - (OS-230)<br/>USEPA<br/>401 "M" Street S.W.<br/>Washington, DC 20460<br/>FTS 382-5746</p>  | <p>12. Edward Kantor<br/>USEPA<br/>EMSL-LV<br/>944 E. Harmon Avenue<br/>Box 93478<br/>Las Vegas, NV 89119</p>  |
| <p>13. Sample Management Office<br/>Viar and Company<br/>P.O. Box 818<br/>Alexandria, VA 22313</p>   |  |

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CASE# 280 28106  
SDG: MONTJ51

Evaluation of Metals Data for the Contract Laboratory Program (CLP)

based on

SOP. 3/90

(SOP Revision XI)

PREPARED BY: Harif Sheikh DATE: 1-30-92  
Harif Sheikh, Quality Assurance Chemist  
Toxic and Hazardous Waste Section

APPROVED BY: Kevin Kubik DATE: 1-30-92  
Kevin Kubik, Chief  
Toxic and Hazardous Waste Section

APPROVED BY: Robert Runyon DATE: 1/31/92  
Robert Runyon, Chief  
Monitoring Management Branch

Title: Evaluation of Metals Data for the  
Contract Laboratory Program

Date: Jan. 1992  
Number: HW-2  
Revision: 11

## 1.0 Scope

1.1 This procedure is applicable to inorganic data obtained from contractor laboratories working for Hazardous Waste Site Contract Laboratory Program (CLP).

1.2 The data validation is based upon analytical and quality assurance requirements specified in Statement of Work (SOW) 3/90 .

2.0 Responsibilities - Data reviewers will complete the following tasks as assigned by the Data Review Coordinator:

2.1. For a total review:

2.1.1 Data Assessment - "Total Review-Inorganics" Checklist Appendix (A.1).  
The reviewer must answer every question on the checklist.

2.1.2 Data Assessment - Data Assessment Narrative (Appendix A.2)  
The answer on the checklist must match the action in the narrative (appendix A.2) and on Form I's. Do not use pencil to write the narrative.

2.1.3 Contract Non-Compliance - SMO Report (Appendix A.3)  
This report is to be completed only when a serious contract violation is encountered, or upon the request of the Data Validation Task Monitor, or Technical Project Officer (TPO). Forward 5 copies: one each for internal files, appropriate Regional TPO, Sample Management Office (SMO) and last two addresses of Mailing List for Data Reviewers (Appendix A.4). In other cases, all contract violations should be appended to the end of the Data Assessment Narrative (Sec. A.2.2).

2.1.4 CLP Data Assessment Summary Forms

2.1.4.1 Appendix A.5

Fill in the total number of analytes analyzed by different analyses and the number of analytes rejected or flagged as estimated due to corresponding quality control criteria. Place an "X" in boxes where analyses were not performed, or criteria do not apply.

2.1.4.2 Appendix A.6

Data reviewer is also required to fill out Inorganic Regional Data Assessment form (Appendix A.7) provided by EPA Headquarters. Codes listed on the form will be used to describe the Data Assessment Summary.

Title: Evaluation of Metals Data for the  
Contract Laboratory Program

Date: Jan. 1992  
Number: HW-2  
Revision: 11

- 
- 2.1.5 Data Review Log: It is recommended that each data reviewer should maintain a log of the reviews completed to include:
- a. date of start of case review
  - b. date of completion of case review
  - c. site
  - d. case number
  - e. contract laboratory
  - f. number of samples
  - g. matrix
  - h. hours worked
  - i. reviewer's initials
- 2.1.6 Telephone Record Log - the data reviewer should enter the bare facts of inquiry, before initiating any phone conversation with CLP laboratory. After the case review has been completed, mail white copy of Telephone Record Log to the laboratory and pink copy to SMO. File yellow copy in the Telephone Record Log folder, and attach a xerox copy of the Telephone Record Log to the completed Data Assessment Narrative (Appendix A.2).
- 2.1.7 Forwarded Paperwork
- 2.1.7.1 Upon completion of review, the following are to be forwarded to the Regional Sample Control Center (RSOC) located in the Surveillance and Monitoring Branch:
- a. data package
  - b. completed data assessment checklist (Appendix A.1, original)
  - c. SMO Contract Compliance Screening (CCS)
  - d. Record of Communication (copy)
  - e. CLP Reanalysis Request/Approval Record (original + 3 copies)
  - f. Appendix A.6 (original).
- 2.1.7.2 Forward 2 copies of completed Data Assessment Narrative (Appendix A.2) along with 2 copies of the Inorganic Data Assessment Form (Appendix A.6) and Telephone Record Log, if any, : one each for appropriate Regional TFO, and the other one to EPA EMSL office in Las Vegas. The addresses of TFOs and EPA office in Las Vegas are given in Appendix A-4.
- 2.1.8 Filed Paperwork - Upon completion of review, the following are to be filed within MMB files:
- a. Two copies of completed Data Assessment Narrative (Appendix A.2) each carrying Appendix A.6.
  - b. Telephone Record Log (copy)
  - c. SMO Report (copy Appendix A-3)
  - d. CLP Reanalysis Request/Approval Record (copy)
- p.11

Title: Evaluation of Metals Data for the  
Contract Laboratory Program

Date: Jan. 1992  
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- 3.0 Data Completeness  
Each data package is checked by a Regional Sample Control Coordinator (RSSC) for completeness. A data package is assumed to be complete when all the deliverables required under the contract are present. If a data package is incomplete, the RSSC would call the laboratory for missing document(s). If the laboratory does not respond within a week, SMO and MMB coordinator of Region II will be notified.
- 4.0 Rejection of Data - All values determined to be unacceptable on the Inorganic Analysis Data Sheet (Form I) must be lined over with a red pencil. As soon as any review criteria causes data to be rejected, that data can be eliminated from any further review or consideration.
- 5.0 Acceptance Criteria - In order that reviews be consistent among reviewers, acceptance criteria as stated in Appendix A.1 (pages 4-25) should be used. Additional guidance can be found in the National Inorganic Functional Guidelines of October 1, 1989.
- 6.0 SMO Contract Compliance Screening (CCS) - This is intended to aid reviewer in locating any problems, both corrected and uncorrected. However, the validation should be carried out even if CCS is not present. Resubmittals received from laboratory in response to CCS must be used by the reviewer.
- 7.0 Request for Reanalysis - Data reviewers must note all items of contract non-compliance within Data Assessment Narrative. If holding times and sample storage times have not been exceeded, TPO may request reanalysis if items of non-compliance are critical to data assessment. Requests are to be made on "CLP Re-Analysis Request/Approval Record".
- 8.0 Record of Communication - Provided by the Regional Sample Control Center (RSSC) to indicate which data packages have been received and are ready to be reviewed.
- 9.0 Rounding off numbers - The data reviewer will follow the standard practice.
- p. 12

Title: Evaluation of Metals Data for the  
 Contract Laboratory Program  
 Appendix A.1: Data Assessment - Contract  
 Compliance (Total Review)

Date: Jan. 1992  
 Number: HW-2  
 Revision: 11

|   | YES                                 | NO                                  | N/A                      |
|---|-------------------------------------|-------------------------------------|--------------------------|
| A.1.1 <u>Contract Compliance Screening Report (CCS)</u> - Present?                        | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/> |
| <u>ACTION:</u> If no, contact RSCC.   |                                     |                                     |                          |
| A.1.2 <u>Record of Communication (from RSCC)</u> - Present?                               | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/> |
| <u>ACTION:</u> If no, request from RSCC.  |                                     |                                     |                          |
| A.1.3 <u>Trip Report</u> - Present and complete?  | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| <u>ACTION:</u> If no, contact RSCC for trip report.                                       |                                     |                                     |                          |
| A.1.4 <u>Sample Traffic Report</u> - Present?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/> |
| Legible?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/> |
| <u>ACTION:</u> If no, request from Regional Sample Control Center (RSCC).                 |                                     |                                     |                          |
| A.1.5 <u>Cover Page</u> - Present?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/> |
| Is cover page properly filled in and signed by the lab manager or the manager's designee? | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/> |
| <u>ACTION:</u> If no, prepare Telephone Record Log, and contact laboratory.               |                                     |                                     |                          |
| Do numbers of samples correspond to numbers on Record of Communication?                   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/> |
| Do sample numbers on cover page agree with sample numbers on:                             |                                     |                                     |                          |
| (a) Traffic Report Sheet?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/> |
| (b) Form I's?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/> |
| <u>ACTION:</u> If no for any of the above, contact RSCC for clarification.                |                                     |                                     |                          |



STANDARD OPERATING PROCEDURE

Title: Evaluation of Metals Data for the  
 Contract Laboratory Program  
 Appendix A.1: Data Assessment - Contract  
 Compliance (Total Review)

Date: Jan. 1992  
 Number: HW-2  
 Revision: 11

A.1.6 Form I to IX Yes No N/A

A.1.6.1 Are all the Form I through Form IX labeled with:

|                  |                                     |   |   |
|------------------|-------------------------------------|---|---|
| Laboratory name? | <input checked="" type="checkbox"/> | — | — |
| Case/SAS number? | <input checked="" type="checkbox"/> | — | — |
| EPA sample No.?  | <input checked="" type="checkbox"/> | — | — |
| SDG No.?         | <input checked="" type="checkbox"/> | — | — |
| Contract No.?    | <input checked="" type="checkbox"/> | — | — |
| Correct units?   | <input checked="" type="checkbox"/> | — | — |
| Matrix?          | <input checked="" type="checkbox"/> | — | — |

ACTION: If no for any of the above, note under Contract Problem/Non-Compliance section of the "Data Assessment Narrative".

A.1.6.2 Do any computation/transcription errors exceed 10% of reported values on Forms I-IX for:

(NOTE: Check all forms against raw data.)

|  |                                     |   |   |
|--|-------------------------------------|---|---|
| (a) all analytes analyzed by ICP?      | <input checked="" type="checkbox"/> | — | — |
| (b) all analytes analyzed by GFAA?     | <input checked="" type="checkbox"/> | — | — |
| (c) all analytes analyzed by AA Flame? | <input checked="" type="checkbox"/> | — | — |
| (d) Mercury?                           | <input checked="" type="checkbox"/> | — | — |
| (e) Cyanide?                           | <input checked="" type="checkbox"/> | — | — |

ACTION: If yes, prepare Telephone Log, contact laboratory for corrected data and correct errors with red pencil and initial.

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|         |  | YES                                 | NO                       | N/A                                 |
|---------|--|-------------------------------------|--------------------------|-------------------------------------|
| A.1.7   | <u>Raw Data</u>  |                                     |                          |                                     |
| A.1.7.1 | Digestion Log* for flame AA/ICP (Form XIII) present?                       | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/>            |
|         | Digestion Log for furnace AA Form XIII present?                            | <input type="checkbox"/>            | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
|         | Distillation Log for mercury Form XIII present?                            | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/>            |
|         | Distillation Log for cyanides Form XIII present?                           | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/>            |
|         | Are pH values (pH<2 for all metals, pH>12 for cyanide) present?            | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/>            |
|         | *Weights, dilutions and volumes used to obtain values.                     |                                     |                          |                                     |
|         | Percent solids calculation present for soils/sediments?                    | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/>            |
|         | Are preparation dates present on sample preparation logs/bench sheets?     | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/>            |
| A.1.7.2 | Measurement read out record present?                                       |                                     |                          |                                     |
|         | ICP  | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/>            |
|         | Flame AA   | <input type="checkbox"/>            | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
|         | Furnace AA   | <input type="checkbox"/>            | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
|         | Mercury  | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/>            |
|         | Cyanides   | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/>            |
| A.1.7.3 | Are all raw data to support all sample analyses and QC operations present? | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/>            |
|         | Legible?   | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/>            |
|         | Properly Labeled?  | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/>            |

**ACTION:** If no for any of the above questions in sections A.1.7.1 through A.1.7.3, write Telephone Record Log and contact laboratory for resubmittals.

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|         |   | YES                                 | NO                                  | N/A |
|---------|---|-------------------------------------|-------------------------------------|-----|
| A.1.8   | <u>Holding Times</u> - (aqueous and soil samples )  |                                     |                                     |     |
|         | (Examine sample traffic reports and digestion/distillation logs.)   |                                     |                                     |     |
|         | Mercury analysis (28 days) . . . . . exceeded?  | —                                   | <input checked="" type="checkbox"/> | —   |
|         | Cyanide distillation (14 days) . . . . . exceeded?  | —                                   | <input checked="" type="checkbox"/> | —   |
|         | Other Metals analysis (6 months) . . . . . exceeded?  | —                                   | <input checked="" type="checkbox"/> | —   |
|         | <u>NOTE:</u> Prepare a list of all samples and analytes for which holding times have been exceeded. Specify the number of days from date of collection to the date of preparation (from raw data). Attach to checklist. |                                     |                                     |     |
|         | <u>ACTION:</u> If yes, reject (red-line) values less than Instrument Detection Limit (IDL) and flag as estimated (J) the values above IDL even though sample(s) was preserved properly.                                 |                                     |                                     |     |
| A.1.8.2 | Is pH of aqueous samples for:   |                                     |                                     |     |
|         | Metals Analysis >2?   | —                                   | <input checked="" type="checkbox"/> | —   |
|         | Cyanides Analysis <12?  | —                                   | <input checked="" type="checkbox"/> | —   |
|         | <u>Action:</u> If yes, flag the associated metals and cyanides data as estimated.   |                                     |                                     |     |
| A.1.9   | <u>Form I (Final Data)</u>  |                                     |                                     |     |
| A.1.9.1 | Are all Form I's present and complete?  | <input checked="" type="checkbox"/> | —                                   | —   |
|         | <u>ACTION:</u> If no, prepare telephone record log and contact laboratory for submittal.  |                                     |                                     |     |
| A.1.9.2 | Are correct units (ug/l for waters and mg/kg for soils) indicated on Form I's?  | <input checked="" type="checkbox"/> | —                                   | —   |
|         | Are soil sample results for each parameter corrected for percent solids?  | <input checked="" type="checkbox"/> | —                                   | —   |
|         | Are all "less than IDL" values properly coded with "U"?   | <input checked="" type="checkbox"/> | —                                   | —   |

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|  | YES                                 | NO                                  | N/A                                 |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| Are the correct concentration qualifiers used with final data?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| <u>ACTION:</u> If no for any of the above, prepare Telephone Record Log, and contact laboratory for corrected data.                |                                     |                                     |                                     |
| A.1.9.3 Are EPA sample # s and corresponding laboratory sample ID # s the same as on the Cover Page, Form I's and in the raw data? | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| Was a brief physical description of samples given on Form I's?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| Was the dilution of any sample diluted beyond the requirements of the contract noted on Form I or Form XIV?                        | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/>            |
| <u>ACTION:</u> If no for any of the above, note under Contract-Problem/Non-Compliance of the "Data Assessment Narrative".          |                                     |                                     |                                     |
| A.1.10 <u>Calibration</u>  |                                     |                                     |                                     |
| A.1.10.1 Is record of at least 2 point calibration present for ICP analysis?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| Is record of 5 point calibration present for Hg analysis?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| Is record of 4 point calibration present for:  |                                     |                                     |                                     |
| Flame AA?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Furnace AA?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Cyanides?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| Is one calibration standard at the CRDL level for all AA and cyanides analyses?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |

ACTION: If no for any of the above, write in the Contract Problem/Non-Compliance section of the "Data Assessment Narrative".

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|  | YES | NO                                  | N/A                                 |
|--|-----|-------------------------------------|-------------------------------------|
| A.1.10.2 Is correlation coefficient less than 0.995 for: |     |                                     |                                     |
| Mercury Analysis?  | —   | <input checked="" type="checkbox"/> | —                                   |
| Cyanide Analysis?  | —   | <input checked="" type="checkbox"/> | —                                   |
| Atomic Absorption Analysis?                              | —   | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |

**ACTION:** If yes, flag the associated data as estimated.

**NOTE:** The data validator shall calculate the correlation coefficient using concentrations of the standards and the corresponding instrument response ( e.g. absorbance, peak area, peak height, etc.).

|          |   |                                     |   |   |
|----------|---|-------------------------------------|---|---|
| A.1.10.3 | In the instance where less than 4 standards are measured in absorbance (or peak area, peak height, etc.) mode, are the remaining standards analyzed in concentration mode immediately after calibration within $\pm 10\%$ of the true values? | <input checked="" type="checkbox"/> | — | — |
|----------|---|-------------------------------------|---|---|

**ACTION:** If no, flag the associated data as estimated if standards are not within  $\pm 10\%$  of true values. Do not flag the data as estimated in linear range indicated by good recovery of standard(s).

**A.1.11 Form II A (Initial and Continuing Calibration Verification)-**

|          |  |                                     |   |                                     |
|----------|--|-------------------------------------|---|-------------------------------------|
| A.1.11.1 | Present and complete for every metal and cyanide?                            | <input checked="" type="checkbox"/> | — | —                                   |
|          | Present and complete for AA and ICP when both are used for the same analyte? | <input type="checkbox"/>            | — | <input checked="" type="checkbox"/> |

**ACTION:** If no for any of the above, prepare Telephone Record Log and contact laboratory.

|          |  |                                     |   |   |
|----------|--|-------------------------------------|---|---|
| A.1.11.2 | Circle on each Form IIA all percent recoveries that are outside the contract windows.<br>Are all calibration standards (initial and continuing) within control limits: |                                     |   |   |
|          | Metals- 90-110%R?  | <input checked="" type="checkbox"/> | — | — |
|          | Hg - 80-120%R?   | <input checked="" type="checkbox"/> | — | — |
|          | Cyanides- 85-115%R?  | <input checked="" type="checkbox"/> | — | — |

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|  | YES | NO | N/A |
|--|-----|----|-----|
| <b>ACTION:</b> Flag as estimated (J) all positive data (not flagged with a "U") analyzed between a calibration standard with %R between 75-89% (65-79% for Hg; 70-84% for CN) or 111-125% (121-135% for Hg; 116-130% for CN) recovery and nearest good calibration standard. Qualify results <IDL as estimated (U) if the ICV or CCV %R is 75-89% (CN, 70-84% ; HG, 65-79%). Reject (red-line) as unacceptable data if recovery of the ICV or CCV is outside the range 75-125% (CN, 70-130%; Hg, 65-135%). Qualify five samples on either side of verification standard out of control limits. |     |    |     |

|  |                                     |   |   |
|--|-------------------------------------|---|---|
| A.1.11.3 Was continuing calibration performed every 10 samples or every 2 hours? | <input checked="" type="checkbox"/> | — | — |
| Was ICV for cyanides distilled?  | <input checked="" type="checkbox"/> | — | — |

**ACTION:** If no for any of the above, write in the Contract-Problem/Non-Compliance section of the "Data Assessment Narrative".

A.1.12 Form II B (CRDL Standards for AA and ICP) -

|  |                                     |   |   |
|--|-------------------------------------|---|---|
| A.1.12.1 Was a CRDL standard (CRA) analyzed after initial calibration for all AA metals?   | <input checked="" type="checkbox"/> | — | — |
| Was a mid-range calib. verification standard distilled and analyzed for cyanide analysis?  | <input checked="" type="checkbox"/> | — | — |
| Was a 2xCRDL ( or 2xIDL when IDL>CRDL) analyzed (CRI) for each ICP run?<br>(Note: CRI for AL,Ba,Ca,Fe,Mg,Na,or K is not required.) | <input checked="" type="checkbox"/> | — | — |

**ACTION:** If no for any of the above, flag as estimated all data falling within the affected ranges. The affected ranges are:  
 AA Analysis - \*\*True Value  $\pm$  CRDL  
 ICP Analysis - \*\*True Value  $\pm$  2CRDL  
 CN Analysis - \*\*True Value  $\pm$  0.5 x True Value.

\*\*True value of CRA, CRI or mid-range standard. Substitute IDL for CRDL when IDL > CRDL. Compute the concentration of the missing mid-range standard from the calibration range.

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|          |  | YES                                 | NO                       | N/A                      |
|----------|--|-------------------------------------|--------------------------|--------------------------|
| A.1.12.2 | Was CRI analyzed after ICV/ICB and before the final CCV/CCB, and twice every eight hours of ICP run? | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |

**ACTION:** If no, write in Contract Problem/Non-Compliance Section of the "Data Assessment Narrative".

A.1.12.3 Circle on each Form IIB all the percent recoveries that are outside the acceptance windows.

Are CRA and CRI standards within control limits:

|        |             |                          |                                     |                          |
|--------|-------------|--------------------------|-------------------------------------|--------------------------|
| Metals | 80 - 120%R? | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
|--------|-------------|--------------------------|-------------------------------------|--------------------------|

Is mid-range standard within control limits:

|         |             |                                     |                          |                          |
|---------|-------------|-------------------------------------|--------------------------|--------------------------|
| Cyanide | 80 - 120%R? | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|---------|-------------|-------------------------------------|--------------------------|--------------------------|

**ACTION:** Flag as estimated all sample results within the affected range if the recovery of the standard is between 50-79%; flag only positive data within the affected range if the recovery is between 121-150%; reject all data within the affected range if the recovery is less than 50%; reject only positive data within the affected range if the recovery is greater than 150%. Qualify 50% of the samples on either side of CRI standard outside the control limits.

**Note:** Flag or reject the final results only when sample raw data are within the affected ranges and the CRDL standards are outside the acceptance windows.

A.1.13 Form III (Initial and Continuing Calibration Blanks)

|          |                       |                                     |                          |                          |
|----------|-----------------------|-------------------------------------|--------------------------|--------------------------|
| A.1.13.1 | Present and complete? | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|----------|-----------------------|-------------------------------------|--------------------------|--------------------------|

|  |  |                                     |                          |                          |
|--|--|-------------------------------------|--------------------------|--------------------------|
|  | For both AA and ICP when both are used for the same analyte? | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|--|-------------------------------------|--------------------------|--------------------------|

|  |  |                                     |                          |                          |
|--|--|-------------------------------------|--------------------------|--------------------------|
|  | Was an initial calibration blank analyzed? | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|--|-------------------------------------|--------------------------|--------------------------|

|  |  |                                     |                          |                          |
|--|--|-------------------------------------|--------------------------|--------------------------|
|  | Was a continuing calibration blank analyzed after every 10 samples or every 2 hours (which ever is more frequent)? | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
|--|--|-------------------------------------|--------------------------|--------------------------|

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|   | <u>YES</u>                          | <u>NO</u>                | <u>N/A</u>                          |
|---|-------------------------------------|--------------------------|-------------------------------------|
| <b><u>ACTION:</u></b> If no, prepare Telephone Record Log, contact laboratory and write in the Contract-Problems/Non-Compliance section of the "Data Assessment Narrative".   |                                     |                          |                                     |
| <b>A.1.13.2</b> Circle on each Form III all calibration blank values that are above CRDL (or 2 x IDL when IDL > CRDL).  |                                     |                          |                                     |
| Are all calibration blanks (when IDL < CRDL) less than or equal to the Contract Required Detection Limits (CRDLs)?  | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/>            |
| Are all calibration blanks less than two times Instrument Detection Limit (when IDL > CRDL)?  | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/>            |
| <b><u>ACTION:</u></b> If no for any of the above, flag as estimated (J) positive sample results when <u>raw sample value</u> is less than or equal to calibration blank value analyzed between calibration blank with value over CRDL (or 2xIDL) and nearest good calibration blank.<br>Flag five samples on either side of the calibration blank outside the control limits. |                                     |                          |                                     |
| <b>A.1.14</b> <u>FORM III (Preparation Blank) -</u><br>(Note: The preparation blank for mercury is the same as the calibration blank.)  |                                     |                          |                                     |
| <b>A.1.14.1</b> Was one prep. blank analyzed for:   |                                     |                          |                                     |
| each Sample Delivery Group (SDG)?   | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/>            |
| each batch of digested samples?   | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/>            |
| each matrix type?   | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/>            |
| both AA and ICP when both are used for the same analyte?  | <input type="checkbox"/>            | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
| <b><u>ACTION:</u></b> If no for any of the above, flag as estimated (J) all the associated positive data <10 x IDLs for which prep. blank was not analyzed.   |                                     |                          |                                     |
| <b><u>NOTE:</u></b> If only one blank was analyzed for more than 20 samples, then first 20 samples analyzed do not have to be flagged as estimated (J).   |                                     |                          |                                     |



STANDARD OPERATING PROCEDURE

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|          |   | YES                                 | NO                                  | N/A                                 |
|----------|---|-------------------------------------|-------------------------------------|-------------------------------------|
| A.1.14.2 | Is concentration of prep. blank value greater than the CRDL when IDL is less than or equal to CRDL?   | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/>            |
|          | If yes, is the concentration of the sample with the least concentrated analyte less than 10 times the prep. blank?                          | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
|          | <b>ACTION:</b> If yes, reject (red-line) all associated data greater than CRDL concentration but less than ten times the prep. blank value. |                                     |                                     |                                     |
| A.1.14.3 | Is concentration of prep. blank value (Form III) less than two times IDL, when IDL is greater than CRDL?                                    | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
|          | <b>ACTION:</b> If no, reject (red-line) all positive sample results when sample raw data are less than 10 times the prep. blank value.      |                                     |                                     |                                     |
| A.1.14.4 | Is concentration of prep. blank below the negative CRDL?  | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/>            |
|          | <b>ACTION:</b> If yes, reject (red-line) all associated sample results less than 10xCRDL.   |                                     |                                     |                                     |
| A.1.15   | <u>Form IV (ICP Interference Check Sample)</u>  |                                     |                                     |                                     |
| A.1.15.1 | Present and complete?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
|          | (NOTE: Not required for furnace AA, flame AA, mercury, cyanide and Ca, Mg, K and Na.)   |                                     |                                     |                                     |
|          | Was ICS analyzed at beginning and end of run (or at least twice every 8 hours)?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
|          | <b>ACTION:</b> If no, flag as estimated (J) all the samples for which Al, Ca, Fe, or Mg is higher than in ICS.                              |                                     |                                     |                                     |
| A.1.15.2 | Circle all values on each Form IV that are more than $\pm 20\%$ of true or established mean value.  |                                     |                                     |                                     |
|          | Are all Interference Check Sample results inside the control limits ( $\pm 20\%$ )?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
|          | If no, is concentration of Al, Ca, Fe, or Mg lower than the respective concentration in ICS?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |

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|   |  | YES                                 | NO                                  | N/A                                 |
|---|--|-------------------------------------|-------------------------------------|-------------------------------------|
| <p><b>ACTION:</b> If no, flag as estimated (J) those positive results for which ICS recovery is between 121-150%; flag all sample results as estimated if ICS recovery falls within 50-79%; reject (red-line) those sample results for which ICS recovery is less than 50%; if ICS recovery is above 150%, reject positive results only (not flagged with a "U").</p> |  |                                     |                                     |                                     |
| A.1.16  | <p><u>Form V A (Spiked Sample Recovery - Pre-Digestion/Pre-Distillation)-</u><br/>( Note: Not required for Ca, Mg, K, and Na (both matrices), Al, and Fe (soil only.)</p>                        |                                     |                                     |                                     |
| A.1.16.1  | Present and complete for: each SDG?  | <input checked="" type="checkbox"/> | ___                                 | ___                                 |
|   | each matrix type?  | <input checked="" type="checkbox"/> | ___                                 | ___                                 |
|   | each conc. range (i.e. low, med., high)?   | <input checked="" type="checkbox"/> | ___                                 | ___                                 |
|   | For both AA and ICP when both are used for the same analyte?   | <input type="checkbox"/>            | ___                                 | <input checked="" type="checkbox"/> |
|   | <p><b>ACTION:</b> If no for any of the above, flag as estimated (J) all the positive data less than four times the spiking levels specified in SCW for which spiked sample was not analyzed.</p> |                                     |                                     |                                     |
|   | <p><b>NOTE:</b> If one spiked sample was analyzed for more than 20 samples, then first 20 samples analyzed do not have to be flagged as estimated (J).</p>                                       |                                     |                                     |                                     |
| A.1.16.2  | Was field blank used for spiked sample?  | ___                                 | <input checked="" type="checkbox"/> | ___                                 |
|   | <p><b>ACTION:</b> If yes, flag all positive data less than 4 x spike added as estimated (J) for which field blank was used as spiked sample.</p>   |                                     |                                     |                                     |
| A.1.16.3  | Circle on each Form VA all spike recoveries that are outside control limits (75% to 125%).   |                                     |                                     |                                     |
|   | Are all recoveries within control limits?  | <input checked="" type="checkbox"/> | ___                                 | ___                                 |
|   | If no, is sample concentration greater than or equal to four times spike concentration?  | <input type="checkbox"/>            | ___                                 | <input checked="" type="checkbox"/> |

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YES      NO      N/A

**ACTION:** If yes, disregard spike recoveries for analytes whose concentrations are greater than or equal to four times spike added. If no, circle those analytes on Form V for which sample concentration is less than four times the spike concentration.

Are results outside the control limits (75-125%) flagged with "N" on Form I's and Form VA?

           1

**ACTION:** If no, write in the Contract - Problem/Non - Compliance section of "Data Assessment Narrative".

A.1.16.4 Aqueous

Are any spike recoveries:

(a) less than 30%?

           4

(b) between 30-74%?

           7

(c) between 126-150%?

           7

(d) greater than 150%?

           7

**ACTION:** If less than 30%, reject all associated aqueous data; if between 30-74%, flag all associated aqueous data as estimated (J); if between 126-150%, flag as estimated (J) all associated aqueous data not flagged with a "U"; if greater than 150%, reject (red-line) all associated aqueous data not flagged with a "U".

A.1.16.5 Soil/Sediment

Are any spike recoveries:

(a) less than 10%?

           1

(b) between 10-74%?

           1

(c) between 126-200%?

           1

(d) greater than 200%?

           1

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|  | <u>YES</u> | <u>NO</u> | <u>N/A</u> |
|--|------------|-----------|------------|
| <b>ACTION:</b> If less than 10%, reject all associated data; if between 10-74%, flag all associated data as estimated; if between 126-200%, flag as estimated all associated data was not flagged with a "U"; if greater than 200%, reject all associated data not flagged with a "U". |            |           |            |

**A.1.17 Form VI (Lab Duplicates)**

|          |                           |  |                                     |   |                                     |
|----------|---------------------------|--|-------------------------------------|---|-------------------------------------|
| A.1.17.1 | Present and complete for: | each SDG?  | <input checked="" type="checkbox"/> | — | —                                   |
|          |                           | each matrix type?  | <input checked="" type="checkbox"/> | — | —                                   |
|          |                           | each concentration range (i.e. low, med., high)?         | <input type="checkbox"/>            | — | —                                   |
|          |                           | both AA and ICP when both are used for the same analyte? | <input type="checkbox"/>            | — | <input checked="" type="checkbox"/> |

**ACTION:** If no for any the above, flag as estimated (J) all the data  $\geq$ CRDL\* for which duplicate sample was not analyzed.

- Note:** 1. If one duplicate sample was analyzed for more than 20 samples, then first 20 samples do not have to be flagged as estimated.  
 2. If percent solids for soil sample and its duplicate differ by more than 1%, prepare a Form VI for each duplicate pair, report concentrations in ug/L on wet weight basis and calculate RPD or Difference for each analyte.

|          |  |   |                                     |   |
|----------|--|---|-------------------------------------|---|
| A.1.17.2 | Was field blank used for duplicate analysis? | — | <input checked="" type="checkbox"/> | — |
|----------|--|---|-------------------------------------|---|

**ACTION:** If yes, flag all data  $\geq$ CRDL\* as estimated (J) for which field blank was used as duplicate.

|          |   |                                     |   |                                     |
|----------|---|-------------------------------------|---|-------------------------------------|
| A.1.17.3 | Are all values within control limits (RPD 20% or difference $\leq$ $\pm$ CRDL)?         | <input checked="" type="checkbox"/> | — | —                                   |
|          | If no, are all results outside the control limits flagged with an * on Form I's and VI? | <input type="checkbox"/>            | — | <input checked="" type="checkbox"/> |

**ACTION:** If no, write in the Contract - Problems/Non-Compliance section of "Data Assessment Narrative".

\* Substitute IDL for CRDL when IDL > CRDL.

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- NOTE:** 1. RPD is not calculable for an analyte of the sample - duplicate pair when both values are less than IDL.  
 2. If the result of lab duplicate analyzed by GFAA is rejectable due to coefficient of correlation of MSA, analytical spike recovery, or duplicate injections criteria, do not apply precision criteria to metals analyzed by GFAA.

YES      NO      N/A

A.1.17.4 Aqueous

Circle on each Form VI all values that are:

RPD > 50%, or  
 Difference > CRDL\*

Is any RPD greater than 50% where sample and duplicate are both greater than or equal to 5 times \*CRDL?

—           

Is any difference\*\* between sample and duplicate greater than \*CRDL where sample and/or duplicate is less than 5 times \*CRDL?

—           

**ACTION:** If yes, flag the associated data as estimated.

A.1.17.5 Soil/Sediment

Circle on each Form VI all values that are:

RPD > 100%, or  
 Difference > 2 x CRDL\*

Is any RPD (where sample and duplicate are both greater than or equal to 5 times \*CRDL) :

> 100%?      —            —

Is any \*\*difference between sample and duplicate (where sample and/or duplicate is less than 5x\*CRDL) :

> 2x\*CRDL?      —            —

\* Substitute IDL for CRDL when IDL > CRDL.

\*\* Use absolute values of sample and duplicate to calculate the difference.

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YES NO N/A

ACTION: If yes, flag the associated data as estimated.

A.1.18 Field Duplicates

A.1.18.1 Were field duplicates analyzed?

[ ] [ ] [ ]

ACTION: If yes, prepare a Form VI for each aqueous field duplicate pair. Prepare a Form VI for each soil duplicate pair, if percent solids for sample and its duplicate differ by more than 1%; report concentrations of soils in ug/l on wet weight basis and calculate RPDs or Difference for each analyte.

- NOTE: 1. Do not calculate RPD when both values are less than IDL. 2. Flag all associated data only for field duplicate pair.

A.1.18.2 Aqueous

Circle all values on self prepared Form VI for field duplicates that are:

RPD > 50%, or Difference > CRDL\*

Is any RPD greater than 50% where sample and duplicate are both greater than or equal to 5 times \*CRDL?

[ ] [ ] [ ]

Is any \*\*difference between sample and duplicate greater than \*CRDL where sample and/or duplicate is less than 5 times \*CRDL?

[ ] [ ] [ ]

ACTION: If yes, flag the associated data as estimated.

\* Substitute IDL for CRDL when IDL > CRDL. \*\* Use absolute values of sample and duplicate to calculate the difference.

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|                                      | YES | NO | N/A |
|--------------------------------------|-----|----|-----|
| <b>A.1.18.3 <u>Soil/Sediment</u></b> |     |    |     |

Circle all values on self prepared Form VI for field duplicates that are:

RPD >100%, or

Difference > 2 x CRDL\*

Is any RPD (where sample and duplicate are both greater than 5 times \*CRDL) :

>100%?

Is any \*\*difference between sample and duplicate (where sample and/or duplicate is less than 5x \*CRDL) :

>2x \*CRDL?

**ACTION:** If yes, flag the associated data as estimated.

**A.1.19 Form VII (Laboratory Control Sample)** (Note: LCS - not required for aqueous Hg and cyanide analyses.)

**A.1.19.1** Was one LCS prepared and analyzed for:

each SDG?

each batch samples digested/distilled?

both AA and ICP when both are used for the same analyte?

**ACTION:** If no for any of the above, prepare Telephone Record Log and contact laboratory for submittal of results of LCS. Flag as estimated (J) all the data for which LCS was not analyzed.

**NOTE:** If only one LCS was analyzed for more than 20 samples, then first 20 samples close to LCS do not have to be flagged as estimated.

\* Substitute IDL for CRDL when IDL > CRDL.  
 \*\* Use absolute values of sample and duplicate to calculate the difference.

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|          |   | <u>YES</u> | <u>NO</u>                           | <u>N/A</u> |
|----------|---|------------|-------------------------------------|------------|
| A.1.19.2 | <u>Aqueous LCS</u>  |            |                                     |            |
|          | Circle on each Form VII the LCS percent recoveries outside control limits (80 - 120%) except for aqueous Ag and Sb.   |            |                                     |            |
|          | Is any LCS recovery:  |            |                                     |            |
|          | less than 50%?  | —          | <input checked="" type="checkbox"/> | —          |
|          | between 50% and 79%?  | —          | <input checked="" type="checkbox"/> | —          |
|          | between 121% and 150%?  | —          | <input checked="" type="checkbox"/> | —          |
|          | greater than 150%?  | —          | <input checked="" type="checkbox"/> | —          |
|          | <u>ACTION:</u> Less than 50%, reject (red-line) all data; between 50% and 79%, flag all associated data as estimated (J); between 121% and 150%, flag all positive (not flagged with a "U") results as estimated; greater than 150%, reject all positive results.   |            |                                     |            |
| A.1.19.3 | <u>Solid LCS</u>  |            |                                     |            |
|          | <u>NOTE:</u> 1. If "Found" value of LCS is rejectable due to duplicate injections or <u>analytical spike</u> recovery criteria, regardless of LCS recovery, flag the associated data as estimated (J).<br>2. If IDL of an analyte is equal to or greater than true value of LCS, disregard the "Action" below even though LCS is out of control limits. |            |                                     |            |
|          | Is LCS "Found" value higher than the control limits on Form VII?  | —          | <input checked="" type="checkbox"/> | —          |
|          | <u>ACTION:</u> If yes, qualify all associated positive data as estimated.   |            |                                     |            |
|          | Is LCS "Found" value lower than the Control limits on Form VII?   | —          | <input checked="" type="checkbox"/> | —          |
|          | <u>ACTION:</u> If yes, qualify all associated data as estimated.  |            |                                     |            |



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|          |  | YES                                 | NO                                  | N/A |
|----------|--|-------------------------------------|-------------------------------------|-----|
| A.1.20   | <u>Form IX (ICP Serial Dilution) -</u><br><br><u>NOTE:</u> Serial dilution analysis is required only for initial concentrations equal to or greater than 10 x IDL.   |                                     |                                     |     |
| A.1.20.1 | Was Serial Dilution analysis performed for:<br>each SDG?<br>each matrix type?<br>each concentration range (i.e. low, med.)?  | <input checked="" type="checkbox"/> | —                                   | —   |
|          | <u>ACTION:</u> If no for any of the above, flag as estimated all the positive data $\geq 10 \times \text{IDLs}$ or $\geq \text{CRDL}$ when $10 \times \text{IDL} \leq \text{CRDL}$ for which Serial Dilution Analysis was not performed. |                                     |                                     |     |
| A.1.20.2 | Was field blank(s) used for Serial Dilution Analysis?  | —                                   | <input checked="" type="checkbox"/> | —   |
|          | <u>ACTION:</u> If yes, flag all associated data $\geq 10 \times \text{IDL}$ as estimated (J). If $10 \times \text{IDL} \leq \text{CRDL}$ , flag all data $\geq \text{CRDL}$ .  |                                     |                                     |     |
| A.1.20.3 | Are results outside control limit flagged with an "E" on Form I's and Form IX when initial concentration on Form IX is equal to 50 times IDL or greater.   | <input checked="" type="checkbox"/> | —                                   | —   |
|          | <u>ACTION:</u> If no, write in the Contract-Problem/Non-Compliance section of the "Data Assessment Narrative".   |                                     |                                     |     |
| A.1.20.4 | Circle on each Form IX all percent difference that are outside the control limits for initial concentrations equal to or greater than 10 x IDLs only.<br><br>Are any % difference values:  |                                     |                                     |     |
|          | > 10%?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | —   |
|          | $\geq 100\%$ ?   | —                                   | <input checked="" type="checkbox"/> | —   |

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|  | YES  | NO                       | N/A                                 |
|--|--|--------------------------|-------------------------------------|
| <p><b>ACTION:</b> Flag as estimated (J) all the associated sample data <math>\geq 10 \times \text{IDLs}</math> (or <math>\geq \text{CRDL}</math> when <math>10 \times \text{IDL} \leq \text{CRDL}</math>) for which percent difference is greater than 10% but less than 100%. Reject (red-line) all the associated sample results equal to or greater than <math>10 \times \text{IDLs}</math> (or <math>\geq \text{CRDL}</math> when <math>10 \times \text{IDL} \leq \text{CRDL}</math>) for which PD is greater than or equal to 100%.</p> |  |                          |                                     |
| <p><b>Note:</b> Flag or reject on Form I's only the sample results whose associated raw data are <math>\geq 10 \times \text{IDL}</math> (or <math>\geq \text{CRDL}</math> when <math>10 \times \text{IDL} \leq \text{CRDL}</math>)</p>   |  |                          |                                     |
| <p><b>A.1.21 <u>Furnace Atomic Absorbtion (AA) QC Analysis</u></b></p>   |  |                          |                                     |
| A.1.21.1   | Are duplicate injections present in furnace raw data (except during full Method of Standard Addition) for each sample analyzed by GFAA?                      | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
| <p><b>ACTION:</b> If no, <u>reject</u> the data on Form I's for which duplicate injections were not performed.</p>   |  |                          |                                     |
| A.1.21.2   | Do the duplicate injection readings agree within 20% Relative Standard Deviation (RSD) or Coefficient of Variation (CV) for concentration greater than CRDL? | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
|  | Was a dilution analyzed for sample with analytical spike recovery less than 40%?   | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
| <p><b>ACTION:</b> If no for any of the above, flag all the associated data as estimated.</p>   |  |                          |                                     |
| A.1.21.3   | Is *analytical spike recovery outside the control limits (85-115%) for any sample?   | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
| <p><b>ACTION:</b> If yes, flag as estimated the affected sample results if the recovery is between 10-84%; if the recovery is between 115-200%, flag the associated positive sample results as estimated; reject the associated sample results if the recovery is less than 10%; reject positive sample results if the recovery is greater than 200%.</p>  |  |                          |                                     |

\* Analytical spike is not required on the pre-digestion spiked sample.

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|   | YES                      | NO                       | N/A                                 |
|---|--------------------------|--------------------------|-------------------------------------|
| <p><b>NOTE:</b> Reject or flag the data only when the affected sample(s) was not subsequently analyzed by Method of Standard Addition.</p>                  |                          |                          |                                     |
| <p><b>A.1.22 Form VIII (Method of Standard Addition Results)</b></p>  |                          |                          |                                     |
| A.1.22.1  |                          |                          |                                     |
| Present?  | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
| If no, is any Form I result coded with "S" or a "+"?  | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
| <p><b>ACTION:</b> If yes, write request on Telephone Record Log and contact laboratory for submittal of Form VIII.</p>                                      |                          |                          |                                     |
| A.1.22.2  |                          |                          |                                     |
| Is coefficient of correlation for MSA less than 0.990 for any sample?   | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
| <p><b>ACTION:</b> If yes, reject (red-line) the affected data.</p>  |                          |                          |                                     |
| A.1.22.3  |                          |                          |                                     |
| Was *MSA required for any sample but not performed?   | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
| Is coefficient of correlation for MSA less than 0.995?  | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
| Are MSA calculations outside the linear range of the calibration curve generated at the beginning of the analytical run?                                    | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
| <p><b>ACTION:</b> If yes for any of the above, flag all the associated data as estimated (J).</p>   |                          |                          |                                     |
| A.1.22.4  |                          |                          |                                     |
| Was proper quantitation procedure followed correctly as outlined in the SOW on page E-23?   | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
| <p><b>ACTION:</b> If no, note exception under Contract Problem/ Non-Compliance section of the "Data Assessment Narrative", and prepare a separate list.</p> |                          |                          |                                     |

\* MSA is not required on LCS and prep. blank.

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|          |  | YES                                 | NO                                  | N/A                                 |
|----------|--|-------------------------------------|-------------------------------------|-------------------------------------|
| A.1.23   | <u>Dissolved/Total or Inorganic/Total Analytes -</u>   |                                     |                                     |                                     |
| A.1.23.1 | Were any analyses performed for dissolved as well as total analytes on the same sample(s).   | —                                   | <input checked="" type="checkbox"/> | —                                   |
|          | Were any analyses performed for inorganic as well as total (organic + inorganic) analytes on the same sample(s)?   | —                                   | <input checked="" type="checkbox"/> | —                                   |
|          | <p><b>NOTE:</b> 1. If yes, prepare a list comparing differences between all dissolved (or inorganic) and total analytes. Compute the differences as a percent of the total analyte only when dissolved concentration is greater than CRDL as well as total concentration.</p> <p>2. Apply the following questions only if inorganic (or dissolved) results are (i) above CRDL, and (ii) greater than total constituents.</p> <p>3. At least one preparation blank, ICS, and LCS should be analyzed in each analytical run.</p> |                                     |                                     |                                     |
| A.1.23.2 | Is the concentration of any dissolved (or inorganic) analyte greater than its total concentration by more than 10%?  | —                                   | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| A.1.23.3 | Is the concentration of any dissolved (or inorganic) analyte greater than its total concentration by more than 50%?  | —                                   | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
|          | <p><b>ACTION:</b> If more than 10%, flag both dissolved (or inorganic) and total values as estimated (J); if more than 50%, reject (red-line) the data for both values.</p>  |                                     |                                     |                                     |
| A.1.24   | <u>Form I (Field Blank) -</u>  |                                     |                                     |                                     |
|          | <u>(Note: Designate "Field Blank" as such on Form I.)</u>  |                                     |                                     |                                     |
| A.1.24.1 | Circle all field blank values on Form I that are greater than CRDL, (or 2 x IDL when IDL > CRDL).  |                                     |                                     |                                     |
|          | Is field blank concentration less than CRDL (or 2 x IDL when IDL > CRDL) for all parameters of associated aqueous and soil samples?  | <input checked="" type="checkbox"/> | —                                   | —                                   |

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|   | YES                                 | NO                                  | N/A                                 |
|---|-------------------------------------|-------------------------------------|-------------------------------------|
| If no, was field blank value already rejected due to other QC criteria?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| <b>ACTION:</b> If no, reject (except field blank results) all associated positive sample data less than or equal to five times the field blank value. Reject on Form I's the soil sample results that when converted to ug/L on wet basis are less than or equal to five times the field blank value in ug/L. |                                     |                                     |                                     |
| <b>A.1.25 <u>Form I, XI, XII (Verification of Instrumental Parameters).</u></b>   |                                     |                                     |                                     |
| <b>A.1.25.1</b> Is verification report present for:   |                                     |                                     |                                     |
| Instrument Detection Limits (quarterly)?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| ICP Interelement Correction Factors (annually)?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| ICP Linear Ranges (quarterly)?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| <b>ACTION:</b> If no, contact TPO of the lab.   |                                     |                                     |                                     |
| <b>A.1.25.2 <u>Form X (Instrument Detection Limits)</u> - (Note: IDL is not required for Cyanide.)</b>  |                                     |                                     |                                     |
| <b>A.1.25.2.1</b> Are IDLs present for:   |                                     |                                     |                                     |
| all the analytes?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| all the instruments used?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| For both AA and ICP when both are used for the same analyte?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| <b>ACTION:</b> If no for any of the above, prepare Telephone Record Log and contact laboratory.   |                                     |                                     |                                     |
| <b>A.1.25.2.2</b> Is IDL greater than CRDL for any analyte?   | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/>            |
| If yes, is the concentration on Form I of the sample analyzed on the instrument whose IDL exceeds CRDL, greater than 5 x IDL.   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |

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YES NO N/A

Action : If no, flag as estimated all values less than five times IDL of the instrument whose IDL exceeds CRDL.

A.1.25.3 Form XI (Linear Ranges)

A.1.25.3.1 Was any sample result higher than high linear range of ICP.

—  —

Was any sample result higher than the highest calibration standard for non-ICP parameters?

—  —

If yes for any of the above, was the sample diluted to obtain the result on Form I?

—

ACTION: If no, flag the result reported on Form I as estimated(J).

A.1.26 Percent Solids of Sediments

A.1.26.1 Are percent solids in sediment(s):

< 50%?

—  —

< 10%?

—  —

ACTION: If yes, qualify as estimated all the results of a sample that has per cent solids between 10%-50% (i.e. moisture content between 50%-90%). Reject all the results of a sample that has per cent solids less than 10% (i.e. moisture content greater than 90%).

NOTE: Reject or flag(J) only the sample results that were not previously rejected or flagged due to other QC criteria.



**SDG NARRATIVE**

USEPA

SDG # MBNJ51

Case # 28706

Contract # 68-W00088

LAB CODE: CHEMED

Chemtech Project # R1115

**A. Number of Samples and Date of Receipt**

1 Aqueous plus 15 Soil Samples were delivered to the laboratory intact on 11/08/00.

**B. Parameters**

Test requested was Metals. This data package contains results for Metals.

**C. Cooler Temp**Indicator Bottle: Presence/Absence

Cooler Temp: 4.0

**D. Detail Documentation (related to Sample Handling  
Shipping, Analytical Problem, Temp of Cooler etc):****E. Corrective Action taken for above:**

No Corrective Action was taken at this time

**F. Analytical Techniques:**

The analysis of Metals is based on CLP Methodology. Mercury by Method ILM4.0 and Cyanide by Method 335.

**G. QA/ QC**

Calibrations met requirements. Blank analyses did not indicate the presence of contamination. Interference Check Sample, Laboratory Control Sample were within Control Limits. Spike Sample recovery met requirements. Serial Dilution met requirements except for the followings: Arsenic, Beryllium, Copper, Potassium, Selenium and Sodium. Duplicate analyses met requirements except for the Selenium.

I certify that the data package is in compliance with the terms and conditions of the contract both technically and for completeness, for other than the conditions detailed above. Release of the data contained in this hard copy data package has been authorized by the Laboratory Director or his designee, as verified by the following signature.

Signature Mildred V. Reyes Name: Mildred V. ReyesDate 11/17/00 Title: QA/QC



INORGANIC ANALYSIS DATA SHEET

MBNJ51

Lab Name: CHEMTECH EDISON Contract: 68-W00-088  
 Lab Code: CHEMED Case No.: 28706 SAS No.: SDG No.: MBNJ51  
 Matrix (soil/water): SOIL Lab Sample ID: R1115-01S  
 Level (low/med): LOW Date Received: 11/08/00  
 % Solids: 66.4

Concentration Units (ug/L or mg/Kg dry weight): MG/KG

| CAS No.   | Analyte   | Concentration | C | Q | M  |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum  | 6770          |   |   | P  |
| 7440-36-0 | Antimony  | 1.8           | U |   | P  |
| 7440-38-2 | Arsenic   | 2.9           | B |   | P  |
| 7440-39-3 | Barium    | 67.3          |   |   | P  |
| 7440-41-7 | Beryllium | 0.45          | B |   | P  |
| 7440-43-9 | Cadmium   | 0.60          | U |   | P  |
| 7440-70-2 | Calcium   | 2140          |   |   | P  |
| 7440-47-3 | Chromium  | 11.5          |   |   | P  |
| 7440-48-4 | Cobalt    | 9.1           | B |   | P  |
| 7440-50-8 | Copper    | 19.7          |   | E | P  |
| 7439-89-6 | Iron      | 12100         |   |   | P  |
| 7439-92-1 | Lead      | 12.6          |   |   | P  |
| 7439-95-4 | Magnesium | 2630          |   |   | P  |
| 7439-96-5 | Manganese | 117           |   |   | P  |
| 7439-97-6 | Mercury   | 0.14          | U |   | CV |
| 7440-02-0 | Nickel    | 27.0          |   |   | P  |
| 7440-09-7 | Potassium | 639           | B | E | P  |
| 7782-49-2 | Selenium  | 1.3           | B |   | P  |
| 7440-22-4 | Silver    | 0.48          | U |   | P  |
| 7440-23-5 | Sodium    | 161           | B |   | P  |
| 7440-28-0 | Thallium  | 1.1           | U |   | P  |
| 7440-62-2 | Vanadium  | 12.8          | B |   | P  |
| 7440-66-6 | Zinc      | 71.4          |   |   | P  |
|           | Cyanide   | 0.22          | U |   | CA |

Color Before: BROWN Clarity Before: Texture: MEDIUM  
 Color After: YELLOW Clarity After: Artifacts:

Comments:

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INORGANIC ANALYSIS DATA SHEET

MBNJ52

Lab Name: CHEMTECH EDISON Contract: 68-W00-088  
 Lab Code: CHEMED Case No.: 28706 SAS No.: SDG No.: MBNJ51  
 Matrix (soil/water): SOIL Lab Sample ID: R1115-02S  
 Level (low/med): LOW Date Received: 11/08/00  
 % Solids: 66.2

Concentration Units (ug/L or mg/Kg dry weight): MG/KG

| CAS No.   | Analyte   | Concentration | C | Q | M  |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum  | 8590          | - |   | P  |
| 7440-36-0 | Antimony  | 1.8           | U |   | P  |
| 7440-38-2 | Arsenic   | 1.7           | U |   | P  |
| 7440-39-3 | Barium    | 97.3          |   |   | P  |
| 7440-41-7 | Beryllium | 0.72          | B |   | P  |
| 7440-43-9 | Cadmium   | 0.60          | U |   | P  |
| 7440-70-2 | Calcium   | 3220          |   |   | P  |
| 7440-47-3 | Chromium  | 14.7          |   |   | P  |
| 7440-48-4 | Cobalt    | 9.7           | B |   | P  |
| 7440-50-8 | Copper    | 20.0          |   | E | P  |
| 7439-89-6 | Iron      | 16200         |   |   | P  |
| 7439-92-1 | Lead      | 15.3          |   |   | P  |
| 7439-95-4 | Magnesium | 3390          |   |   | P  |
| 7439-96-5 | Manganese | 138           |   |   | P  |
| 7439-97-6 | Mercury   | 0.15          | U |   | CV |
| 7440-02-0 | Nickel    | 27.3          |   |   | P  |
| 7440-09-7 | Potassium | 938           | B | E | P  |
| 7782-49-2 | Selenium  | 1.2           | U |   | P  |
| 7440-22-4 | Silver    | 0.48          | U |   | P  |
| 7440-23-5 | Sodium    | 77.5          | U |   | P  |
| 7440-28-0 | Thallium  | 1.1           | U |   | P  |
| 7440-62-2 | Vanadium  | 14.4          | B |   | P  |
| 7440-66-6 | Zinc      | 72.1          |   |   | P  |
|           | Cyanide   | 0.22          | U |   | CA |

Color Before: BROWN Clarity Before: Texture: MEDIUM  
 Color After: YELLOW Clarity After: Artifacts:

Comments:

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0.39

INORGANIC ANALYSIS DATA SHEET

MBNJ53

Lab Name: CHEMTECH EDISON

Contract: 68-W00-088

Lab Code: CHEMED

Case No.: 28706

SAS No.:

SDG No.: MBNJ51

Matrix (soil/water): SOIL

Lab Sample ID: R1115-03S

Level (low/med): LOW

Date Received: 11/08/00

% Solids: 79.1

Concentration Units (ug/L or mg/Kg dry weight): MG/KG

| CAS No.   | Analyte   | Concentration | C | Q | M  |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum  | 15900         | - |   | P  |
| 7440-36-0 | Antimony  | 1.5           | U |   | P  |
| 7440-38-2 | Arsenic   | 4.1           |   |   | P  |
| 7440-39-3 | Barium    | 187           |   |   | P  |
| 7440-41-7 | Beryllium | 0.73          | B |   | P  |
| 7440-43-9 | Cadmium   | 0.50          | U |   | P  |
| 7440-70-2 | Calcium   | 2640          |   |   | P  |
| 7440-47-3 | Chromium  | 26.2          |   |   | P  |
| 7440-48-4 | Cobalt    | 20.3          |   |   | P  |
| 7440-50-8 | Copper    | 21.0          |   | E | P  |
| 7439-89-6 | Iron      | 30100         |   |   | P  |
| 7439-92-1 | Lead      | 21.6          |   |   | P  |
| 7439-95-4 | Magnesium | 5510          |   |   | P  |
| 7439-96-5 | Manganese | 1050          |   |   | P  |
| 7439-97-6 | Mercury   | 0.13          | U |   | CV |
| 7440-02-0 | Nickel    | 48.7          |   |   | P  |
| 7440-09-7 | Potassium | 1080          | B | E | P  |
| 7782-49-2 | Selenium  | 0.99          | U |   | P  |
| 7440-22-4 | Silver    | 0.88          | B |   | P  |
| 7440-23-5 | Sodium    | 64.2          | U |   | P  |
| 7440-28-0 | Thallium  | 0.92          | U |   | P  |
| 7440-62-2 | Vanadium  | 24.0          |   |   | P  |
| 7440-66-6 | Zinc      | 128           |   |   | P  |
|           | Cyanide   | 0.18          | U |   | CA |

Color Before: BROWN

Clarity Before:

Texture: MEDIUM

Color After: YELLOW

Clarity After:

Artifacts:

Comments:

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0.40

INORGANIC ANALYSIS DATA SHEET

MBNJ54

Lab Name: CHEMTECH EDISON Contract: 68-W00-088  
 Lab Code: CHEMED Case No.: 28706 SAS No.: SDG No.: MBNJ51  
 Matrix (soil/water): SOIL Lab Sample ID: R1115-05S  
 Level (low/med): LOW Date Received: 11/08/00  
 % Solids: 96.7

Concentration Units (ug/L or mg/Kg dry weight): MG/KG

| CAS No.   | Analyte   | Concentration | C | Q | M  |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum  | 686           | - |   | P  |
| 7440-36-0 | Antimony  | 1.2           | U |   | P  |
| 7440-38-2 | Arsenic   | 1.1           | U |   | P  |
| 7440-39-3 | Barium    | 9.4           | B |   | P  |
| 7440-41-7 | Beryllium | 0.22          | U |   | P  |
| 7440-43-9 | Cadmium   | 0.40          | U |   | P  |
| 7440-70-2 | Calcium   | 936           | B |   | P  |
| 7440-47-3 | Chromium  | 6.2           |   |   | P  |
| 7440-48-4 | Cobalt    | 0.99          | B |   | P  |
| 7440-50-8 | Copper    | 12.5          |   | E | P  |
| 7439-89-6 | Iron      | 7390          |   |   | P  |
| 7439-92-1 | Lead      | 15.2          |   |   | P  |
| 7439-95-4 | Magnesium | 244           | B |   | P  |
| 7439-96-5 | Manganese | 56.1          |   |   | P  |
| 7439-97-6 | Mercury   | 0.09          |   |   | CV |
| 7440-02-0 | Nickel    | 3.6           | B |   | P  |
| 7440-09-7 | Potassium | 90.1          | B | E | P  |
| 7782-49-2 | Selenium  | 0.80          | U |   | P  |
| 7440-22-4 | Silver    | 0.32          | U |   | P  |
| 7440-23-5 | Sodium    | 52.0          | U |   | P  |
| 7440-28-0 | Thallium  | 0.74          | U |   | P  |
| 7440-62-2 | Vanadium  | 3.0           | B |   | P  |
| 7440-66-6 | Zinc      | 23.8          |   |   | P  |
|           | Cyanide   | 0.15          | U |   | CA |

Color Before: BROWN Clarity Before: Texture: MEDIUM  
 Color After: YELLOW Clarity After: Artifacts:

Comments:

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*p. 4/1*

INORGANIC ANALYSIS DATA SHEET

MBNJ55

Lab Name: CHEMTECH EDISON

Contract: 68-W00-088

Lab Code: CHEMED

Case No.: 28706

SAS No.:

SDG No.: MBNJ51

Matrix (soil/water): SOIL

Lab Sample ID: R1115-06S

Level (low/med): LOW

Date Received: 11/08/00

% Solids: 90.0

Concentration Units (ug/L or mg/Kg dry weight): MG/KG

| CAS No.   | Analyte   | Concentration | C | Q | M  |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum  | 4170          |   |   | P  |
| 7440-36-0 | Antimony  | 1.3           | U |   | P  |
| 7440-38-2 | Arsenic   | 1.8           | B |   | P  |
| 7440-39-3 | Barium    | 35.1          | B |   | P  |
| 7440-41-7 | Beryllium | 0.24          | U |   | P  |
| 7440-43-9 | Cadmium   | 0.44          | U |   | P  |
| 7440-70-2 | Calcium   | 60300         |   |   | P  |
| 7440-47-3 | Chromium  | 8.0           |   |   | P  |
| 7440-48-4 | Cobalt    | 2.4           | B |   | P  |
| 7440-50-8 | Copper    | 14.3          |   | E | P  |
| 7439-89-6 | Iron      | 9500          |   |   | P  |
| 7439-92-1 | Lead      | 24.6          |   |   | P  |
| 7439-95-4 | Magnesium | 2490          |   |   | P  |
| 7439-96-5 | Manganese | 145           |   |   | P  |
| 7439-97-6 | Mercury   | 0.14          |   |   | CV |
| 7440-02-0 | Nickel    | 7.7           | B |   | P  |
| 7440-09-7 | Potassium | 425           | B | E | P  |
| 7782-49-2 | Selenium  | 0.89          | U |   | P  |
| 7440-22-4 | Silver    | 0.36          | U |   | P  |
| 7440-23-5 | Sodium    | 88.4          | B |   | P  |
| 7440-28-0 | Thallium  | 0.82          | U |   | P  |
| 7440-62-2 | Vanadium  | 6.0           | B |   | P  |
| 7440-66-6 | Zinc      | 46.4          |   |   | P  |
|           | Cyanide   | 0.16          | U |   | CA |

Color Before: BROWN

Clarity Before:

Texture: MEDIUM

Color After: YELLOW

Clarity After:

Artifacts:

Comments:

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P. 42

INORGANIC ANALYSIS DATA SHEET

MBNJ56

Lab Name: CHEMTECH EDISON

Contract: 68-W00-088

Lab Code: CHEMED

Case No.: 28706

SAS No.:

SDG No.: MBNJ51

Matrix (soil/water): SOIL

Lab Sample ID: R1115-07S

Level (low/med): LOW

Date Received: 11/08/00

% Solids: 83.2

Concentration Units (ug/L or mg/Kg dry weight): MG/KG

| CAS No.   | Analyte   | Concentration | C | Q | M  |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum  | 12300         |   |   | P  |
| 7440-36-0 | Antimony  | 1.4           | U |   | P  |
| 7440-38-2 | Arsenic   | 4.2           |   |   | P  |
| 7440-39-3 | Barium    | 110           |   |   | P  |
| 7440-41-7 | Beryllium | 0.66          | B |   | P  |
| 7440-43-9 | Cadmium   | 0.46          | U |   | P  |
| 7440-70-2 | Calcium   | 1900          |   |   | P  |
| 7440-47-3 | Chromium  | 18.9          |   |   | P  |
| 7440-48-4 | Cobalt    | 10.1          | B |   | P  |
| 7440-50-8 | Copper    | 21.3          |   | E | P  |
| 7439-89-6 | Iron      | 24800         |   |   | P  |
| 7439-92-1 | Lead      | 18.4          |   |   | P  |
| 7439-95-4 | Magnesium | 4430          |   |   | P  |
| 7439-96-5 | Manganese | 217           |   |   | P  |
| 7439-97-6 | Mercury   | 0.12          | U |   | CV |
| 7440-02-0 | Nickel    | 34.5          |   |   | P  |
| 7440-09-7 | Potassium | 847           | B | E | P  |
| 7782-49-2 | Selenium  | 0.92          | U |   | P  |
| 7440-22-4 | Silver    | 0.53          | B |   | P  |
| 7440-23-5 | Sodium    | 228           | B |   | P  |
| 7440-28-0 | Thallium  | 0.86          | U |   | P  |
| 7440-62-2 | Vanadium  | 19.7          |   |   | P  |
| 7440-66-6 | Zinc      | 91.0          |   |   | P  |
|           | Cyanide   | 0.17          | U |   | CA |

Color Before: BROWN

Clarity Before:

Texture: MEDIUM

Color After: YELLOW

Clarity After:

Artifacts:

Comments:

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0.43

INORGANIC ANALYSIS DATA SHEET

MBNJ57

Lab Name: CHEMTECH EDISON

Contract: 68-W00-088

Lab Code: CHEMED

Case No.: 28706

SAS No.:

SDG No.: MBNJ51

Matrix (soil/water): SOIL

Lab Sample ID: R1115-08S

Level (low/med): LOW

Date Received: 11/08/00

% Solids: 80.4

Concentration Units (ug/L or mg/Kg dry weight): MG/KG

| CAS No.   | Analyte   | Concentration | C | Q | M  |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum  | 15100         | - |   | P  |
| 7440-36-0 | Antimony  | 1.5           | U |   | P  |
| 7440-38-2 | Arsenic   | 16.9          |   |   | P  |
| 7440-39-3 | Barium    | 166           |   |   | P  |
| 7440-41-7 | Beryllium | 1.0           | B |   | P  |
| 7440-43-9 | Cadmium   | 0.49          | U |   | P  |
| 7440-70-2 | Calcium   | 2630          |   |   | P  |
| 7440-47-3 | Chromium  | 23.7          |   |   | P  |
| 7440-48-4 | Cobalt    | 12.7          |   |   | P  |
| 7440-50-8 | Copper    | 41.9          |   | E | P  |
| 7439-89-6 | Iron      | 52800         |   |   | P  |
| 7439-92-1 | Lead      | 25.2          |   |   | P  |
| 7439-95-4 | Magnesium | 5450          |   |   | P  |
| 7439-96-5 | Manganese | 197           |   |   | P  |
| 7439-97-6 | Mercury   | 0.11          | U |   | CV |
| 7440-02-0 | Nickel    | 48.9          |   |   | P  |
| 7440-09-7 | Potassium | 1030          | B | E | P  |
| 7782-49-2 | Selenium  | 0.98          | U |   | P  |
| 7440-22-4 | Silver    | 1.3           | B |   | P  |
| 7440-23-5 | Sodium    | 63.2          | U |   | P  |
| 7440-28-0 | Thallium  | 0.90          | U |   | P  |
| 7440-62-2 | Vanadium  | 24.0          |   |   | P  |
| 7440-66-6 | Zinc      | 127           |   |   | P  |
|           | Cyanide   | 0.18          | U |   | CA |

Color Before: BROWN

Clarity Before:

Texture: MEDIUM

Color After: YELLOW

Clarity After:

Artifacts:

Comments:

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P.44

INORGANIC ANALYSIS DATA SHEET

MBNJ58

Lab Name: CHEMTECH EDISON

Contract: 68-W00-088

Lab Code: CHEMED

Case No.: 28706

SAS No.:

SDG No.: MBNJ51

Matrix (soil/water): SOIL

Lab Sample ID: R1115-12S

Level (low/med): LOW

Date Received: 11/08/00

% Solids: 78.3

Concentration Units (ug/L or mg/Kg dry weight): MG/KG

| CAS No.   | Analyte   | Concentration | C | Q | M  |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum  | 13800         | - |   | P  |
| 7440-36-0 | Antimony  | 1.5           | U |   | P  |
| 7440-38-2 | Arsenic   | 5.4           |   |   | P  |
| 7440-39-3 | Barium    | 129           |   |   | P  |
| 7440-41-7 | Beryllium | 0.72          | B |   | P  |
| 7440-43-9 | Cadmium   | 0.50          | U |   | P  |
| 7440-70-2 | Calcium   | 2210          |   |   | P  |
| 7440-47-3 | Chromium  | 20.4          |   |   | P  |
| 7440-48-4 | Cobalt    | 12.6          |   |   | P  |
| 7440-50-8 | Copper    | 23.4          |   | E | P  |
| 7439-89-6 | Iron      | 26700         |   |   | P  |
| 7439-92-1 | Lead      | 19.5          |   |   | P  |
| 7439-95-4 | Magnesium | 4730          |   |   | P  |
| 7439-96-5 | Manganese | 364           |   |   | P  |
| 7439-97-6 | Mercury   | 0.12          | U |   | CV |
| 7440-02-0 | Nickel    | 38.8          |   |   | P  |
| 7440-09-7 | Potassium | 1200          | B | E | P  |
| 7782-49-2 | Selenium  | 1.8           |   |   | P  |
| 7440-22-4 | Silver    | 0.56          | B |   | P  |
| 7440-23-5 | Sodium    | 64.3          | U |   | P  |
| 7440-28-0 | Thallium  | 2.3           | B |   | P  |
| 7440-62-2 | Vanadium  | 20.6          |   |   | P  |
| 7440-66-6 | Zinc      | 108           |   |   | P  |
|           | Cyanide   | 0.19          | U |   | CA |

Color Before: BROWN

Clarity Before:

Texture: MEDIUM

Color After: YELLOW

Clarity After:

Artifacts:

Comments:

0.45



INORGANIC ANALYSIS DATA SHEET

MBNJ59

Lab Name: CHEMTECH EDISON

Contract: 68-W00-088

Lab Code: CHEMED

Case No.: 28706

SAS No.:

SDG No.: MBNJ51

Matrix (soil/water): SOIL

Lab Sample ID: R1115-13S

Level (low/med): LOW

Date Received: 11/08/00

% Solids: 74.1

Concentration Units (ug/L or mg/Kg dry weight): MG/KG

| CAS No.   | Analyte   | Concentration | C | Q | M  |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum  | 16200         |   |   | P  |
| 7440-36-0 | Antimony  | 1.6           | U |   | P  |
| 7440-38-2 | Arsenic   | 4.8           |   |   | P  |
| 7440-39-3 | Barium    | 157           |   |   | P  |
| 7440-41-7 | Beryllium | 0.84          | B |   | P  |
| 7440-43-9 | Cadmium   | 0.53          | U |   | P  |
| 7440-70-2 | Calcium   | 2260          |   |   | P  |
| 7440-47-3 | Chromium  | 24.0          |   |   | P  |
| 7440-48-4 | Cobalt    | 14.9          |   |   | P  |
| 7440-50-8 | Copper    | 25.1          |   | E | P  |
| 7439-89-6 | Iron      | 31000         |   |   | P  |
| 7439-92-1 | Lead      | 28.8          |   |   | P  |
| 7439-95-4 | Magnesium | 5190          |   |   | P  |
| 7439-96-5 | Manganese | 391           |   |   | P  |
| 7439-97-6 | Mercury   | 0.14          | U |   | CV |
| 7440-02-0 | Nickel    | 42.2          |   |   | P  |
| 7440-09-7 | Potassium | 1420          |   | E | P  |
| 7782-49-2 | Selenium  | 1.5           |   |   | P  |
| 7440-22-4 | Silver    | 0.73          | B |   | P  |
| 7440-23-5 | Sodium    | 138           | B |   | P  |
| 7440-28-0 | Thallium  | 0.98          | U |   | P  |
| 7440-62-2 | Vanadium  | 24.0          |   |   | P  |
| 7440-66-6 | Zinc      | 134           |   |   | P  |
|           | Cyanide   | 0.20          | U |   | CA |

Color Before: BROWN

Clarity Before:

Texture: MEDIUM

Color After: YELLOW

Clarity After:

Artifacts:

Comments:

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INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

MBNJ60

Lab Name: CHEMTECH EDISON

Contract: 68-W00-088

Lab Code: CHEMED

Case No.: 28706

SAS No.:

SDG No.: MBNJ51

Matrix (soil/water): SOIL

Lab Sample ID: R1115-14S

Level (low/med): LOW

Date Received: 11/08/00

% Solids:

76.6

Concentration Units (ug/L or mg/Kg dry weight): MG/KG

| CAS No.   | Analyte   | Concentration | C | Q | M  |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum  | 18300         | - |   | P  |
| 7440-36-0 | Antimony  | 1.6           | U |   | P  |
| 7440-38-2 | Arsenic   | 1.9           | B |   | P  |
| 7440-39-3 | Barium    | 179           |   |   | P  |
| 7440-41-7 | Beryllium | 0.89          | B |   | P  |
| 7440-43-9 | Cadmium   | 0.52          | U |   | P  |
| 7440-70-2 | Calcium   | 2250          |   |   | P  |
| 7440-47-3 | Chromium  | 26.1          |   |   | P  |
| 7440-48-4 | Cobalt    | 11.0          | B |   | P  |
| 7440-50-8 | Copper    | 19.9          |   | E | P  |
| 7439-89-6 | Iron      | 27300         |   |   | P  |
| 7439-92-1 | Lead      | 17.8          |   |   | P  |
| 7439-95-4 | Magnesium | 5830          |   |   | P  |
| 7439-96-5 | Manganese | 261           |   |   | P  |
| 7439-97-6 | Mercury   | 0.12          | U |   | CV |
| 7440-02-0 | Nickel    | 45.9          |   |   | P  |
| 7440-09-7 | Potassium | 1940          |   | E | P  |
| 7782-49-2 | Selenium  | 1.8           |   |   | P  |
| 7440-22-4 | Silver    | 0.43          | B |   | P  |
| 7440-23-5 | Sodium    | 161           | B |   | P  |
| 7440-28-0 | Thallium  | 1.2           | B |   | P  |
| 7440-62-2 | Vanadium  | 24.7          |   |   | P  |
| 7440-66-6 | Zinc      | 134           |   |   | P  |
|           | Cyanide   | 0.19          | U |   | CA |

Color Before: BROWN

Clarity Before:

Texture: MEDIUM

Color After: YELLOW

Clarity After:

Artifacts:

Comments:

p.47

INORGANIC ANALYSIS DATA SHEET

MBNJ62

Lab Name: CHEMTECH EDISON Contract: 68-W00-088  
 Lab Code: CHEMED Case No.: 28706 SAS No.: SDG No.: MBNJ51  
 Matrix (soil/water): SOIL Lab Sample ID: R1115-09S  
 Level (low/med): LOW Date Received: 11/08/00  
 % Solids: 76.3

Concentration Units (ug/L or mg/Kg dry weight): MG/KG

| CAS No.   | Analyte   | Concentration | C | Q | M  |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum  | 15400         |   |   | P  |
| 7440-36-0 | Antimony  | 1.5           | U |   | P  |
| 7440-38-2 | Arsenic   | 4.3           |   |   | P  |
| 7440-39-3 | Barium    | 182           |   |   | P  |
| 7440-41-7 | Beryllium | 0.85          | B |   | P  |
| 7440-43-9 | Cadmium   | 0.51          | U |   | P  |
| 7440-70-2 | Calcium   | 5220          |   |   | P  |
| 7440-47-3 | Chromium  | 21.8          |   |   | P  |
| 7440-48-4 | Cobalt    | 13.2          |   |   | P  |
| 7440-50-8 | Copper    | 34.4          |   | E | P  |
| 7439-89-6 | Iron      | 26000         |   |   | P  |
| 7439-92-1 | Lead      | 24.0          |   |   | P  |
| 7439-95-4 | Magnesium | 4810          |   |   | P  |
| 7439-96-5 | Manganese | 484           |   |   | P  |
| 7439-97-6 | Mercury   | 0.13          | U |   | CV |
| 7440-02-0 | Nickel    | 44.8          |   |   | P  |
| 7440-09-7 | Potassium | 1870          |   | E | P  |
| 7782-49-2 | Selenium  | 1.4           |   |   | P  |
| 7440-22-4 | Silver    | 0.42          | B |   | P  |
| 7440-23-5 | Sodium    | 66.6          | U |   | P  |
| 7440-28-0 | Thallium  | 0.95          | U |   | P  |
| 7440-62-2 | Vanadium  | 23.8          |   |   | P  |
| 7440-66-6 | Zinc      | 139           |   |   | P  |
|           | Cyanide   | 0.19          | U |   | CA |

Color Before: BROWN Clarity Before: Texture: MEDIUM  
 Color After: YELLOW Clarity After: Artifacts:

Comments:

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U.S. EPA - CLP

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EPA SAMPLE NO.

INORGANIC ANALYSIS DATA SHEET

MBNJ63

Lab Name: CHEMTECH EDISON

Contract: 68-W00-088

Lab Code: CHEMED

Case No.: 28706

SAS No.:

SDG No.: MBNJ51

Matrix (soil/water): SOIL

Lab Sample ID: R1115-10S

Level (low/med): LOW

Date Received: 11/08/00

% Solids: 78.2

Concentration Units (ug/L or mg/Kg dry weight): MG/KG

| CAS No.   | Analyte   | Concentration | C | Q | M  |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum  | 13500         |   |   | P  |
| 7440-36-0 | Antimony  | 1.5           | U |   | P  |
| 7440-38-2 | Arsenic   | 6.9           |   |   | P  |
| 7440-39-3 | Barium    | 114           |   |   | P  |
| 7440-41-7 | Beryllium | 0.76          | B |   | P  |
| 7440-43-9 | Cadmium   | 0.51          | U |   | P  |
| 7440-70-2 | Calcium   | 2450          |   |   | P  |
| 7440-47-3 | Chromium  | 19.2          |   |   | P  |
| 7440-48-4 | Cobalt    | 15.2          |   |   | P  |
| 7440-50-8 | Copper    | 23.9          |   | E | P  |
| 7439-89-6 | Iron      | 29800         |   |   | P  |
| 7439-92-1 | Lead      | 18.8          |   |   | P  |
| 7439-95-4 | Magnesium | 4250          |   |   | P  |
| 7439-96-5 | Manganese | 427           |   |   | P  |
| 7439-97-6 | Mercury   | 0.12          | U |   | CV |
| 7440-02-0 | Nickel    | 36.5          |   |   | P  |
| 7440-09-7 | Potassium | 1420          |   | E | P  |
| 7782-49-2 | Selenium  | 2.3           |   |   | P  |
| 7440-22-4 | Silver    | 0.58          | B |   | P  |
| 7440-23-5 | Sodium    | 108           | B |   | P  |
| 7440-28-0 | Thallium  | 0.94          | U |   | P  |
| 7440-62-2 | Vanadium  | 21.8          |   |   | P  |
| 7440-66-6 | Zinc      | 96.9          |   |   | P  |
|           | Cyanide   | 0.19          | U |   | CA |

Color Before: BROWN

Clarity Before:

Texture: MEDIUM

Color After: YELLOW

Clarity After:

Artifacts:

Comments:

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*p. 49*

INORGANIC ANALYSIS DATA SHEET

MBNJ64

Lab Name: CHEMTECH EDISON Contract: 68-W00-088  
 Lab Code: CHEMED Case No.: 28706 SAS No.: SDG No.: MBNJ51  
 Matrix (soil/water): SOIL Lab Sample ID: R1115-11S  
 Level (low/med): LOW Date Received: 11/08/00  
 % Solids: 79.7

Concentration Units (ug/L or mg/Kg dry weight): MG/KG

| CAS No.   | Analyte   | Concentration | C | Q | M  |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum  | 12200         | - |   | P  |
| 7440-36-0 | Antimony  | 1.5           | U |   | P  |
| 7440-38-2 | Arsenic   | 8.0           |   |   | P  |
| 7440-39-3 | Barium    | 89.8          |   |   | P  |
| 7440-41-7 | Beryllium | 0.58          | B |   | P  |
| 7440-43-9 | Cadmium   | 0.50          | U |   | P  |
| 7440-70-2 | Calcium   | 2000          |   |   | P  |
| 7440-47-3 | Chromium  | 17.6          |   |   | P  |
| 7440-48-4 | Cobalt    | 13.0          |   |   | P  |
| 7440-50-8 | Copper    | 22.0          |   | E | P  |
| 7439-89-6 | Iron      | 28300         |   |   | P  |
| 7439-92-1 | Lead      | 14.1          |   |   | P  |
| 7439-95-4 | Magnesium | 4000          |   |   | P  |
| 7439-96-5 | Manganese | 457           |   |   | P  |
| 7439-97-6 | Mercury   | 0.13          | U |   | CV |
| 7440-02-0 | Nickel    | 32.5          |   |   | P  |
| 7440-09-7 | Potassium | 1140          | B | E | P  |
| 7782-49-2 | Selenium  | 2.2           |   |   | P  |
| 7440-22-4 | Silver    | 0.50          | B |   | P  |
| 7440-23-5 | Sodium    | 92.9          | B |   | P  |
| 7440-28-0 | Thallium  | 2.2           | B |   | P  |
| 7440-62-2 | Vanadium  | 18.5          |   |   | P  |
| 7440-66-6 | Zinc      | 88.5          |   |   | P  |
|           | Cyanide   | 0.18          | U |   | CA |

Color Before: BROWN Clarity Before: Texture: MEDIUM  
 Color After: YELLOW Clarity After: Artifacts:

Comments:

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p. 50

INORGANIC ANALYSIS DATA SHEET

MBNJ66

Lab Name: CHEMTECH EDISON

Contract: 68-W00-088

Lab Code: CHEMED

Case No.: 28706

SAS No.:

SDG No.: MBNJ51

Matrix (soil/water): SOIL

Lab Sample ID: R1115-04S

Level (low/med): LOW

Date Received: 11/08/00

% Solids: 78.9

Concentration Units (ug/L or mg/Kg dry weight): MG/KG

| CAS No.   | Analyte   | Concentration | C | Q | M  |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum  | 18500         | - |   | P  |
| 7440-36-0 | Antimony  | 1.5           | U |   | P  |
| 7440-38-2 | Arsenic   | 2.2           | B |   | P  |
| 7440-39-3 | Barium    | 200           |   |   | P  |
| 7440-41-7 | Beryllium | 0.76          | B |   | P  |
| 7440-43-9 | Cadmium   | 0.50          | U |   | P  |
| 7440-70-2 | Calcium   | 2940          |   |   | P  |
| 7440-47-3 | Chromium  | 27.2          |   |   | P  |
| 7440-48-4 | Cobalt    | 14.6          |   |   | P  |
| 7440-50-8 | Copper    | 22.7          |   | E | P  |
| 7439-89-6 | Iron      | 27700         |   |   | P  |
| 7439-92-1 | Lead      | 18.1          |   |   | P  |
| 7439-95-4 | Magnesium | 5910          |   |   | P  |
| 7439-96-5 | Manganese | 754           |   |   | P  |
| 7439-97-6 | Mercury   | 0.12          | U |   | CV |
| 7440-02-0 | Nickel    | 47.4          |   |   | P  |
| 7440-09-7 | Potassium | 1630          |   | E | P  |
| 7782-49-2 | Selenium  | 1.0           | U |   | P  |
| 7440-22-4 | Silver    | 0.63          | B |   | P  |
| 7440-23-5 | Sodium    | 65.0          | U |   | P  |
| 7440-28-0 | Thallium  | 0.94          | B |   | P  |
| 7440-62-2 | Vanadium  | 23.6          |   |   | P  |
| 7440-66-6 | Zinc      | 136           |   |   | P  |
|           | Cyanide   | 0.18          | U |   | CA |

Color Before: BROWN

Clarity Before:

Texture: MEDIUM

Color After: YELLOW

Clarity After:

Artifacts:

Comments:

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p. 51

INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

Field Blank

MBNJ67

Lab Name: CHEMTECH EDISON

Contract: 68-W00-088

Lab Code: CHEMED

Case No.: 28706

SAS No.:

SDG No.: MBNJ51

Matrix (soil/water): WATER

Lab Sample ID: R1115-15S

Level (low/med): LOW

Date Received: 11/08/00

% Solids: 0.0

Concentration Units (ug/L or mg/Kg dry weight): UG/L

| CAS No.   | Analyte   | Concentration | C | Q | M  |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum  | 34.8          | U |   | P  |
| 7440-36-0 | Antimony  | 6.0           | U |   | P  |
| 7440-38-2 | Arsenic   | 5.7           | U |   | P  |
| 7440-39-3 | Barium    | 1.4           | U |   | P  |
| 7440-41-7 | Beryllium | 1.1           | U |   | P  |
| 7440-43-9 | Cadmium   | 2.0           | U |   | P  |
| 7440-70-2 | Calcium   | 21.1          | U |   | P  |
| 7440-47-3 | Chromium  | 3.7           | U |   | P  |
| 7440-48-4 | Cobalt    | 2.3           | U |   | P  |
| 7440-50-8 | Copper    | 1.0           | U |   | P  |
| 7439-89-6 | Iron      | 17.4          | U |   | P  |
| 7439-92-1 | Lead      | 3.0           | U |   | P  |
| 7439-95-4 | Magnesium | 26.5          | U |   | P  |
| 7439-96-5 | Manganese | 0.70          | U |   | P  |
| 7439-97-6 | Mercury   | 0.20          | U |   | CV |
| 7440-02-0 | Nickel    | 4.1           | U |   | P  |
| 7440-09-7 | Potassium | 17.8          | U |   | P  |
| 7782-49-2 | Selenium  | 4.0           | U |   | P  |
| 7440-22-4 | Silver    | 1.6           | U |   | P  |
| 7440-23-5 | Sodium    | 259           | U |   | P  |
| 7440-28-0 | Thallium  | 6.1           | B |   | P  |
| 7440-62-2 | Vanadium  | 3.6           | U |   | P  |
| 7440-66-6 | Zinc      | 2.5           | U |   | P  |
|           | Cyanide   | 2.9           | U |   | CA |

Color Before: COLORLESS

Clarity Before: CLEAR

Texture:

Color After: COLORLESS

Clarity After: CLEAR

Artifacts:

Comments:

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P. 52

# SAMPLE DELIVERY GROUP (SDG) TRAFFIC REPORT (TR) COVER SHEET

Lab Name: CHEMTECH EDISON Contract No.: 68-W00-088

Lab Code: CHEMED Case No.: 28706

Full Sample Analysis Price in Contract: \_\_\_\_\_

SDG No./First Sample in SDG. MBNJ51 Sample Receipt Date: 11/08/00  
(Lowest EPA Sample Number in first shipment of (MM/DD/YY)  
samples received under SDG.)

Last Sample in SDG: MBNJ67 Sample Receipt Date: 11/08/00  
(Highest EPA Sample Number in last shipment of (MM/DD/YY)  
samples received under SDG.)

EPA Sample Numbers in the SDG (listed in alphanumeric order)

|    |        |    |        |
|----|--------|----|--------|
| 1  | MBNJ51 | 12 | MBNJ63 |
| 2  | MBNJ52 | 13 | MBNJ64 |
| 3  | MBNJ53 | 14 | MBNJ66 |
| 4  | MBNJ54 | 15 | MBNJ67 |
| 5  | MBNJ55 | 16 |        |
| 6  | MBNJ56 | 17 |        |
| 7  | MBNJ57 | 18 |        |
| 8  | MBNJ58 | 19 |        |
| 9  | MBNJ59 | 20 |        |
| 10 | MBNJ60 | 21 |        |
| 11 | MBNJ62 | 22 |        |

Note: There are a maximum of 20 field samples in an SDG.

Attach Traffic Reports to this form in alphanumeric order  
(i.e.. the order listed on this form).

Sunny Patel  
Signature

11/08/00  
Date



REFERENCE NO. 15



**Roy F. Weston, Inc.**  
Raritan Plaza III, Suite 2B  
101 Fieldcrest Avenue  
Edison, New Jersey 08837-3622  
732-417-5800 • Fax 732-417-5801  
www.rfweston.com

8 January 2000

Mr. Dennis Munhall  
U.S. Environmental Protection Agency  
290 Broadway - 18th Floor  
New York, NY 10007-1866

**DOCUMENT CONTROL NO: SAT-1002.19**  
**SUBJECT: SAMPLING TRIP REPORT -**  
**LACKAWANNA FOUNDRY**

Dear Mr. Munhall:

Enclosed please find the Sampling Trip Report for samples collected during the Brownfields investigation conducted at the Lackawanna Foundry site on 5 December 2000.

If you have any questions, do not hesitate to call me at (732) 417-5806.

Very truly yours,

ROY F. WESTON, INC.

Donna Janda  
Project Manager

enclosure

cc: J. Bulich, Lackawanna Foundry  
cc: [unclear]



## SAMPLING TRIP REPORT

**SITE NAME:** Lackawanna Foundry  
DCN #: SAT-1002.19

**EPA I.D. NO.:** NYSFN0204209

**SAMPLING DATE:** December 5, 2000

1. Site Location: Refer to Figure 1
2. Sample Locations: Refer to Figure 2
3. Sample Descriptions: Refer to Table 1
4. Laboratories Receiving Samples:

| <u>Sample Type</u>                     | <u>Name and Address of Laboratory</u>                                      |
|--|--|
| Target Compound List<br>(TCL) Organics | U.S. EPA DESA Laboratory<br>2890 Woodbridge Avenue, Bldg 209<br>Edison, NJ |
| Target Analyte List (TAL)              | Same as above  |

5. Sample Dispatch Data:

Six aqueous samples for TCL volatiles, BNA, pesticide/PCB, and TAL inorganics analyses and one aqueous sample for volatile organic analysis only were hand delivered to the EPA DESA laboratory in Edison, NJ on 12/07/00.

6. On-Site Personnel:

| <u>Name</u>   | <u>Company</u> | <u>Duties on Site</u>                   |
|---------------|----------------|---|
| David Lewitt  | Region II SAT  | Site Health and Safety Officer, Sampler |
| Ernie Salazar | Region II SAT  | Sample Management Officer, Sampler      |

7. Additional Comments:

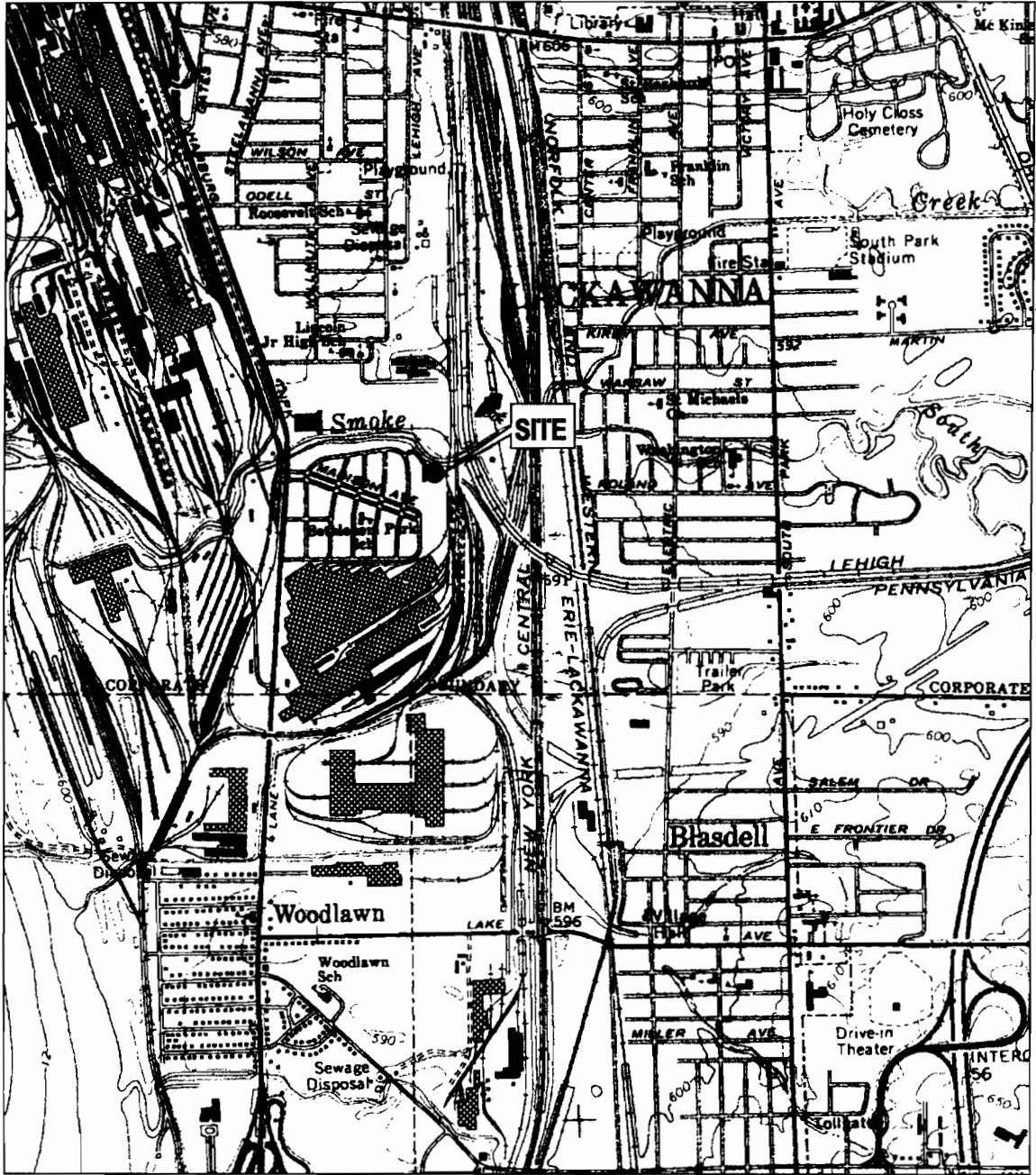
On December 5, 2000, the Region II Site Assessment Team (SAT) collected groundwater samples from five on-site monitoring wells at the Lackawanna Foundry site. These samples were collected as part of a Brownfields investigation conducted on site. All samples collected by Region II SAT

were analyzed for Target Compound List (TCL) and Target Analyte List (TAL), parameters by the U.S. EPA Region II Division of Environmental Science Assessment (DESA) Laboratory. The U.S. EPA Region 2 Laboratory Chain of Custody/Field Data Form is presented in Attachment 1.

8. Report Prepared by: *Dorna Janda* Date: 1/3/01  
Dorna Janda

9. Report Approved by: *W. S. Butterfield* Date: 1/8/01  
W. S. Butterfield, CHMM

**SITE LOCATION AND SAMPLE  
LOCATION MAPS**

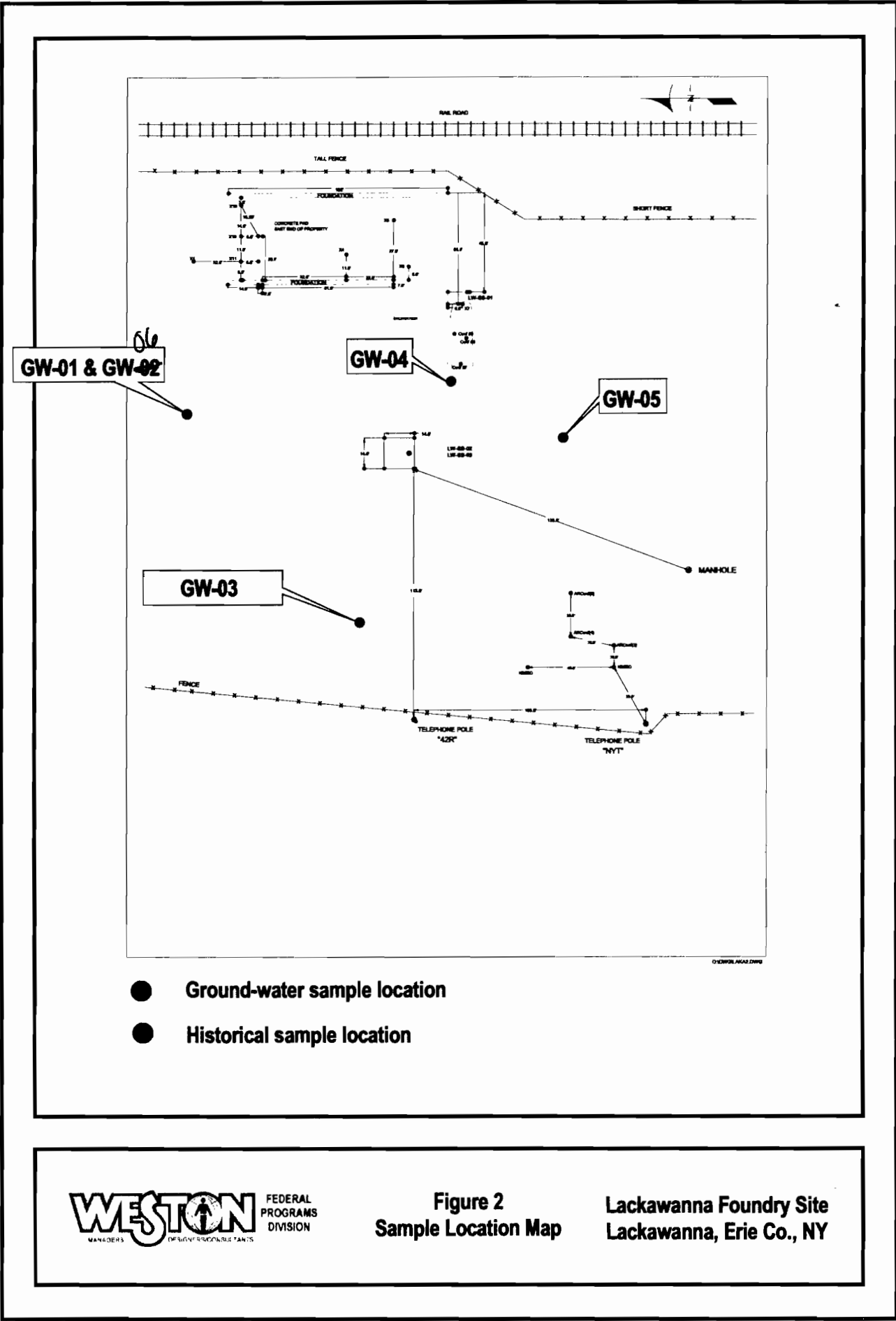


U.S. Geological Survey  
 7.5 Minute Series Topographical Map  
 Buffalo Southeast, NY Quadrangle  
 1948, Revised 1965



Lackawanna Foundry Site  
 Lackawanna, Erie Co., NY  
 Figure 1  
 Site Location Map

● Site Location



**Figure 2**  
**Sample Location Map**

**Lackawanna Foundry Site**  
**Lackawanna, Erie Co., NY**

**TABLE 1  
SAMPLE DESCRIPTIONS  
LACKAWANNA FOUNDRY  
LACKAWANNA, ERIE COUNTY, NEW YORK**

| <b>SAMPLE<br/>NUMBER<br/>1002-</b> | <b>DATE</b> | <b>TIME</b> | <b>COMMENTS</b>   |
|------------------------------------|-------------|-------------|---|
| GW01                               | 12/05/00    | 1400        | Groundwater sample collected from monitoring well MW-01. MS/MSD collected also. |
| GW03                               | 12/05/00    | 0855        | Groundwater sample collected from monitoring well MW-03.                        |
| GW04                               | 12/05/00    | 1100        | Groundwater sample collected from monitoring well MW-04.                        |
| GW05                               | 12/05/00    | 0955        | Groundwater sample collected from monitoring well MW-05.                        |
| GW06                               | 12/05/00    | 1400        | Duplicate sample of GW01.   |
| RB02                               | 12/05/00    | 1500        | Rinsate blank collected from decontaminated sampling pump (QA/QC sample).       |
| TB01                               | 12/04/00    | NA          | Trip blank (QA/QC sample).  |



**ATTACHMENT 1**

**TRAFFIC REPORTS/CHAIN OF CUSTODY RECORDS**

US EPA REGION 2 LABORATORY  
**CHAIN OF CUSTODY/ FIELD DATA FORM**

JRVEY NAME & LOCALITY Lock Haven County, Lock Haven, NY PROJECT LEADER Renee Daniels  
 PROGRAM: SF  : SITE ID \_\_\_\_\_ OPERABLE UNIT \_\_\_\_\_ PROGRAM RESULTS CODE \_\_\_\_\_  
 Permit # \_\_\_\_\_ RCRA  NPDES  SDWA  AM  CAA  TSCA  ENFORCEMENT: CRIMINAL  CIVIL

| LAB ID/ FIELD ID | CONTAINERS<br># OF | MATRIX | SPECIAL REQUIREMENTS?    | DESCRIPTION & INSTRUCTIONS INCLUDING LOCATION, ESTIMATED CONCENTRATIONS, SPECIAL REPORTING LIMITS, SPECIAL TEST REQUIREMENTS & ALIQUOTING | Preservative (circle) | Collection Time (24hr clock) |       | Collection Date mm/dd/yy |
|------------------|--------------------|--------|--------------------------|---|-----------------------|------------------------------|-------|--------------------------|
|                  |                    |        |                          |   |                       | . Begin                      | . End |                          |
| 002 - G2001      | 2                  | A      | <input type="checkbox"/> | VOA, BNA, PCB/Pest, Total Metals, Cyanide, 17.5/7.5/D   | 120050789             | 1400                         | 1405  | 12/15/00                 |
| 002 - G2003      | 8                  | A      | <input type="checkbox"/> | VOA, BNA, PCB/Pest, Total Metals, Cyanide   | 120050789             | 0855                         | 0900  | 12/15/00                 |
| 002 - G2004      | 8                  | A      | <input type="checkbox"/> | VOA, BNA, PCB/Pest, Total Metals, Cyanide   | 120050789             | 1100                         | 1105  | 12/15/00                 |
| 002 - G2005      | 8                  | A      | <input type="checkbox"/> | VOA, BNA, PCB/Pest, Total Metals, Cyanide   | 120050789             | 0955                         | 1000  | 12/15/00                 |
| 002 - G2006      | 8                  | A      | <input type="checkbox"/> | VOA, BNA, PCB/Pest, Total Metals, Cyanide   | 120050789             | 1400                         | 1405  | 12/15/00                 |
| 002 - Q2002      | 8                  | A      | <input type="checkbox"/> | VOA, BNA, PCB/Pest, Total Metals, Cyanide   | 120050789             | 1500                         | 1505  | 12/15/00                 |
| 002 - T2001      | 2                  | A      | <input type="checkbox"/> | VOA   | 123456789             | N/A                          | N/A   | 12/14/00                 |
|                  |                    |        | <input type="checkbox"/> |   | 123456789             |                              |       |                          |
|                  |                    |        | <input type="checkbox"/> |   | 123456789             |                              |       |                          |
|                  |                    |        | <input type="checkbox"/> |   | 123456789             |                              |       |                          |

COMMENTS: Roy Feinston

Preservative  
 1=ice  
 2=H2SO4 pH<2  
 3=HNO3 pH<2  
 4=HCl pH<2  
 5=Na2S2O3  
 6=NaOH pH>9  
 7=Ascorbic Acid  
 8 = FAS  
 9=ZnAc

Person Assuming Responsibility for Sample(s):

|                       |                  |              |
|-----------------------|------------------|--------------|
| aqueous               | Relinquished By: | Received By: |
| aqueous (chlorinated) | Relinquished By: | Received By: |
| soil                  | Relinquished By: | Received By: |
| sediment              | Relinquished By: | Received By: |
| sludge                | Relinquished By: | Received By: |

Survey Complete? Y  N

REFERENCE NO. 16

***NYS Department of Environmental Conservation - Home - Site Map - Search***Division of Water**Technical & Operational Guidance Series (TOGS)**

The documents listed on this page are in Adobe Acrobat PDF format. Viewing these PDF files requires that you have Adobe Acrobat Reader software (available at no cost from Adobe) installed on your computer. If you need a paper copy of these documents, call (518) 457-7464 or follow the Ordering Information at the bottom of this page.

**Table Of Contents**SPDES   Groundwater   Water Quality   Other Technical Subjects**1.0 SPDES****1.1 Standards**

- \* 1.1.1 Ambient Water Quality Standards and Guidance Values and Groundwater Effluent Limitations. 6/98 Ed. pdf file size 464k - Print Copy from DEC  
Cost \$15.50.  
Errata Sheet for above added 1/99  
April 2000 addendum
- \* 1.1.3 Procedures for Derivation of Site-Specific Standards and Guidance Values for Protection of Aquatic Life.  
2/98 Ed. pdf file size 73k - Print Copy from DEC  
Cost \$0.50
- \* 1.1.4 Procedures for Derivation of Bioaccumulation Factors.  
2/98 Ed. pdf file size 52k - Print Copy from DEC  
Cost \$2.50
- \* 1.1.5 Procedures for Deriving Ambient Water Quality Standards and Guidance Values for the Protection of Wildlife.  
2/98 Ed. pdf file size 126k - Print Copy from DEC  
Cost \$1.75

**1.2 Permit Administration**

- \* 1.2.1 Industrial Permit Writing  
2/98 Ed. pdf file size 373k - Print Copy from DEC  
Cost \$15.00
- 1.2.2 SPDES Permit Program Priorities and Definitions  
5/90 Ed. pdf file size 48k - Print Copy from DEC  
Cost \$3.00
- 1.2.3 Decentralization of Technical Permit Drafting Authority  
4/90 Ed. pdf file size 11k - Print Copy from DEC  
Cost \$0.50
- 1.2.4 Individual Sewage Treatment System Discharges to Surface Waters  
10/90 Ed. pdf file size 10k - Print Copy from DEC  
Cost \$0.50

**1.3 Permit Development**

- \* 1.3.1 Total Maximum Daily Loads & Water Quality-Based Effluent Limits  
2/98 Ed. pdf file size 60k - Print Copy from DEC  
Cost \$2.00
- 1.3.1A Amendment - Organic Substances

REFERENCE NO. 17



UNITED STATES ENVIRONMENTAL PROTECTION AGENCY  
REGION II  
EDISON, NEW JERSEY 08837

30 JAN 2001

Roy F. Weston Inc.  
Raritan Plaza 3, Suite 2B  
101 Fieldcrest Avenue  
Edison, NJ 08837  
Attn: Donna Janda

JAN 31 2001

Dear Ms. Janda:

Enclosed are the results of the Lackawanna Foundry sampling survey conducted by your firm during the week of December 4, 2000. These data include the results of seven aqueous samples for metals, cyanide and volatile and non-volatile organic analyses. Any correspondence concerning these results should refer to our internal project numbers, 00120005 and 00120006, to uniquely identify the data. Please refer to the first page of each report for a description of any remark codes used as data qualifiers. It should be noted that all data are considered to be EPA-validated.

If you have any questions you can contact me by phone at (732) 906-6886, by fax at (732) 321-6613 or via the Internet at "birri.john@epa.gov".

Sincerely,

A handwritten signature in cursive script that reads "John Birri".

John Birri  
Special Projects Coordinator  
Laboratory Branch

Enclosure



**Data Report: Lackawanna Foundry**

**Project Number: 00120005**

Program: Y206

Project Leader: Donna Janda

| Codes | Explanation  |
|-------|--|
| B     | RESULTS BASED UPON COLONY COUNTS OUTSIDE ACCEPTABLE RANGE        |
| J     | ESTIMATED VALUE  |
| K     | ACTUAL VALUE KNOWN TO BE LESS THAN VALUE GIVEN                   |
| L     | ACTUAL VALUE KNOWN TO BE GREATER THAN VALUE GIVEN                |
| N     | SEE NARRATIVE COMMENTS   |
| V     | SAMPLE RECEIVED BUT NOT ANALYZED DUE TO LAB ACCIDENT             |
| U     | REPORTING LIMIT  |
| QD    | ACCURACY CHECK SAMPLE ABOVE UPPER ACCEPTANCE LIMIT               |
| QE    | ACCURACY CHECK SAMPLE BELOW LOWER ACCEPTANCE LIMIT               |
| QF    | PRECISION OF CALIBRATION CURVE LESS THAN ACCEPTANCE CRITERIA     |
| QJ    | REPORTING LIMIT ESTIMATED DUE TO INTERFERENCE                    |
| QG    | CONTINUING CALIBRATION CHECK DOES NOT MEET ACCEPTANCE CRITERIA   |
| QS    | SPIKE RECOVERIES ABOVE UPPER ACCEPTANCE LIMIT                    |
| QR    | SPIKE RECOVERIES BELOW LOWER ACCEPTANCE LIMIT                    |
| QP    | SAMPLE REPLICATE PRECISION DOES NOT MEET ACCEPTANCE CRITERIA     |
| QH    | RECOMMENDED HOLDING TIME EXCEEDED                                |
| QT    | TENTATIVELY IDENTIFIED COMPOUND                                  |
| QM    | PRESENCE OF MATERIAL VERIFIED BUT NOT QUANTIFIED                 |
| QB    | BLANK CONTAMINATED WITH ANALYTE IN EXCESS OF ACCEPTANCE CRITERIA |
| QQ    | SAMPLE IMPROPERLY MAINTAINED                                     |



**Survey Name: Lackawanna Foundry**

Project Number: 00120005

\*Sorted By Sample ID

**AB03259**

Field/Station ID: 1002-GW01

Coll. Ending Date/Time: 12/5/00 14:05

Matrix: Aqueous

Collection Begin Date/Time: 12/05/00 14:00

Sample Description:

**Analysis Type: VOA TCL GCMS AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u>             | <u>Result</u> | <u>Units</u> | <u>Codes</u> | <u>Remark</u> |
|-------------------|---------------------------------|---------------|--------------|--------------|---------------|
| 000074873         | CHLOROMETHANE                   | 10            | ug/L         | U            |               |
| 000075014         | VINYL CHLORIDE                  | 10            | ug/L         | U            |               |
| 000074839         | BROMOMETHANE                    | 10            | ug/L         | U            |               |
| 000075003         | CHLOROETHANE                    | 10            | ug/L         | U            |               |
| 000075354         | 1,1-DICHLOROETHENE              | 10            | ug/L         | U            |               |
| 000075150         | CARBON DISULFIDE                | 10            | ug/L         | U            |               |
| 000067641         | ACETONE                         | 10            | ug/L         | U            |               |
| 000075092         | METHYLENE CHLORIDE              | 10            | ug/L         | U            |               |
| 000156605         | TRANS-1,2-DICHLOROETHENE        | 10            | ug/L         | U            |               |
| 000156592         | CIS-1,2-DICHLOROETHENE          | 10            | ug/L         | U            |               |
| 000075343         | 1,1-DICHLOROETHANE              | 10            | ug/L         | U            |               |
| 000078933         | 2-BUTANONE                      | 10            | ug/L         | U            |               |
| 000067663         | CHLOROFORM                      | 10            | ug/L         | U            |               |
| 000107062         | 1,2-DICHLOROETHANE              | 10            | ug/L         | U            |               |
| 000071556         | 1,1,1-TRICHLOROETHANE           | 10            | ug/L         | U            |               |
| 000056235         | CARBON TETRACHLORIDE            | 10            | ug/L         | U            |               |
| 000071432         | BENZENE                         | 10            | ug/L         | U            |               |
| 025323891         | TRICHLOROETHENE                 | 10            | ug/L         | U            |               |
| 000078875         | 1,2-DICHLOROPROPANE             | 10            | ug/L         | U            |               |
| 000074975         | BROMODICHLOROMETHANE            | 10            | ug/L         | U            |               |
| 010061015         | CIS-1,3-DICHLOROPROPENE         | 10            | ug/L         | U            |               |
| 010061026         | TRANS-1,3-DICHLOROPROPENE       | 10            | ug/L         | U            |               |
| 000079005         | 1,1,2-TRICHLOROETHANE           | 10            | ug/L         | U            |               |
| 000124481         | DIBROMOCHLOROMETHANE            | 10            | ug/L         | U            |               |
| 000075252         | BROMOFORM                       | 10            | ug/L         | U            |               |
| 000108101         | 4-METHYL-2-PENTANONE            | 10            | ug/L         | U            |               |
| 000108883         | TOLUENE                         | 10            | ug/L         | U            |               |
| 000127184         | TETRACHLOROETHENE               | 10            | ug/L         | U            |               |
| 000591786         | 2-HEXANONE                      | 10            | ug/L         | U            |               |
| 000108907         | CHLOROBENZENE                   | 10            | ug/L         | U            |               |
| 000100414         | ETHYLBENZENE                    | 10            | ug/L         | U            |               |
| 001330207         | M,P-XYLENE                      | 10            | ug/L         | U            |               |
| 000095476         | O-XYLENE                        | 10            | ug/L         | U            |               |
| 000100425         | STYRENE                         | 10            | ug/L         | U            |               |
| 000079345         | 1,1,2,2-TETRACHLOROETHANE       | 10            | ug/L         | U            |               |
|                   | UNKNOWN HYDROCARBON ,RT =3.04   | 19            | ug/L         | J QT         |               |
|                   | UNKNOWN HYDROCARBON ,RT =3.48   | 27            | ug/L         | J QT         |               |
|                   | BUTANE, 2-METHYL- ,RT =4.63     | 31            | ug/L         | J QT         |               |
|                   | PENTANE ,RT =5.14               | 6.0           | ug/L         | J QT         |               |
|                   | BUTANE, 2,3-DIMETHYL- ,RT =6.73 | 9.0           | ug/L         | J QT         |               |
|                   | UNKNOWN ,RT =7.23               | 6.0           | ug/L         | J QT         |               |





**Survey Name: Lackawanna Foundry**

Project Number: 00120005

\*Sorted By Sample

**AB03259** Field/Station ID: 1002-GW01  
Matrix: Aqueous  
Sample Description:

Coll. Ending Date/Time: 12/5/00 14:05  
Collection Begin Date/Time: 12/05/00 14:00

**Analysis Type: VOA TCL GCMS AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|---------------------|---------------|--------------|---------------------|
|                   | UNKNOWN ,RT =8.59   | 5.0           | ug/L         | J QT                |

**Analysis Type: NVOA GCMS AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u>         | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|-----------------------------|---------------|--------------|---------------------|
| 000108952         | PHENOL                      | 4.0           | ug/L         | U                   |
| 000111444         | BIS(2-CHLOROETHYL)ETHER     | 4.0           | ug/L         | U                   |
| 000095578         | 2-CHLOROPHENOL              | 4.0           | ug/L         | U                   |
| 000541731         | 1,3-DICHLOROBENZENE         | 4.0           | ug/L         | U                   |
| 000106467         | 1,4-DICHLOROBENZENE         | 4.0           | ug/L         | U                   |
| 000095501         | 1,2-DICHLOROBENZENE         | 4.0           | ug/L         | U                   |
| 000100516         | BENZYL ALCOHOL              | 10            | ug/L         | U                   |
| 000095487         | 2-METHYLPHENOL              | 4.0           | ug/L         | U                   |
| 000108601         | BIS(2-CHLOROISOPROPYL)ETHER | 4.0           | ug/L         | U                   |
| 000106445         | 4-METHYLPHENOL              | 4.0           | ug/L         | U                   |
| 000621647         | N-NITROSO-DI-N-PROPYLAMINE  | 4.0           | ug/L         | U                   |
| 000067721         | HEXACHLOROETHANE            | 4.0           | ug/L         | U                   |
| 000098953         | NITROBENZENE                | 4.0           | ug/L         | U                   |
| 000078591         | ISOPHORONE                  | 4.0           | ug/L         | U                   |
| 000088755         | 2-NITROPHENOL               | 4.0           | ug/L         | U                   |
| 000105679         | 2,4-DIMETHYLPHENOL          | 4.0           | ug/L         | U                   |
| 000111911         | BIS(2-CHLOROETHOXY)METHANE  | 4.0           | ug/L         | U                   |
| 000120832         | 2,4-DICHLOROPHENOL          | 4.0           | ug/L         | U                   |
| 000120821         | 1,2,4-TRICHLOROBENZENE      | 4.0           | ug/L         | U                   |
| 000065850         | BENZOIC ACID                | 49            | ug/L         | U                   |
| 000091203         | NAPHTHALENE                 | 4.0           | ug/L         | U                   |
| 000106478         | 4-CHLOROANILINE             | 49            | ug/L         | U                   |
| 000087683         | HEXACHLOROBUTADIENE         | 4.0           | ug/L         | U                   |
| 000059507         | 4-CHLORO-3-METHYLPHENOL     | 4.0           | ug/L         | U                   |
| 000091576         | 2-METHYL NAPHTHALENE        | 4.0           | ug/L         | U                   |
| 000077474         | HEXACHLOROCYCLOPENTADIENE   | 4.0           | ug/L         | U                   |
| 000088062         | 2,4,6-TRICHLOROPHENOL       | 4.0           | ug/L         | U                   |
| 000095954         | 2,4,5-TRICHLOROPHENOL       | 4.0           | ug/L         | U                   |
| 000091587         | 2-CHLORONAPHTHALENE         | 4.0           | ug/L         | U                   |
| 000088744         | 2-NITROANILINE              | 4.0           | ug/L         | U                   |
| 000131113         | DIMETHYL PHTHALATE          | 4.0           | ug/L         | U                   |
| 000208968         | ACENAPHTHYLENE              | 4.0           | ug/L         | U                   |
| 000606202         | 2,6-DINITROTOLUENE          | 4.0           | ug/L         | U                   |
| 000099092         | 3-NITROANILINE              | 49            | ug/L         | U                   |
| 000083329         | ACENAPHTHENE                | 4.0           | ug/L         | U                   |
| 000051285         | 2,4-DINITROPHENOL           | 10            | ug/L         | U                   |
| 000100027         | 4-NITROPHENOL               | 4.0           | ug/L         | U                   |
| 000132649         | DIBENZOFURAN                | 4.0           | ug/L         | U                   |



**Survey Name: Lackawanna Foundry**

Project Number: 00120005

\*Sorted By Sample ID

**AB03259** Field/Station ID: 1002-GW01

Coll. Ending Date/Time: 12/5/00 14:05

Matrix: Aqueous

Collection Begin Date/Time: 12/05/00 14:00

Sample Description:

**Analysis Type: NVOA GCMS AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u>                   | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|---------------------------------------|---------------|--------------|---------------------|
| 000121142         | 2,4-DINITROTOLUENE                    | 4.0           | ug/L         | U                   |
| 000086737         | FLUORENE                              | 4.0           | ug/L         | U                   |
| 000084662         | DIETHYLPHTHALATE                      | 4.0           | ug/L         | U                   |
| 007005723         | 4-CHLOROPHENYL PHENYL ETHER           | 4.0           | ug/L         | U                   |
| 000100016         | 4-NITROANILINE                        | 4.0           | ug/L         | U                   |
| 000534521         | 4,6-DINITRO-2-METHYLPHENOL            | 4.0           | ug/L         | U                   |
| 000086306         | N-NITROSODIPHENYLAMINE                | 4.0           | ug/L         | U                   |
| 000103333         | DIAZENE, DIPHENYL                     | 4.0           | ug/L         | U                   |
| 000101553         | 4-BROMOPHENYL PHENYL ETHER            | 4.0           | ug/L         | U                   |
| 000118741         | HEXACHLORO BENZENE                    | 4.0           | ug/L         | U                   |
| 000087865         | PENTACHLOROPHENOL                     | 10            | ug/L         | U                   |
| 000085018         | PHENANTHRENE                          | 4.0           | ug/L         | U                   |
| 000120127         | ANTHRACENE                            | 4.0           | ug/L         | U                   |
| 000084742         | DI-N-BUTYL PHTHALATE                  | 4.0           | ug/L         | U                   |
| 000206440         | FLUORANTHENE                          | 4.0           | ug/L         | U                   |
| 000129000         | PYRENE                                | 4.0           | ug/L         | U                   |
| 000085687         | BUTYL BENZYL PHTHALATE                | 4.0           | ug/L         | U                   |
| 000056553         | BENZO(A)ANTHRACENE                    | 4.0           | ug/L         | U                   |
| 000218019         | CHRYSENE                              | 4.0           | ug/L         | U                   |
| 000117817         | BIS(2-ETHYLHEXYL)PHTHALATE            | 4.0           | ug/L         | U                   |
| 000117840         | DI-N-OCTYL PHTHALATE                  | 4.0           | ug/L         | U                   |
| 000205992         | BENZO(B)FLUORANTHENE                  | 4.0           | ug/L         | U                   |
| 000207089         | BENZO(K)FLUORANTHENE                  | 4.0           | ug/L         | U                   |
| 000050328         | BENZO(A)PYRENE                        | 4.0           | ug/L         | U                   |
| 000193395         | INDENO(1,2,3-CD)PYRENE                | 4.0           | ug/L         | U                   |
| 000053703         | DIBENZO(A,H)ANTHRACENE                | 4.0           | ug/L         | U                   |
| 000191242         | BENZO(G,H,I)PERYLENE                  | 4.0           | ug/L         | U                   |
|                   | CYCLOHEXANOL, 1-METHYL-, RT =7.44     | 3.0           | ug/L         | J QT                |
|                   | UNKNOWN ,RT =8.45                     | 4.0           | ug/L         | J QT                |
|                   | UNKNOWN ,RT =8.95                     | 4.0           | ug/L         | J QT                |
|                   | UNKNOWN SUBSTITUTED BENZENE ,RT =9.48 | 2.0           | ug/L         | J QT                |
|                   | UNKNOWN ,RT =9.53                     | 2.0           | ug/L         | J QT                |

**Analysis Type: PESTICIDES TCL GC AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|---------------------|---------------|--------------|---------------------|
| 319846            | ALPHA-BHC           | 0.010         | ug/L         | U                   |
| 58899             | GAMMA-BHC           | 0.010         | ug/L         | U                   |
| 319857            | BETA-BHC            | 0.010         | ug/L         | U                   |
| 319868            | DELTA-BHC           | 0.010         | ug/L         | U                   |
| 76448             | HEPTACHLOR          | 0.010         | ug/L         | U                   |
| 309002            | ALDRIN              | 0.010         | ug/L         | U                   |
| 1024573           | HEPTACHLOR EPOXIDE  | 0.010         | ug/L         | U                   |



**Survey Name: Lackawanna Foundry**

Project Number: 00120005

\*Sorted By Sample

**AB03259**

Field/Station ID: 1002-GW01

Coll. Ending Date/Time: 12/5/00 14:05

Matrix: Aqueous

Collection Begin Date/Time: 12/05/00 14:00

Sample Description:

**Analysis Type: PESTICIDES TCL GC AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|---------------------|---------------|--------------|---------------------|
| 5566347           | GAMMA-CHLORDANE     | 0.010         | ug/L         | U                   |
| 5103719           | ALPHA-CHLORDANE     | 0.010         | ug/L         | U                   |
| 115297            | ENDOSULFAN I        | 0.010         | ug/L         | U                   |
| 72559             | 4,4'-DDE            | 0.010         | ug/L         | U                   |
| 60571             | DIELDRIN            | 0.010         | ug/L         | U                   |
| 72208             | ENDRIN              | 0.010         | ug/L         | U                   |
| 72548             | 4,4'-DDD            | 0.010         | ug/L         | U                   |
| 33213659          | ENDOSULFAN II       | 0.010         | ug/L         | U                   |
| 50293             | 4,4'-DDT            | 0.010         | ug/L         | U                   |
| 7421934           | ENDRIN ALDEHYDE     | 0.010         | ug/L         | U                   |
| 72435             | METHOXYCHLOR        | 0.050         | ug/L         | U                   |
| 1031078           | ENDOSULFAN SULFATE  | 0.010         | ug/L         | U                   |
| 53494705          | ENDRIN KETONE       | 0.010         | ug/L         | U                   |
| 57749             | CHLORDANE           | 0.10          | ug/L         | U                   |
| 8001352           | TOXAPHENE           | 0.10          | ug/L         | U                   |

**Analysis Type: PCBS TCL GC AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|---------------------|---------------|--------------|---------------------|
| 12674-11-2        | AROCLOR 1016        | 0.10          | ug/L         | U                   |
| 011104282         | AROCLOR 1221        | 0.20          | ug/L         | U                   |
| 011141165         | AROCLOR 1232        | 0.10          | ug/L         | U                   |
| 53469-21-9        | AROCLOR 1242        | 0.10          | ug/L         | U                   |
| 012672296         | AROCLOR 1248        | 0.10          | ug/L         | U                   |
| 11097-69-1        | AROCLOR 1254        | 0.10          | ug/L         | U                   |
| 11096-82-5        | AROCLOR 1260        | 0.10          | ug/L         | U                   |

**Analysis Type: METALS TAL ICP AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|---------------------|---------------|--------------|---------------------|
| 007440224         | SILVER              | 10            | ug/L         | U                   |
| 007429905         | ALUMINUM            | 200           | ug/L         | U                   |
| 007440393         | BARIUM              | 380           | ug/L         | U                   |
| 007440417         | BERYLLIUM           | 5.0           | ug/L         | U                   |
| 007440702         | CALCIUM             | 250,000       | ug/L         | U                   |
| 007440439         | CADMIUM             | 5.0           | ug/L         | U                   |
| 007440484         | COBALT              | 50            | ug/L         | U                   |
| 007440473         | CHROMIUM            | 10            | ug/L         | U                   |
| 007440508         | COPPER              | 25            | ug/L         | U                   |
| 007439896         | IRON                | 13,000        | ug/L         | U                   |
| 007440097         | POTASSIUM           | 8,100         | ug/L         | U                   |
| 007439954         | MAGNESIUM           | 130,000       | ug/L         | U                   |
| 007439965         | MANGANESE           | 300           | ug/L         | U                   |
| 007440235         | SODIUM              | 540,000       | ug/L         | U                   |
| 007440020         | NICKEL              | 40            | ug/L         | U                   |



**Survey Name: Lackawanna Foundry**

Project Number: 00120005

\*Sorted By Sample ID

**AB03259**

Field/Station ID: 1002-GW01

Coll. Ending Date/Time: 12/5/00 14:05

Matrix: Aqueous

Collection Begin Date/Time: 12/05/00 14:00

Sample Description:

**Analysis Type: METALS TAL ICP AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Codes</u> | <u>Remark</u> |
|-------------------|---------------------|---------------|--------------|--------------|---------------|
| 007440360         | ANTIMONY            | 60            | ug/L         | U            |               |
| 007440622         | VANADIUM            | 50            | ug/L         | U            |               |
| 007440666         | ZINC                | 20            | ug/L         | U            |               |

**Single Component Analyses**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Codes</u> | <u>Remark</u> |
|-------------------|---------------------|---------------|--------------|--------------|---------------|
| 7440-38-2         | ARSENIC             | 10            | ug/L         | U            |               |
| 7439-92-1         | LEAD                | 3.0           | ug/L         | U            |               |
| 7782-49-2         | SELENIUM            | 5.0           | ug/L         | U            |               |
| 7440-28-0         | THALLIUM            | 10            | ug/L         | U            | QR            |

**Single Component Analyses**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Codes</u> | <u>Remark</u> |
|-------------------|---------------------|---------------|--------------|--------------|---------------|
| 007439976         | MERCURY (ESAT)      | 0.20          | ug/L         | U            | QR            |

**AB03260**

Field/Station ID: 1002-GW03

Coll. Ending Date/Time: 12/5/00 09:00

Matrix: Aqueous

Collection Begin Date/Time: 12/05/00 08:55

Sample Description:

**Analysis Type: VOA TCL GCMS AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u>       | <u>Result</u> | <u>Units</u> | <u>Codes</u> | <u>Remark</u> |
|-------------------|---------------------------|---------------|--------------|--------------|---------------|
| 000074873         | CHLOROMETHANE             | 10            | ug/L         | U            |               |
| 000075014         | VINYL CHLORIDE            | 10            | ug/L         | U            |               |
| 000074839         | BROMOMETHANE              | 10            | ug/L         | U            |               |
| 000075003         | CHLOROETHANE              | 10            | ug/L         | U            |               |
| 000075354         | 1,1-DICHLOROETHENE        | 10            | ug/L         | U            |               |
| 000075150         | CARBON DISULFIDE          | 10            | ug/L         | U            |               |
| 000067641         | ACETONE                   | 11            | ug/L         |              |               |
| 000075092         | METHYLENE CHLORIDE        | 10            | ug/L         | U            |               |
| 000156605         | TRANS-1,2-DICHLOROETHENE  | 10            | ug/L         | U            |               |
| 000156592         | CIS-1,2-DICHLOROETHENE    | 10            | ug/L         | U            |               |
| 000075343         | 1,1-DICHLOROETHANE        | 10            | ug/L         | U            |               |
| 000078933         | 2-BUTANONE                | 10            | ug/L         | U            |               |
| 000067663         | CHLOROFORM                | 10            | ug/L         | U            |               |
| 000107062         | 1,2-DICHLOROETHANE        | 10            | ug/L         | U            |               |
| 000071556         | 1,1,1-TRICHLOROETHANE     | 10            | ug/L         | U            |               |
| 000056235         | CARBON TETRACHLORIDE      | 10            | ug/L         | U            |               |
| 000071432         | BENZENE                   | 10            | ug/L         | U            |               |
| 025323891         | TRICHLOROETHENE           | 10            | ug/L         | U            |               |
| 000078875         | 1,2-DICHLOROPROPANE       | 10            | ug/L         | U            |               |
| 000074975         | BROMODICHLOROMETHANE      | 10            | ug/L         | U            |               |
| 010061015         | CIS-1,3-DICHLOROPROPENE   | 10            | ug/L         | U            |               |
| 010061026         | TRANS-1,3-DICHLOROPROPENE | 10            | ug/L         | U            |               |



Survey Name: Lackawanna Foundry

Project Number: 00120005

\*Sorted By Sample

**AB03260**

Field/Station ID: 1002-GW03

Coll. Ending Date/Time: 12/5/00 09:00

Matrix: Aqueous

Collection Begin Date/Time: 12/05/00 08:55

Sample Description:

**Analysis Type: VOA TCL GCMS AQUEOUS**

| CAS Number | Analyte Name              | Result | Units | Remark Codes |
|------------|---------------------------|--------|-------|--------------|
| 000079005  | 1,1,2-TRICHLOROETHANE     | 10     | ug/L  | U            |
| 000124481  | DIBROMOCHLOROMETHANE      | 10     | ug/L  | U            |
| 000075252  | BROMOFORM                 | 10     | ug/L  | U            |
| 000108101  | 4-METHYL-2-PENTANONE      | 10     | ug/L  | U            |
| 000108883  | TOLUENE                   | 10     | ug/L  | U            |
| 000127184  | TETRACHLOROETHENE         | 10     | ug/L  | U            |
| 000591786  | 2-HEXANONE                | 10     | ug/L  | U            |
| 000108907  | CHLOROBENZENE             | 10     | ug/L  | U            |
| 000100414  | ETHYLBENZENE              | 10     | ug/L  | U            |
| 001330207  | M,P-XYLENE                | 10     | ug/L  | U            |
| 000095476  | O-XYLENE                  | 10     | ug/L  | U            |
| 000100425  | STYRENE                   | 10     | ug/L  | U            |
| 000079345  | 1,1,2,2-TETRACHLOROETHANE | 10     | ug/L  | U            |

**Analysis Type: NVOA GCMS AQUEOUS**

| CAS Number | Analyte Name                | Result | Units | Remark Codes |
|------------|-----------------------------|--------|-------|--------------|
| 000108952  | PHENOL                      | 4.0    | ug/L  | U            |
| 000111444  | BIS(2-CHLOROETHYL)ETHER     | 4.0    | ug/L  | U            |
| 000095578  | 2-CHLOROPHENOL              | 4.0    | ug/L  | U            |
| 000541731  | 1,3-DICHLOROBENZENE         | 4.0    | ug/L  | U            |
| 000106467  | 1,4-DICHLOROBENZENE         | 4.0    | ug/L  | U            |
| 000095501  | 1,2-DICHLOROBENZENE         | 4.0    | ug/L  | U            |
| 000100516  | BENZYL ALCOHOL              | 10     | ug/L  | U            |
| 000095487  | 2-METHYLPHENOL              | 4.0    | ug/L  | U            |
| 000108601  | BIS(2-CHLOROISOPROPYL)ETHER | 4.0    | ug/L  | U            |
| 000106445  | 4-METHYLPHENOL              | 4.0    | ug/L  | U            |
| 000621647  | N-NITROSO-DI-N-PROPYLAMINE  | 4.0    | ug/L  | U            |
| 000067721  | HEXACHLOROETHANE            | 4.0    | ug/L  | U            |
| 000098953  | NITROBENZENE                | 4.0    | ug/L  | U            |
| 000078591  | ISOPHORONE                  | 4.0    | ug/L  | U            |
| 000088755  | 2-NITROPHENOL               | 4.0    | ug/L  | U            |
| 000105679  | 2,4-DIMETHYLPHENOL          | 4.0    | ug/L  | U            |
| 000111911  | BIS(2-CHLOROETHOXY)METHANE  | 4.0    | ug/L  | U            |
| 000120832  | 2,4-DICHLOROPHENOL          | 4.0    | ug/L  | U            |
| 000120821  | 1,2,4-TRICHLOROBENZENE      | 4.0    | ug/L  | U            |
| 000065850  | BENZOIC ACID                | 48     | ug/L  | U            |
| 000091203  | NAPHTHALENE                 | 4.0    | ug/L  | U            |
| 000106478  | 4-CHLOROANILINE             | 48     | ug/L  | U            |
| 000087683  | HEXACHLOROBUTADIENE         | 4.0    | ug/L  | U            |
| 000059507  | 4-CHLORO-3-METHYLPHENOL     | 4.0    | ug/L  | U            |
| 000091576  | 2-METHYL NAPHTHALENE        | 4.0    | ug/L  | U            |
| 000077474  | HEXACHLOROCYCLOPENTADIENE   | 4.0    | ug/L  | U            |



U.S. EPA Region 2 Laboratory  
Data Report

Survey Name: Lackawanna Foundry

Project Number: 00120005

\*Sorted By Sample ID

**AB03260** Field/Station ID: 1002-GW03

Coll. Ending Date/Time: 12/5/00 09:00

Matrix: Aqueous

Collection Begin Date/Time: 12/05/00 08:55

Sample Description:

**Analysis Type: NVOA GCMS AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u>                | <u>Result</u> | <u>Units</u> | <u>Codes</u> | <u>Remark</u> |
|-------------------|------------------------------------|---------------|--------------|--------------|---------------|
| 000088062         | 2,4,6-TRICHLOROPHENOL              | 4.0           | ug/L         | U            |               |
| 000095954         | 2,4,5-TRICHLOROPHENOL              | 4.0           | ug/L         | U            |               |
| 000091587         | 2-CHLORONAPHTHALENE                | 4.0           | ug/L         | U            |               |
| 000088744         | 2-NITROANILINE                     | 4.0           | ug/L         | U            |               |
| 000131113         | DIMETHYL PHTHALATE                 | 4.0           | ug/L         | U            |               |
| 000208968         | ACENAPHTHYLENE                     | 4.0           | ug/L         | U            |               |
| 000606202         | 2,6-DINITROTOLUENE                 | 4.0           | ug/L         | U            |               |
| 000099092         | 3-NITROANILINE                     | 48            | ug/L         | U            |               |
| 000083329         | ACENAPHTHENE                       | 4.0           | ug/L         | U            |               |
| 000051285         | 2,4-DINITROPHENOL                  | 10            | ug/L         | U            |               |
| 000100027         | 4-NITROPHENOL                      | 4.0           | ug/L         | U            |               |
| 000132649         | DIBENZOFURAN                       | 4.0           | ug/L         | U            |               |
| 000121142         | 2,4-DINITROTOLUENE                 | 4.0           | ug/L         | U            |               |
| 000086737         | FLUORENE                           | 4.0           | ug/L         | U            |               |
| 000084662         | DIETHYLPHTHALATE                   | 4.0           | ug/L         | U            |               |
| 007005723         | 4-CHLOROPHENYL PHENYL ETHER        | 4.0           | ug/L         | U            |               |
| 000100016         | 4-NITROANILINE                     | 4.0           | ug/L         | U            |               |
| 000534521         | 4,6-DINITRO-2-METHYLPHENOL         | 4.0           | ug/L         | U            |               |
| 000086306         | N-NITROSODIPHENYLAMINE             | 4.0           | ug/L         | U            |               |
| 000103333         | DIAZENE, DIPHENYL                  | 4.0           | ug/L         | U            |               |
| 000101553         | 4-BROMOPHENYL PHENYL ETHER         | 4.0           | ug/L         | U            |               |
| 000118741         | HEXACHLOROBENZENE                  | 4.0           | ug/L         | U            |               |
| 000087865         | PENTACHLOROPHENOL                  | 10            | ug/L         | U            |               |
| 000085018         | PHENANTHRENE                       | 4.0           | ug/L         | U            |               |
| 000120127         | ANTHRACENE                         | 4.0           | ug/L         | U            |               |
| 000084742         | DI-N-BUTYL PHTHALATE               | 4.0           | ug/L         | U            |               |
| 000206440         | FLUORANTHENE                       | 4.0           | ug/L         | U            |               |
| 000129000         | PYRENE                             | 4.0           | ug/L         | U            |               |
| 000085687         | BUTYL BENZYL PHTHALATE             | 4.0           | ug/L         | U            |               |
| 000056553         | BENZO(A)ANTHRACENE                 | 4.0           | ug/L         | U            |               |
| 000218019         | CHRYSENE                           | 4.0           | ug/L         | U            |               |
| 000117817         | BIS(2-ETHYLHEXYL)PHTHALATE         | 4.0           | ug/L         | U            |               |
| 000117840         | DI-N-OCTYL PHTHALATE               | 4.0           | ug/L         | U            |               |
| 000205992         | BENZO(B)FLUORANTHENE               | 4.0           | ug/L         | U            |               |
| 000207089         | BENZO(K)FLUORANTHENE               | 4.0           | ug/L         | U            |               |
| 000050328         | BENZO(A)PYRENE                     | 4.0           | ug/L         | U            |               |
| 000193395         | INDENO(1,2,3-CD)PYRENE             | 4.0           | ug/L         | U            |               |
| 000053703         | DIBENZO(A,H)ANTHRACENE             | 4.0           | ug/L         | U            |               |
| 000191242         | BENZO(G,H,I)PERYLENE               | 4.0           | ug/L         | U            |               |
|                   | UNKNOWN SUBSTITUTED ACID ,RT =14.5 | 7.0           | ug/L         | J Q1         |               |
|                   | UNKNOWN ,RT =15.92                 | 2.0           | ug/L         | J QT         |               |



**Survey Name: Lackawanna Foundry**

Project Number: 00120005

\*Sorted By Sample

**AB03260** Field/Station ID: I002-GW03  
Matrix: Aqueous  
Sample Description:

Coll. Ending Date/Time: 12/5/00 09:00  
Collection Begin Date/Time: 12/05/00 08:55

**Analysis Type: NVOA GCMS AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u>                      | <u>Result</u> | <u>Units</u> | <u>Codes</u> | <u>Remark</u> |
|-------------------|--|---------------|--------------|--------------|---------------|
|                   | UNKNOWN SUBSTITUTED BENZENE ,RT =17.49   | 3.0           | ug/L         | J QT         |               |
|                   | 2,5-CYCLOHEXADIENE-1,4-DIONE, 2, ,RT =18 | 3.0           | ug/L         | J QT         |               |
|                   | UNKNOWN ,RT =18.78                       | 3.0           | ug/L         | J QT         |               |

**Analysis Type: PESTICIDES TCL GC AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Codes</u> | <u>Remark</u> |
|-------------------|---------------------|---------------|--------------|--------------|---------------|
| 319846            | ALPHA-BHC           | 0.010         | ug/L         | U            |               |
| 58899             | GAMMA-BHC           | 0.010         | ug/L         | U            |               |
| 319857            | BETA-BHC            | 0.010         | ug/L         | U            |               |
| 319868            | DELTA-BHC           | 0.010         | ug/L         | U            |               |
| 76448             | HEPTACHLOR          | 0.010         | ug/L         | U            |               |
| 309002            | ALDRIN              | 0.010         | ug/L         | U            |               |
| 1024573           | HEPTACHLOR EPOXIDE  | 0.010         | ug/L         | U            |               |
| 5566347           | GAMMA-CHLORDANE     | 0.010         | ug/L         | U            |               |
| 5103719           | ALPHA-CHLORDANE     | 0.010         | ug/L         | U            |               |
| 115297            | ENDOSULFAN I        | 0.010         | ug/L         | U            |               |
| 72559             | 4,4'-DDE            | 0.010         | ug/L         | U            |               |
| 60571             | DIELDRIN            | 0.010         | ug/L         | U            |               |
| 72208             | ENDRIN              | 0.010         | ug/L         | U            |               |
| 72548             | 4,4'-DDD            | 0.010         | ug/L         | U            |               |
| 33213659          | ENDOSULFAN II       | 0.010         | ug/L         | U            |               |
| 50293             | 4,4'-DDT            | 0.010         | ug/L         | U            |               |
| 7421934           | ENDRIN ALDEHYDE     | 0.010         | ug/L         | U            |               |
| 72435             | METHOXYCHLOR        | 0.050         | ug/L         | U            |               |
| 1031078           | ENDOSULFAN SULFATE  | 0.010         | ug/L         | U            |               |
| 53494705          | ENDRIN KETONE       | 0.010         | ug/L         | U            |               |
| 57749             | CHLORDANE           | 0.10          | ug/L         | U            |               |
| 8001352           | TOXAPHENE           | 0.10          | ug/L         | U            |               |

**Analysis Type: PCBS TCL GC AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Codes</u> | <u>Remark</u> |
|-------------------|---------------------|---------------|--------------|--------------|---------------|
| 12674-11-2        | AROCLOR 1016        | 0.10          | ug/L         | U            |               |
| 011104282         | AROCLOR 1221        | 0.20          | ug/L         | U            |               |
| 011141165         | AROCLOR 1232        | 0.10          | ug/L         | U            |               |
| 53469-21-9        | AROCLOR 1242        | 0.10          | ug/L         | U            |               |
| 012672296         | AROCLOR 1248        | 0.10          | ug/L         | U            |               |
| 11097-69-1        | AROCLOR 1254        | 0.10          | ug/L         | U            |               |
| 11096-82-5        | AROCLOR 1260        | 0.10          | ug/L         | U            |               |

**Analysis Type: METALS TAL ICP AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Codes</u> | <u>Remark</u> |
|-------------------|---------------------|---------------|--------------|--------------|---------------|
| 007440224         | SILVER              | 10            | ug/L         | U            |               |
| 007429905         | ALUMINUM            | 200           | ug/L         | U            |               |



U.S. EPA Region 2 Laboratory  
Data Report

Survey Name: Lackawanna Foundry

Project Number: 00120005

\*Sorted By Sample ID

**AB03260** Field/Station ID: 1002-GW03

Coll. Ending Date/Time: 12/5/00 09:00

Matrix: Aqueous

Collection Begin Date/Time: 12/05/00 08:55

Sample Description:

**Analysis Type: METALS TAL ICP AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|---------------------|---------------|--------------|---------------------|
| 007440393         | BARIUM              | 240           | ug/L         |                     |
| 007440417         | BERYLLIUM           | 5.0           | ug/L         | U                   |
| 007440702         | CALCIUM             | 100,000       | ug/L         |                     |
| 007440439         | CADMIUM             | 5.0           | ug/L         | U                   |
| 007440484         | COBALT              | 50            | ug/L         | U                   |
| 007440473         | CHROMIUM            | 10            | ug/L         | U                   |
| 007440508         | COPPER              | 25            | ug/L         | U                   |
| 007439896         | IRON                | 6,200         | ug/L         |                     |
| 007440097         | POTASSIUM           | 5,000         | ug/L         | U                   |
| 007439954         | MAGNESIUM           | 22,000        | ug/L         |                     |
| 007439965         | MANGANESE           | 1,700         | ug/L         |                     |
| 007440235         | SODIUM              | 22,000        | ug/L         |                     |
| 007440020         | NICKEL              | 40            | ug/L         | U                   |
| 007440360         | ANTIMONY            | 60            | ug/L         | U                   |
| 007440622         | VANADIUM            | 50            | ug/L         | U                   |
| 007440666         | ZINC                | 33            | ug/L         |                     |

**Single Component Analyses**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|---------------------|---------------|--------------|---------------------|
| 7440-38-2         | ARSENIC             | 10            | ug/L         | U                   |
| 7439-92-1         | LEAD                | 3.0           | ug/L         | U                   |
| 7782-49-2         | SELENIUM            | 5.0           | ug/L         | U                   |
| 7440-28-0         | THALLIUM            | 10            | ug/L         | U                   |

**Single Component Analyses**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|---------------------|---------------|--------------|---------------------|
| 007439976         | MERCURY (ESAT)      | 0.20          | ug/L         | U                   |

**AB03261** Field/Station ID: 1002-GW04

Coll. Ending Date/Time: 12/5/00 11:05

Matrix: Aqueous

Collection Begin Date/Time: 12/05/00 11:00

Sample Description:

**Analysis Type: VOA TCL GCMS AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u>      | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|--------------------------|---------------|--------------|---------------------|
| 000074873         | CHLOROMETHANE            | 10            | ug/L         | U                   |
| 000075014         | VINYL CHLORIDE           | 10            | ug/L         | U                   |
| 000074839         | BROMOMETHANE             | 10            | ug/L         | U                   |
| 000075003         | CHLOROETHANE             | 10            | ug/L         | U                   |
| 000075354         | 1,1-DICHLOROETHENE       | 10            | ug/L         | U                   |
| 000075150         | CARBON DISULFIDE         | 10            | ug/L         | U                   |
| 000067641         | ACETONE                  | 9.0           | ug/L         | JQM                 |
| 000075092         | METHYLENE CHLORIDE       | 10            | ug/L         | U                   |
| 000156605         | TRANS-1,2-DICHLOROETHENE | 10            | ug/L         | U                   |





**Survey Name: Lackawanna Foundry**

Project Number: 00120005

\*Sorted By Sample

**AB03261** Field/Station ID: 1002-GW04  
Matrix: Aqueous  
Sample Description:

Coll. Ending Date/Time: 12/5/00 11:05  
Collection Begin Date/Time: 12/05/00 11:00

**Analysis Type: VOA TCL GCMS AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u>       | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|---------------------------|---------------|--------------|---------------------|
| 000156592         | CIS-1,2-DICHLOROETHENE    | 10            | ug/L         | U                   |
| 000075343         | 1,1-DICHLOROETHANE        | 10            | ug/L         | U                   |
| 000078933         | 2-BUTANONE                | 10            | ug/L         | U                   |
| 000067663         | CHLOROFORM                | 10            | ug/L         | U                   |
| 000107062         | 1,2-DICHLOROETHANE        | 10            | ug/L         | U                   |
| 000071556         | 1,1,1-TRICHLOROETHANE     | 10            | ug/L         | U                   |
| 000056235         | CARBON TETRACHLORIDE      | 10            | ug/L         | U                   |
| 000071432         | BENZENE                   | 10            | ug/L         | U                   |
| 025323891         | TRICHLOROETHENE           | 10            | ug/L         | U                   |
| 000078875         | 1,2-DICHLOROPROPANE       | 10            | ug/L         | U                   |
| 000074975         | BROMODICHLOROMETHANE      | 10            | ug/L         | U                   |
| 010061015         | CIS-1,3-DICHLOROPROPENE   | 10            | ug/L         | U                   |
| 010061026         | TRANS-1,3-DICHLOROPROPENE | 10            | ug/L         | U                   |
| 000079005         | 1,1,2-TRICHLOROETHANE     | 10            | ug/L         | U                   |
| 000124481         | DIBROMOCHLOROMETHANE      | 10            | ug/L         | U                   |
| 000075252         | BROMOFORM                 | 10            | ug/L         | U                   |
| 000108101         | 4-METHYL-2-PENTANONE      | 10            | ug/L         | U                   |
| 000108883         | TOLUENE                   | 10            | ug/L         | U                   |
| 000127184         | TETRACHLOROETHENE         | 10            | ug/L         | U                   |
| 000591786         | 2-HEXANONE                | 10            | ug/L         | U                   |
| 000108907         | CHLOROBENZENE             | 10            | ug/L         | U                   |
| 000100414         | ETHYLBENZENE              | 10            | ug/L         | U                   |
| 001330207         | M,P-XYLENE                | 10            | ug/L         | U                   |
| 000095476         | O-XYLENE                  | 10            | ug/L         | U                   |
| 000100425         | STYRENE                   | 10            | ug/L         | U                   |
| 000079345         | 1,1,2,2-TETRACHLOROETHANE | 10            | ug/L         | U                   |

**Analysis Type: NVOA GCMS AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u>         | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|-----------------------------|---------------|--------------|---------------------|
| 000108952         | PHENOL                      | 4.0           | ug/L         | U                   |
| 000111444         | BIS(2-CHLOROETHYL)ETHER     | 4.0           | ug/L         | U                   |
| 000095578         | 2-CHLOROPHENOL              | 4.0           | ug/L         | U                   |
| 000541731         | 1,3-DICHLOROBENZENE         | 4.0           | ug/L         | U                   |
| 000106467         | 1,4-DICHLOROBENZENE         | 4.0           | ug/L         | U                   |
| 000095501         | 1,2-DICHLOROBENZENE         | 4.0           | ug/L         | U                   |
| 000100516         | BENZYL ALCOHOL              | 10            | ug/L         | U                   |
| 000095487         | 2-METHYLPHENOL              | 4.0           | ug/L         | U                   |
| 000108601         | BIS(2-CHLOROISOPROPYL)ETHER | 4.0           | ug/L         | U                   |
| 000106445         | 4-METHYLPHENOL              | 4.0           | ug/L         | U                   |
| 000621647         | N-NITROSO-DI-N-PROPYLAMINE  | 4.0           | ug/L         | U                   |
| 000067721         | HEXACHLOROETHANE            | 4.0           | ug/L         | U                   |
| 000098953         | NITROBENZENE                | 4.0           | ug/L         | U                   |



U.S. EPA Region 2 Laboratory  
Data Report

Survey Name: Lackawanna Foundry

Project Number: 00120005

\*Sorted By Sample ID

**AB03261**

Field/Station ID: 1002-GW04

Coll. Ending Date/Time: 12/5/00 11:05

Matrix: Aqueous

Collection Begin Date/Time: 12/05/00 11:00

Sample Description:

**Analysis Type: NVOA GCMS AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u>         | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|-----------------------------|---------------|--------------|---------------------|
| 000078591         | ISOPHORONE                  | 4.0           | ug/L         | U                   |
| 000088755         | 2-NITROPHENOL               | 4.0           | ug/L         | U                   |
| 000105679         | 2,4-DIMETHYLPHENOL          | 4.0           | ug/L         | U                   |
| 000111911         | BIS(2-CHLOROETHOXY)METHANE  | 4.0           | ug/L         | U                   |
| 000120832         | 2,4-DICHLOROPHENOL          | 4.0           | ug/L         | U                   |
| 000120821         | 1,2,4-TRICHLOROBENZENE      | 4.0           | ug/L         | U                   |
| 000065850         | BENZOIC ACID                | 50            | ug/L         | U                   |
| 000091203         | NAPHTHALENE                 | 4.0           | ug/L         | U                   |
| 000106478         | 4-CHLOROANILINE             | 50            | ug/L         | U                   |
| 000087683         | HEXACHLOROBUTADIENE         | 4.0           | ug/L         | U                   |
| 000059507         | 4-CHLORO-3-METHYLPHENOL     | 4.0           | ug/L         | U                   |
| 000091576         | 2-METHYL NAPHTHALENE        | 4.0           | ug/L         | U                   |
| 000077474         | HEXACHLOROCYCLOPENTADIENE   | 4.0           | ug/L         | U                   |
| 000088062         | 2,4,6-TRICHLOROPHENOL       | 4.0           | ug/L         | U                   |
| 000095954         | 2,4,5-TRICHLOROPHENOL       | 4.0           | ug/L         | U                   |
| 000091587         | 2-CHLORONAPHTHALENE         | 4.0           | ug/L         | U                   |
| 000088744         | 2-NITROANILINE              | 4.0           | ug/L         | U                   |
| 000131113         | DIMETHYL PHTHALATE          | 4.0           | ug/L         | U                   |
| 000208968         | ACENAPHTHYLENE              | 4.0           | ug/L         | U                   |
| 000606202         | 2,6-DINITROTOLUENE          | 4.0           | ug/L         | U                   |
| 000099092         | 3-NITROANILINE              | 50            | ug/L         | U                   |
| 000083329         | ACENAPHTHENE                | 4.0           | ug/L         | U                   |
| 000051285         | 2,4-DINITROPHENOL           | 10            | ug/L         | U                   |
| 000100027         | 4-NITROPHENOL               | 4.0           | ug/L         | U                   |
| 000132649         | DIBENZOFURAN                | 4.0           | ug/L         | U                   |
| 000121142         | 2,4-DINITROTOLUENE          | 4.0           | ug/L         | U                   |
| 000086737         | FLUORENE                    | 4.0           | ug/L         | U                   |
| 000084662         | DIETHYLPHTHALATE            | 4.0           | ug/L         | U                   |
| 007005723         | 4-CHLOROPHENYL PHENYL ETHER | 4.0           | ug/L         | U                   |
| 000100016         | 4-NITROANILINE              | 4.0           | ug/L         | U                   |
| 000534521         | 4,6-DINITRO-2-METHYLPHENOL  | 4.0           | ug/L         | U                   |
| 000086306         | N-NITROSODIPHENYLAMINE      | 4.0           | ug/L         | U                   |
| 000103333         | DIAZENE, DIPHENYL           | 4.0           | ug/L         | U                   |
| 000101553         | 4-BROMOPHENYL PHENYL ETHER  | 4.0           | ug/L         | U                   |
| 000118741         | HEXACHLOROBENZENE           | 4.0           | ug/L         | U                   |
| 000087865         | PENTACHLOROPHENOL           | 10            | ug/L         | U                   |
| 000085018         | PHENANTHRENE                | 4.0           | ug/L         | U                   |
| 000120127         | ANTHRACENE                  | 4.0           | ug/L         | U                   |
| 000084742         | DI-N-BUTYL PHTHALATE        | 4.0           | ug/L         | U                   |
| 000206440         | FLUORANTHENE                | 4.0           | ug/L         | U                   |
| 000129000         | PYRENE                      | 4.0           | ug/L         | U                   |



**Survey Name: Lackawanna Foundry**

Project Number: 00120005

\*Sorted By Sample

**AB03261**

Field/Station ID: 1002-GW04

Coll. Ending Date/Time: 12/5/00 11:05

Matrix: Aqueous

Collection Begin Date/Time: 12/05/00 11:00

Sample Description:

**Analysis Type: NVOA GCMS AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u>        | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|----------------------------|---------------|--------------|---------------------|
| 000085687         | BUTYL BENZYL PHTHALATE     | 4.0           | ug/L         | U                   |
| 000056553         | BENZO(A)ANTHRACENE         | 4.0           | ug/L         | U                   |
| 000218019         | CHRYSENE                   | 4.0           | ug/L         | U                   |
| 000117817         | BIS(2-ETHYLHEXYL)PHTHALATE | 4.0           | ug/L         | U                   |
| 000117840         | DI-N-OCTYL PHTHALATE       | 4.0           | ug/L         | U                   |
| 000205992         | BENZO(B)FLUORANTHENE       | 4.0           | ug/L         | U                   |
| 000207089         | BENZO(K)FLUORANTHENE       | 4.0           | ug/L         | U                   |
| 000050328         | BENZO(A)PYRENE             | 4.0           | ug/L         | U                   |
| 000193395         | INDENO(1,2,3-CD)PYRENE     | 4.0           | ug/L         | U                   |
| 000053703         | DIBENZO(A,H)ANTHRACENE     | 4.0           | ug/L         | U                   |
| 000191242         | BENZO(G,H,I)PERYLENE       | 4.0           | ug/L         | U                   |
|                   | UNKNOWN ,RT =13.55         | 2.0           | ug/L         | J QT                |

**Analysis Type: PESTICIDES TCL GC AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|---------------------|---------------|--------------|---------------------|
| 319846            | ALPHA-BHC           | 0.010         | ug/L         | U                   |
| 58899             | GAMMA-BHC           | 0.010         | ug/L         | U                   |
| 319857            | BETA-BHC            | 0.010         | ug/L         | U                   |
| 319868            | DELTA-BHC           | 0.010         | ug/L         | U                   |
| 76448             | HEPTACHLOR          | 0.010         | ug/L         | U                   |
| 309002            | ALDRIN              | 0.010         | ug/L         | U                   |
| 1024573           | HEPTACHLOR EPOXIDE  | 0.010         | ug/L         | U                   |
| 5566347           | GAMMA-CHLORDANE     | 0.010         | ug/L         | U                   |
| 5103719           | ALPHA-CHLORDANE     | 0.010         | ug/L         | U                   |
| 115297            | ENDOSULFAN I        | 0.010         | ug/L         | U                   |
| 72559             | 4,4'-DDE            | 0.010         | ug/L         | U                   |
| 60571             | DIELDRIN            | 0.010         | ug/L         | U                   |
| 72208             | ENDRIN              | 0.010         | ug/L         | U                   |
| 72548             | 4,4'-DDD            | 0.010         | ug/L         | U                   |
| 33213659          | ENDOSULFAN II       | 0.010         | ug/L         | U                   |
| 50293             | 4,4'-DDT            | 0.010         | ug/L         | U                   |
| 7421934           | ENDRIN ALDEHYDE     | 0.010         | ug/L         | U                   |
| 72435             | METHOXYCHLOR        | 0.050         | ug/L         | U                   |
| 1031078           | ENDOSULFAN SULFATE  | 0.010         | ug/L         | U                   |
| 53494705          | ENDRIN KETONE       | 0.010         | ug/L         | U                   |
| 57749             | CHLORDANE           | 0.10          | ug/L         | U                   |
| 8001352           | TOXAPHENE           | 0.10          | ug/L         | U                   |

**Analysis Type: PCBS TCL GC AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|---------------------|---------------|--------------|---------------------|
| 12674-11-2        | AROCLOR 1016        | 0.10          | ug/L         | U                   |
| 011104282         | AROCLOR 1221        | 0.20          | ug/L         | U                   |
| 011141165         | AROCLOR 1232        | 0.10          | ug/L         | U                   |



**Survey Name: Lackawanna Foundry**

Project Number: 00120005

\*Sorted By Sample ID

**AB03261**

Field/Station ID: 1002-GW04

Coll. Ending Date/Time: 12/5/00 11:05

Matrix: Aqueous

Collection Begin Date/Time: 12/05/00 11:00

Sample Description:

**Analysis Type: PCBS TCL GC AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Codes</u> | <u>Remark</u> |
|-------------------|---------------------|---------------|--------------|--------------|---------------|
| 53469-21-9        | AROCLOR 1242        | 0.10          | ug/L         | U            |               |
| 012672296         | AROCLOR 1248        | 0.10          | ug/L         | U            |               |
| 11097-69-1        | AROCLOR 1254        | 0.10          | ug/L         | U            |               |
| 11096-82-5        | AROCLOR 1260        | 0.10          | ug/L         | U            |               |

**Analysis Type: METALS TAL ICP AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Codes</u> | <u>Remark</u> |
|-------------------|---------------------|---------------|--------------|--------------|---------------|
| 007440224         | SILVER              | 10            | ug/L         | U            |               |
| 007429905         | ALUMINUM            | 3,500         | ug/L         |              |               |
| 007440393         | BARIUM              | 220           | ug/L         |              |               |
| 007440417         | BERYLLIUM           | 5.0           | ug/L         | U            |               |
| 007440702         | CALCIUM             | 60,000        | ug/L         |              |               |
| 007440439         | CADMIUM             | 5.0           | ug/L         | U            |               |
| 007440484         | COBALT              | 50            | ug/L         | U            |               |
| 007440473         | CHROMIUM            | 10            | ug/L         | U            |               |
| 007440508         | COPPER              | 25            | ug/L         | U            |               |
| 007439896         | IRON                | 42,000        | ug/L         |              |               |
| 007440097         | POTASSIUM           | 5,000         | ug/L         | U            |               |
| 007439954         | MAGNESIUM           | 14,000        | ug/L         |              |               |
| 007439965         | MANGANESE           | 1,700         | ug/L         |              |               |
| 007440235         | SODIUM              | 25,000        | ug/L         |              |               |
| 007440020         | NICKEL              | 40            | ug/L         | U            |               |
| 007440360         | ANTIMONY            | 60            | ug/L         | U            |               |
| 007440622         | VANADIUM            | 50            | ug/L         | U            |               |
| 007440666         | ZINC                | 76            | ug/L         |              |               |

**Single Component Analyses**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Codes</u> | <u>Remark</u> |
|-------------------|---------------------|---------------|--------------|--------------|---------------|
| 7440-38-2         | ARSENIC             | 13            | ug/L         |              |               |
| 7439-92-1         | LEAD                | 3.0           | ug/L         | U            |               |
| 7782-49-2         | SELENIUM            | 5.0           | ug/L         | U            |               |
| 7440-28-0         | THALLIUM            | 10            | ug/L         | U            |               |

**Single Component Analyses**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Codes</u> | <u>Remark</u> |
|-------------------|---------------------|---------------|--------------|--------------|---------------|
| 007439976         | MERCURY (ESAT)      | 0.20          | ug/L         | U            |               |

**AB03262**

Field/Station ID: 1002-GW05

Coll. Ending Date/Time: 12/5/00 10:00

Matrix: Aqueous

Collection Begin Date/Time: 12/05/00 09:55

Sample Description:

**Analysis Type: VOA TCL GCMS AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Codes</u> | <u>Remark</u> |
|-------------------|---------------------|---------------|--------------|--------------|---------------|
| 000074873         | CHLOROMETHANE       | 10            | ug/L         | U            |               |



**Survey Name: Lackawanna Foundry**

Project Number: 00120005

\*Sorted By Sample

**AB03262**

Field/Station ID: 1002-GW05

Coll. Ending Date/Time: 12/5/00 10:00

Matrix: Aqueous

Collection Begin Date/Time: 12/05/00 09:55

Sample Description:

**Analysis Type: VOA TCL GCMS AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u>       | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|---------------------------|---------------|--------------|---------------------|
| 000075014         | VINYL CHLORIDE            | 10            | ug/L         | U                   |
| 000074839         | BROMOMETHANE              | 10            | ug/L         | U                   |
| 000075003         | CHLOROETHANE              | 10            | ug/L         | U                   |
| 000075354         | 1,1-DICHLOROETHENE        | 10            | ug/L         | U                   |
| 000075150         | CARBON DISULFIDE          | 10            | ug/L         | U                   |
| 000067641         | ACETONE                   | 6.0           | ug/L         | J QM                |
| 000075092         | METHYLENE CHLORIDE        | 10            | ug/L         | U                   |
| 000156605         | TRANS-1,2-DICHLOROETHENE  | 10            | ug/L         | U                   |
| 000156592         | CIS-1,2-DICHLOROETHENE    | 10            | ug/L         | U                   |
| 000075343         | 1,1-DICHLOROETHANE        | 10            | ug/L         | U                   |
| 000078933         | 2-BUTANONE                | 10            | ug/L         | U                   |
| 000067663         | CHLOROFORM                | 10            | ug/L         | U                   |
| 000107062         | 1,2-DICHLOROETHANE        | 10            | ug/L         | U                   |
| 000071556         | 1,1,1-TRICHLOROETHANE     | 10            | ug/L         | U                   |
| 000056235         | CARBON TETRACHLORIDE      | 10            | ug/L         | U                   |
| 000071432         | BENZENE                   | 10            | ug/L         | U                   |
| 025323891         | TRICHLOROETHENE           | 10            | ug/L         | U                   |
| 000078875         | 1,2-DICHLOROPROPANE       | 10            | ug/L         | U                   |
| 000074975         | BROMODICHLOROMETHANE      | 10            | ug/L         | U                   |
| 010061015         | CIS-1,3-DICHLOROPROPENE   | 10            | ug/L         | U                   |
| 010061026         | TRANS-1,3-DICHLOROPROPENE | 10            | ug/L         | U                   |
| 000079005         | 1,1,2-TRICHLOROETHANE     | 10            | ug/L         | U                   |
| 000124481         | DIBROMOCHLOROMETHANE      | 10            | ug/L         | U                   |
| 000075252         | BROMOFORM                 | 10            | ug/L         | U                   |
| 000108101         | 4-METHYL-2-PENTANONE      | 10            | ug/L         | U                   |
| 000108883         | TOLUENE                   | 10            | ug/L         | U                   |
| 000127184         | TETRACHLOROETHENE         | 10            | ug/L         | U                   |
| 000591786         | 2-HEXANONE                | 10            | ug/L         | U                   |
| 000108907         | CHLOROBENZENE             | 10            | ug/L         | U                   |
| 000100414         | ETHYLBENZENE              | 10            | ug/L         | U                   |
| 001330207         | M,P-XYLENE                | 10            | ug/L         | U                   |
| 000095476         | O-XYLENE                  | 10            | ug/L         | U                   |
| 000100425         | STYRENE                   | 10            | ug/L         | U                   |
| 000079345         | 1,1,2,2-TETRACHLOROETHANE | 10            | ug/L         | U                   |

**Analysis Type: NVOA GCMS AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u>     | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|-------------------------|---------------|--------------|---------------------|
| 000108952         | PHENOL                  | 4.0           | ug/L         | U                   |
| 000111444         | BIS(2-CHLOROETHYL)ETHER | 4.0           | ug/L         | U                   |
| 000095578         | 2-CHLOROPHENOL          | 4.0           | ug/L         | U                   |
| 000541731         | 1,3-DICHLOROBENZENE     | 4.0           | ug/L         | U                   |
| 000106467         | 1,4-DICHLOROBENZENE     | 4.0           | ug/L         | U                   |



U.S. EPA Region 2 Laboratory  
Data Report

Survey Name: Lackawanna Foundry

Project Number: 00120005

\*Sorted By Sample ID

AB03262

Field/Station ID: 1002-GW05

Coll. Ending Date/Time: 12/5/00 10:00

Matrix: Aqueous

Collection Begin Date/Time: 12/05/00 09:55

Sample Description:

Analysis Type: NVOA GCMS AQUEOUS

| CAS Number | Analyte Name                | Result | Units | Remarks Codes |
|------------|-----------------------------|--------|-------|---------------|
| 000095501  | 1,2-DICHLOROBENZENE         | 4.0    | ug/L  | U             |
| 000100516  | BENZYL ALCOHOL              | 10     | ug/L  | U             |
| 000095487  | 2-METHYLPHENOL              | 4.0    | ug/L  | U             |
| 000108601  | BIS(2-CHLOROISOPROPYL)ETHER | 4.0    | ug/L  | U             |
| 000106445  | 4-METHYLPHENOL              | 4.0    | ug/L  | U             |
| 000621647  | N-NITROSO-DI-N-PROPYLAMINE  | 4.0    | ug/L  | U             |
| 000067721  | HEXACHLOROETHANE            | 4.0    | ug/L  | U             |
| 000098953  | NITROBENZENE                | 4.0    | ug/L  | U             |
| 000078591  | ISOPHORONE                  | 4.0    | ug/L  | U             |
| 000088755  | 2-NITROPHENOL               | 4.0    | ug/L  | U             |
| 000105679  | 2,4-DIMETHYLPHENOL          | 4.0    | ug/L  | U             |
| 000111911  | BIS(2-CHLOROETHOXY)METHANE  | 4.0    | ug/L  | U             |
| 000120832  | 2,4-DICHLOROPHENOL          | 4.0    | ug/L  | U             |
| 000120821  | 1,2,4-TRICHLOROBENZENE      | 4.0    | ug/L  | U             |
| 000065850  | BENZOIC ACID                | 48     | ug/L  | U             |
| 000091203  | NAPHTHALENE                 | 4.0    | ug/L  | U             |
| 000106478  | 4-CHLOROANILINE             | 48     | ug/L  | U             |
| 000087683  | HEXACHLOROBUTADIENE         | 4.0    | ug/L  | U             |
| 000059507  | 4-CHLORO-3-METHYLPHENOL     | 4.0    | ug/L  | U             |
| 000091576  | 2-METHYL NAPHTHALENE        | 4.0    | ug/L  | U             |
| 000077474  | HEXACHLOROCYCLOPENTADIENE   | 4.0    | ug/L  | U             |
| 000088062  | 2,4,6-TRICHLOROPHENOL       | 4.0    | ug/L  | U             |
| 000095954  | 2,4,5-TRICHLOROPHENOL       | 4.0    | ug/L  | U             |
| 000091587  | 2-CHLORONAPHTHALENE         | 4.0    | ug/L  | U             |
| 000088744  | 2-NITROANILINE              | 4.0    | ug/L  | U             |
| 000131113  | DIMETHYL PHTHALATE          | 4.0    | ug/L  | U             |
| 000208968  | ACENAPHTHYLENE              | 4.0    | ug/L  | U             |
| 000606202  | 2,6-DINITROTOLUENE          | 4.0    | ug/L  | U             |
| 000099092  | 3-NITROANILINE              | 48     | ug/L  | U             |
| 000083329  | ACENAPHTHENE                | 4.0    | ug/L  | U             |
| 000051285  | 2,4-DINITROPHENOL           | 10     | ug/L  | U             |
| 000100027  | 4-NITROPHENOL               | 4.0    | ug/L  | U             |
| 000132649  | DIBENZOFURAN                | 4.0    | ug/L  | U             |
| 000121142  | 2,4-DINITROTOLUENE          | 4.0    | ug/L  | U             |
| 000086737  | FLUORENE                    | 4.0    | ug/L  | U             |
| 000084662  | DIETHYLPHTHALATE            | 4.0    | ug/L  | U             |
| 007005723  | 4-CHLOROPHENYL PHENYL ETHER | 4.0    | ug/L  | U             |
| 000100016  | 4-NITROANILINE              | 4.0    | ug/L  | U             |
| 000534521  | 4,6-DINITRO-2-METHYLPHENOL  | 4.0    | ug/L  | U             |
| 000086306  | N-NITROSODIPHENYLAMINE      | 4.0    | ug/L  | U             |
| 000103333  | DIAZENE, DIPHENYL           | 4.0    | ug/L  | U             |



**Survey Name: Lackawanna Foundry**

Project Number: 00120005

\*Sorted By Sample

**AB03262** Field/Station ID: 1002-GW05

Coll. Ending Date/Time: 12/5/00 10:00

Matrix: Aqueous

Collection Begin Date/Time: 12/05/00 09:55

Sample Description:

**Analysis Type: NVOA GCMS AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u>        | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|----------------------------|---------------|--------------|---------------------|
| 000101553         | 4-BROMOPHENYL PHENYL ETHER | 4.0           | ug/L         | U                   |
| 000118741         | HEXACHLOROBENZENE          | 4.0           | ug/L         | U                   |
| 000087865         | PENTACHLOROPHENOL          | 10            | ug/L         | U                   |
| 000085018         | PHENANTHRENE               | 4.0           | ug/L         | U                   |
| 000120127         | ANTHRACENE                 | 4.0           | ug/L         | U                   |
| 000084742         | DI-N-BUTYL PHTHALATE       | 4.0           | ug/L         | U                   |
| 000206440         | FLUORANTHENE               | 4.0           | ug/L         | U                   |
| 000129000         | PYRENE                     | 4.0           | ug/L         | U                   |
| 000085687         | BUTYL BENZYL PHTHALATE     | 4.0           | ug/L         | U                   |
| 000056553         | BENZO(A)ANTHRACENE         | 4.0           | ug/L         | U                   |
| 000218019         | CHRYSENE                   | 4.0           | ug/L         | U                   |
| 000117817         | BIS(2-ETHYLHEXYL)PHTHALATE | 4.0           | ug/L         | U                   |
| 000117840         | DI-N-OCTYL PHTHALATE       | 4.0           | ug/L         | U                   |
| 000205992         | BENZO(B)FLUORANTHENE       | 4.0           | ug/L         | U                   |
| 000207089         | BENZO(K)FLUORANTHENE       | 4.0           | ug/L         | U                   |
| 000050328         | BENZO(A)PYRENE             | 4.0           | ug/L         | U                   |
| 000193395         | INDENO(1,2,3-CD)PYRENE     | 4.0           | ug/L         | U                   |
| 000053703         | DIBENZO(A,H)ANTHRACENE     | 4.0           | ug/L         | U                   |
| 000191242         | BENZO(G,H,I)PERYLENE       | 4.0           | ug/L         | U                   |

**Analysis Type: PESTICIDES TCL GC AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|---------------------|---------------|--------------|---------------------|
| 319846            | ALPHA-BHC           | 0.010         | ug/L         | U                   |
| 58899             | GAMMA-BHC           | 0.010         | ug/L         | U                   |
| 319857            | BETA-BHC            | 0.010         | ug/L         | U                   |
| 319868            | DELTA-BHC           | 0.010         | ug/L         | U                   |
| 76448             | HEPTACHLOR          | 0.010         | ug/L         | U                   |
| 309002            | ALDRIN              | 0.010         | ug/L         | U                   |
| 1024573           | HEPTACHLOR EPOXIDE  | 0.010         | ug/L         | U                   |
| 5566347           | GAMMA-CHLORDANE     | 0.010         | ug/L         | U                   |
| 5103719           | ALPHA-CHLORDANE     | 0.010         | ug/L         | U                   |
| 115297            | ENDOSULFAN I        | 0.010         | ug/L         | U                   |
| 72559             | 4,4'-DDE            | 0.010         | ug/L         | U                   |
| 60571             | DIELDRIN            | 0.010         | ug/L         | U                   |
| 72208             | ENDRIN              | 0.010         | ug/L         | U                   |
| 72548             | 4,4'-DDD            | 0.010         | ug/L         | U                   |
| 33213659          | ENDOSULFAN II       | 0.010         | ug/L         | U                   |
| 50293             | 4,4'-DDT            | 0.010         | ug/L         | U                   |
| 7421934           | ENDRIN ALDEHYDE     | 0.010         | ug/L         | U                   |
| 72435             | METHOXYCHLOR        | 0.050         | ug/L         | U                   |
| 1031078           | ENDOSULFAN SULFATE  | 0.010         | ug/L         | U                   |
| 53494705          | ENDRIN KETONE       | 0.010         | ug/L         | U                   |



**Survey Name: Lackawanna Foundry**

Project Number: 00120005

\*Sorted By Sample ID

**AB03262**

Field/Station ID: 1002-GW05

Coll. Ending Date/Time: 12/5/00 10:00

Matrix: Aqueous

Collection Begin Date/Time: 12/05/00 09:55

Sample Description:

**Analysis Type: PESTICIDES TCL GC AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Codes</u> | <u>Remark</u> |
|-------------------|---------------------|---------------|--------------|--------------|---------------|
| 57749             | CHLORDANE           | 0.10          | ug/L         | U            |               |
| 8001352           | TOXAPHENE           | 0.10          | ug/L         | U            |               |

**Analysis Type: PCBS TCL GC AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Codes</u> | <u>Remark</u> |
|-------------------|---------------------|---------------|--------------|--------------|---------------|
| 12674-11-2        | AROCLOR 1016        | 0.10          | ug/L         | U            |               |
| 011104282         | AROCLOR 1221        | 0.20          | ug/L         | U            |               |
| 011141165         | AROCLOR 1232        | 0.10          | ug/L         | U            |               |
| 53469-21-9        | AROCLOR 1242        | 0.10          | ug/L         | U            |               |
| 012672296         | AROCLOR 1248        | 0.10          | ug/L         | U            |               |
| 11097-69-1        | AROCLOR 1254        | 0.10          | ug/L         | U            |               |
| 11096-82-5        | AROCLOR 1260        | 0.10          | ug/L         | U            |               |

**Analysis Type: METALS TAL ICP AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Codes</u> | <u>Remark</u> |
|-------------------|---------------------|---------------|--------------|--------------|---------------|
| 007440224         | SILVER              | 10            | ug/L         | U            |               |
| 007429905         | ALUMINUM            | 2,500         | ug/L         |              |               |
| 007440393         | BARIUM              | 300           | ug/L         |              |               |
| 007440417         | BERYLLIUM           | 5.0           | ug/L         | U            |               |
| 007440702         | CALCIUM             | 87,000        | ug/L         |              |               |
| 007440439         | CADMIUM             | 5.0           | ug/L         | U            |               |
| 007440484         | COBALT              | 50            | ug/L         | U            |               |
| 007440473         | CHROMIUM            | 23            | ug/L         |              |               |
| 007440508         | COPPER              | 25            | ug/L         | U            |               |
| 007439896         | IRON                | 34,000        | ug/L         |              |               |
| 007440097         | POTASSIUM           | 5,000         | ug/L         | U            |               |
| 007439954         | MAGNESIUM           | 22,000        | ug/L         |              |               |
| 007439965         | MANGANESE           | 2,200         | ug/L         |              |               |
| 007440235         | SODIUM              | 20,000        | ug/L         |              |               |
| 007440020         | NICKEL              | 40            | ug/L         | U            |               |
| 007440360         | ANTIMONY            | 60            | ug/L         | U            |               |
| 007440622         | VANADIUM            | 50            | ug/L         | U            |               |
| 007440666         | ZINC                | 20            | ug/L         | U            |               |

**Single Component Analyses**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Codes</u> | <u>Remark</u> |
|-------------------|---------------------|---------------|--------------|--------------|---------------|
| 7440-38-2         | ARSENIC             | 12            | ug/L         |              |               |
| 7439-92-1         | LEAD                | 3.4           | ug/L         |              |               |
| 7782-49-2         | SELENIUM            | 5.0           | ug/L         | U            |               |
| 7440-28-0         | THALLIUM            | 10            | ug/L         | U            |               |

**Single Component Analyses**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Codes</u> | <u>Remark</u> |
|-------------------|---------------------|---------------|--------------|--------------|---------------|
| 007439976         | MERCURY (ESAT)      | 0.20          | ug/L         | U            |               |





Survey Name: Lackawanna Foundry

Project Number: 00120005

\*Sorted By Sample

AB03263

Field/Station ID: 1002-GW06

Coll. Ending Date/Time: 12/5/00 14:05

Matrix: Aqueous

Collection Begin Date/Time: 12/05/00 14:00

Sample Description:

Analysis Type: VOA TCL GCMS AQUEOUS

| CAS Number | Analyte Name                   | Result | Units | Remark Codes |
|------------|--------------------------------|--------|-------|--------------|
| 000074873  | CHLOROMETHANE                  | 10     | ug/L  | U            |
| 000075014  | VINYL CHLORIDE                 | 10     | ug/L  | U            |
| 000074839  | BROMOMETHANE                   | 10     | ug/L  | U            |
| 000075003  | CHLOROETHANE                   | 10     | ug/L  | U            |
| 000075354  | 1,1-DICHLOROETHENE             | 10     | ug/L  | U            |
| 000075150  | CARBON DISULFIDE               | 10     | ug/L  | U            |
| 000067641  | ACETONE                        | 10     | ug/L  | U            |
| 000075092  | METHYLENE CHLORIDE             | 10     | ug/L  | U            |
| 000156605  | TRANS-1,2-DICHLOROETHENE       | 10     | ug/L  | U            |
| 000156592  | CIS-1,2-DICHLOROETHENE         | 10     | ug/L  | U            |
| 000075343  | 1,1-DICHLOROETHANE             | 10     | ug/L  | U            |
| 000078933  | 2-BUTANONE                     | 10     | ug/L  | U            |
| 000067663  | CHLOROFORM                     | 10     | ug/L  | U            |
| 000107062  | 1,2-DICHLOROETHANE             | 10     | ug/L  | U            |
| 000071556  | 1,1,1-TRICHLOROETHANE          | 10     | ug/L  | U            |
| 000056235  | CARBON TETRACHLORIDE           | 10     | ug/L  | U            |
| 000071432  | BENZENE                        | 10     | ug/L  | U            |
| 025323891  | TRICHLOROETHENE                | 10     | ug/L  | U            |
| 000078875  | 1,2-DICHLOROPROPANE            | 10     | ug/L  | U            |
| 000074975  | BROMODICHLOROMETHANE           | 10     | ug/L  | U            |
| 010061015  | CIS-1,3-DICHLOROPROPENE        | 10     | ug/L  | U            |
| 010061026  | TRANS-1,3-DICHLOROPROPENE      | 10     | ug/L  | U            |
| 000079005  | 1,1,2-TRICHLOROETHANE          | 10     | ug/L  | U            |
| 000124481  | DIBROMOCHLOROMETHANE           | 10     | ug/L  | U            |
| 000075252  | BROMOFORM                      | 10     | ug/L  | U            |
| 000108101  | 4-METHYL-2-PENTANONE           | 10     | ug/L  | U            |
| 000108883  | TOLUENE                        | 10     | ug/L  | U            |
| 000127184  | TETRACHLOROETHENE              | 10     | ug/L  | U            |
| 000591786  | 2-HEXANONE                     | 10     | ug/L  | U            |
| 000108907  | CHLOROBENZENE                  | 10     | ug/L  | U            |
| 000100414  | ETHYLBENZENE                   | 10     | ug/L  | U            |
| 0001330207 | M,P-XYLENE                     | 10     | ug/L  | U            |
| 000095476  | O-XYLENE                       | 10     | ug/L  | U            |
| 000100425  | STYRENE                        | 10     | ug/L  | U            |
| 000079345  | 1,1,2,2-TETRACHLOROETHANE      | 10     | ug/L  | U            |
|            | UNKNOWN HYDROCARBON ,RT =3.03  | 16     | ug/L  | J QT         |
|            | UNKNOWN HYDROCARBON ,RT =3.47  | 28     | ug/L  | J QT         |
|            | BUTANE, 2-METHYL-,RT =4.61     | 29     | ug/L  | J QT         |
|            | PENTANE ,RT =5.13              | 5.0    | ug/L  | J            |
|            | BUTANE, 2,3-DIMETHYL-,RT =6.71 | 9.0    | ug/L  | J QT         |
|            | UNKNOWN HYDROCARBON ,RT =11.57 | 5.0    | ug/L  | J QT         |



**Survey Name: Lackawanna Foundry**

Project Number: 00120005

\*Sorted By Sample ID

**AB03263**

Field/Station ID: 1002-GW06

Coll. Ending Date/Time: 12/5/00 14:05

Matrix: Aqueous

Collection Begin Date/Time: 12/05/00 14:00

Sample Description:

**Analysis Type: NVOA GCMS AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u>         | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|-----------------------------|---------------|--------------|---------------------|
| 000108952         | PHENOL                      | 4.0           | ug/L         | U                   |
| 000111444         | BIS(2-CHLOROETHYL)ETHER     | 4.0           | ug/L         | U                   |
| 000095578         | 2-CHLOROPHENOL              | 4.0           | ug/L         | U                   |
| 000541731         | 1,3-DICHLOROBENZENE         | 4.0           | ug/L         | U                   |
| 000106467         | 1,4-DICHLOROBENZENE         | 4.0           | ug/L         | U                   |
| 000095501         | 1,2-DICHLOROBENZENE         | 4.0           | ug/L         | U                   |
| 000100516         | BENZYL ALCOHOL              | 10            | ug/L         | U                   |
| 000095487         | 2-METHYLPHENOL              | 4.0           | ug/L         | U                   |
| 000108601         | BIS(2-CHLOROISOPROPYL)ETHER | 4.0           | ug/L         | U                   |
| 000106445         | 4-METHYLPHENOL              | 4.0           | ug/L         | U                   |
| 000621647         | N-NITROSO-DI-N-PROPYLAMINE  | 4.0           | ug/L         | U                   |
| 000067721         | HEXACHLOROETHANE            | 4.0           | ug/L         | U                   |
| 000098953         | NITROBENZENE                | 4.0           | ug/L         | U                   |
| 000078591         | ISOPHORONE                  | 4.0           | ug/L         | U                   |
| 000088755         | 2-NITROPHENOL               | 4.0           | ug/L         | U                   |
| 000105679         | 2,4-DIMETHYLPHENOL          | 4.0           | ug/L         | U                   |
| 000111911         | BIS(2-CHLOROETHOXY)METHANE  | 4.0           | ug/L         | U                   |
| 000120832         | 2,4-DICHLOROPHENOL          | 4.0           | ug/L         | U                   |
| 000120821         | 1,2,4-TRICHLOROBENZENE      | 4.0           | ug/L         | U                   |
| 000065850         | BENZOIC ACID                | 49            | ug/L         | U                   |
| 000091203         | NAPHTHALENE                 | 4.0           | ug/L         | U                   |
| 000106478         | 4-CHLOROANILINE             | 49            | ug/L         | U                   |
| 000087683         | HEXACHLOROBUTADIENE         | 4.0           | ug/L         | U                   |
| 000059507         | 4-CHLORO-3-METHYLPHENOL     | 4.0           | ug/L         | U                   |
| 000091576         | 2-METHYL NAPHTHALENE        | 4.0           | ug/L         | U                   |
| 000077474         | HEXACHLOROCYCLOPENTADIENE   | 4.0           | ug/L         | U                   |
| 000088062         | 2,4,6-TRICHLOROPHENOL       | 4.0           | ug/L         | U                   |
| 000095954         | 2,4,5-TRICHLOROPHENOL       | 4.0           | ug/L         | U                   |
| 000091587         | 2-CHLORONAPHTHALENE         | 4.0           | ug/L         | U                   |
| 000088744         | 2-NITROANILINE              | 4.0           | ug/L         | U                   |
| 000131113         | DIMETHYL PHTHALATE          | 4.0           | ug/L         | U                   |
| 000208968         | ACENAPHTHYLENE              | 4.0           | ug/L         | U                   |
| 000606202         | 2,6-DINITROTOLUENE          | 4.0           | ug/L         | U                   |
| 000099092         | 3-NITROANILINE              | 49            | ug/L         | U                   |
| 000083329         | ACENAPHTHENE                | 4.0           | ug/L         | U                   |
| 000051285         | 2,4-DINITROPHENOL           | 10            | ug/L         | U                   |
| 000100027         | 4-NITROPHENOL               | 4.0           | ug/L         | U                   |
| 000132649         | DIBENZOFURAN                | 4.0           | ug/L         | U                   |
| 000121142         | 2,4-DINITROTOLUENE          | 4.0           | ug/L         | U                   |
| 000086737         | FLUORENE                    | 4.0           | ug/L         | U                   |
| 000084662         | DIETHYLPHTHALATE            | 4.0           | ug/L         | U                   |
| 007005723         | 4-CHLOROPHENYL PHENYL ETHER | 4.0           | ug/L         | U                   |



U.S. EPA Region 2 Laboratory  
Data Report

Survey Name: Lackawanna Foundry

Project Number: 00120005

\*Sorted By Sample

**AB03263**

Field/Station ID: 1002-GW06

Coll. Ending Date/Time: 12/5/00 14:05

Matrix: Aqueous

Collection Begin Date/Time: 12/05/00 14:00

Sample Description:

**Analysis Type: NVOA GCMS AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u>               | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|-----------------------------------|---------------|--------------|---------------------|
| 000100016         | 4-NITROANILINE                    | 4.0           | ug/L         | U                   |
| 000534521         | 4,6-DINITRO-2-METHYLPHENOL        | 4.0           | ug/L         | U                   |
| 000086306         | N-NITROSODIPHENYLAMINE            | 4.0           | ug/L         | U                   |
| 000103333         | DIAZENE, DIPHENYL                 | 4.0           | ug/L         | U                   |
| 000101553         | 4-BROMOPHENYL PHENYL ETHER        | 4.0           | ug/L         | U                   |
| 000118741         | HEXACHLOROBENZENE                 | 4.0           | ug/L         | U                   |
| 000087865         | PENTACHLOROPHENOL                 | 10            | ug/L         | U                   |
| 000085018         | PHENANTHRENE                      | 4.0           | ug/L         | U                   |
| 000120127         | ANTHRACENE                        | 4.0           | ug/L         | U                   |
| 000084742         | DI-N-BUTYL PHTHALATE              | 4.0           | ug/L         | U                   |
| 000206440         | FLUORANTHENE                      | 4.0           | ug/L         | U                   |
| 000129000         | PYRENE                            | 4.0           | ug/L         | U                   |
| 000085687         | BUTYL BENZYL PHTHALATE            | 4.0           | ug/L         | U                   |
| 000056553         | BENZO(A)ANTHRACENE                | 4.0           | ug/L         | U                   |
| 000218019         | CHRYSENE                          | 4.0           | ug/L         | U                   |
| 000117817         | BIS(2-ETHYLHEXYL)PHTHALATE        | 4.0           | ug/L         | U                   |
| 000117840         | DI-N-OCTYL PHTHALATE              | 4.0           | ug/L         | U                   |
| 000205992         | BENZO(B)FLUORANTHENE              | 4.0           | ug/L         | U                   |
| 000207089         | BENZO(K)FLUORANTHENE              | 4.0           | ug/L         | U                   |
| 000050328         | BENZO(A)PYRENE                    | 4.0           | ug/L         | U                   |
| 000193395         | INDENO(1,2,3-CD)PYRENE            | 4.0           | ug/L         | U                   |
| 000053703         | DIBENZO(A,H)ANTHRACENE            | 4.0           | ug/L         | U                   |
| 000191242         | BENZO(G,H,I)PERYLENE              | 4.0           | ug/L         | U                   |
|                   | CYCLOHEXANOL, 1-METHYL-, RT =7.44 | 3.0           | ug/L         | J QT                |
|                   | UNKNOWN ,RT =8.45                 | 4.0           | ug/L         | J QT                |
|                   | UNKNOWN ,RT =8.95                 | 4.0           | ug/L         | J QT                |
|                   | UNKNOWN ,RT =9.53                 | 2.0           | ug/L         | J QT                |

**Analysis Type: PESTICIDES TCL GC AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|---------------------|---------------|--------------|---------------------|
| 319846            | ALPHA-BHC           | 0.010         | ug/L         | U                   |
| 58899             | GAMMA-BHC           | 0.010         | ug/L         | U                   |
| 319857            | BETA-BHC            | 0.010         | ug/L         | U                   |
| 319868            | DELTA-BHC           | 0.010         | ug/L         | U                   |
| 76448             | HEPTACHLOR          | 0.010         | ug/L         | U                   |
| 309002            | ALDRIN              | 0.010         | ug/L         | U                   |
| 1024573           | HEPTACHLOR EPOXIDE  | 0.010         | ug/L         | U                   |
| 5566347           | GAMMA-CHLORDANE     | 0.010         | ug/L         | U                   |
| 5103719           | ALPHA-CHLORDANE     | 0.010         | ug/L         | U                   |
| 115297            | ENDOSULFAN I        | 0.010         | ug/L         | U                   |
| 72559             | 4,4'-DDE            | 0.010         | ug/L         | U                   |
| 60571             | DIELDRIN            | 0.010         | ug/L         | U                   |



**Survey Name: Lackawanna Foundry**

Project Number: 00120005

\*Sorted By Sample ID

**AB03263**

Field/Station ID: 1002-GW06

Coll. Ending Date/Time: 12/5/00 14:05

Matrix: Aqueous

Collection Begin Date/Time: 12/05/00 14:00

Sample Description:

**Analysis Type: PESTICIDES TCL GC AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Codes</u> | <u>Remark</u> |
|-------------------|---------------------|---------------|--------------|--------------|---------------|
| 72208             | ENDRIN              | 0.010         | ug/L         | U            |               |
| 72548             | 4,4'-DDD            | 0.010         | ug/L         | U            |               |
| 33213659          | ENDOSULFAN II       | 0.010         | ug/L         | U            |               |
| 50293             | 4,4'-DDT            | 0.010         | ug/L         | U            |               |
| 7421934           | ENDRIN ALDEHYDE     | 0.010         | ug/L         | U            |               |
| 72435             | METHOXYCHLOR        | 0.050         | ug/L         | U            |               |
| 1031078           | ENDOSULFAN SULFATE  | 0.010         | ug/L         | U            |               |
| 53494705          | ENDRIN KETONE       | 0.010         | ug/L         | U            |               |
| 57749             | CHLORDANE           | 0.10          | ug/L         | U            |               |
| 8001352           | TOXAPHENE           | 0.10          | ug/L         | U            |               |

**Analysis Type: PCBS TCL GC AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Codes</u> | <u>Remark</u> |
|-------------------|---------------------|---------------|--------------|--------------|---------------|
| 12674-11-2        | AROCLOR 1016        | 0.10          | ug/L         | U            |               |
| 011104282         | AROCLOR 1221        | 0.20          | ug/L         | U            |               |
| 011141165         | AROCLOR 1232        | 0.10          | ug/L         | U            |               |
| 53469-21-9        | AROCLOR 1242        | 0.10          | ug/L         | U            |               |
| 012672296         | AROCLOR 1248        | 0.10          | ug/L         | U            |               |
| 11097-69-1        | AROCLOR 1254        | 0.10          | ug/L         | U            |               |
| 11096-82-5        | AROCLOR 1260        | 0.10          | ug/L         | U            |               |

**Analysis Type: METALS TAL ICP AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Codes</u> | <u>Remark</u> |
|-------------------|---------------------|---------------|--------------|--------------|---------------|
| 007440224         | SILVER              | 10            | ug/L         | U            |               |
| 007429905         | ALUMINUM            | 200           | ug/L         | U            |               |
| 007440393         | BARIUM              | 380           | ug/L         |              |               |
| 007440417         | BERYLLIUM           | 5.0           | ug/L         | U            |               |
| 007440702         | CALCIUM             | 250,000       | ug/L         |              |               |
| 007440439         | CADMIUM             | 5.0           | ug/L         | U            |               |
| 007440484         | COBALT              | 50            | ug/L         | U            |               |
| 007440473         | CHROMIUM            | 10            | ug/L         | U            |               |
| 007440508         | COPPER              | 25            | ug/L         | U            |               |
| 007439896         | IRON                | 13,000        | ug/L         |              |               |
| 007440097         | POTASSIUM           | 8,000         | ug/L         |              |               |
| 007439954         | MAGNESIUM           | 120,000       | ug/L         |              |               |
| 007439965         | MANGANESE           | 300           | ug/L         |              |               |
| 007440235         | SODIUM              | 550,000       | ug/L         |              |               |
| 007440020         | NICKEL              | 40            | ug/L         | U            |               |
| 007440360         | ANTIMONY            | 60            | ug/L         | U            |               |
| 007440622         | VANADIUM            | 50            | ug/L         | U            |               |
| 007440666         | ZINC                | 20            | ug/L         | U            |               |



**Survey Name: Lackawanna Foundry**

Project Number: 00120005

\*Sorted By Sample

**AB03263** Field/Station ID: 1002-GW06

Coll. Ending Date/Time: 12/5/00 14:05

Matrix: Aqueous

Collection Begin Date/Time: 12/05/00 14:00

Sample Description:

**Single Component Analyses**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|---------------------|---------------|--------------|---------------------|
| 7440-38-2         | ARSENIC             | 10            | ug/L         | U                   |
| 7439-92-1         | LEAD                | 3.0           | ug/L         | U                   |
| 7782-49-2         | SELENIUM            | 5.0           | ug/L         | U                   |
| 7440-28-0         | THALLIUM            | 10            | ug/L         | U QR                |

**Single Component Analyses**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|---------------------|---------------|--------------|---------------------|
| 007439976         | MERCURY (ESAT)      | 0.20          | ug/L         | U                   |

**AB03264** Field/Station ID: 1002-RB02

Coll. Ending Date/Time: 12/5/00 15:05

Matrix: Aqueous

Collection Begin Date/Time: 12/05/00 15:00

Sample Description:

**Analysis Type: VOA TCL GCMS AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u>       | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|---------------------------|---------------|--------------|---------------------|
| 000074873         | CHLOROMETHANE             | 10            | ug/L         | U                   |
| 000075014         | VINYL CHLORIDE            | 10            | ug/L         | U                   |
| 000074839         | BROMOMETHANE              | 10            | ug/L         | U                   |
| 000075003         | CHLOROETHANE              | 10            | ug/L         | U                   |
| 000075354         | 1,1-DICHLOROETHENE        | 10            | ug/L         | U                   |
| 000075150         | CARBON DISULFIDE          | 10            | ug/L         | U                   |
| 000067641         | ACETONE                   | 10            | ug/L         | U                   |
| 000075092         | METHYLENE CHLORIDE        | 10            | ug/L         | U                   |
| 000156605         | TRANS-1,2-DICHLOROETHENE  | 10            | ug/L         | U                   |
| 000156592         | CIS-1,2-DICHLOROETHENE    | 10            | ug/L         | U                   |
| 000075343         | 1,1-DICHLOROETHANE        | 10            | ug/L         | U                   |
| 000078933         | 2-BUTANONE                | 10            | ug/L         | U                   |
| 000067663         | CHLOROFORM                | 10            | ug/L         | U                   |
| 000107062         | 1,2-DICHLOROETHANE        | 10            | ug/L         | U                   |
| 000071556         | 1,1,1-TRICHLOROETHANE     | 10            | ug/L         | U                   |
| 000056235         | CARBON TETRACHLORIDE      | 10            | ug/L         | U                   |
| 000071432         | BENZENE                   | 10            | ug/L         | U                   |
| 025323891         | TRICHLOROETHENE           | 10            | ug/L         | U                   |
| 000078875         | 1,2-DICHLOROPROPANE       | 10            | ug/L         | U                   |
| 000074975         | BROMODICHLOROMETHANE      | 10            | ug/L         | U                   |
| 010061015         | CIS-1,3-DICHLOROPROPENE   | 10            | ug/L         | U                   |
| 010061026         | TRANS-1,3-DICHLOROPROPENE | 10            | ug/L         | U                   |
| 000079005         | 1,1,2-TRICHLOROETHANE     | 10            | ug/L         | U                   |
| 000124481         | DIBROMOCHLOROMETHANE      | 10            | ug/L         | U                   |
| 000075252         | BROMOFORM                 | 10            | ug/L         | U                   |
| 000108101         | 4-METHYL-2-PENTANONE      | 10            | ug/L         | U                   |
| 000108883         | TOLUENE                   | 10            | ug/L         | U                   |
| 000127184         | TETRACHLOROETHENE         | 10            | ug/L         | U                   |



U.S. EPA Region 2 Laboratory  
Data Report

Survey Name: Lackawanna Foundry

Project Number: 00120005

\*Sorted By Sample ID

AB03264

Field/Station ID: 1002-RB02

Coll. Ending Date/Time: 12/5/00 15:05

Matrix: Aqueous

Collection Begin Date/Time: 12/05/00 15:00

Sample Description:

Analysis Type: VOA TCL GCMS AQUEOUS

| CAS Number | Analyte Name              | Result | Units | Codes | Remark |
|------------|---------------------------|--------|-------|-------|--------|
| 000591786  | 2-HEXANONE                | 10     | ug/L  | U     |        |
| 000108907  | CHLOROBENZENE             | 10     | ug/L  | U     |        |
| 000100414  | ETHYLBENZENE              | 10     | ug/L  | U     |        |
| 001330207  | M,P-XYLENE                | 10     | ug/L  | U     |        |
| 000095476  | O-XYLENE                  | 10     | ug/L  | U     |        |
| 000100425  | STYRENE                   | 10     | ug/L  | U     |        |
| 000079345  | 1,1,2,2-TETRACHLOROETHANE | 10     | ug/L  | U     |        |

Analysis Type: NVOA GCMS AQUEOUS

| CAS Number | Analyte Name                | Result | Units | Codes | Remark |
|------------|-----------------------------|--------|-------|-------|--------|
| 000108952  | PHENOL                      | 4.0    | ug/L  | U     |        |
| 000111444  | BIS(2-CHLOROETHYL)ETHER     | 4.0    | ug/L  | U     |        |
| 000095578  | 2-CHLOROPHENOL              | 4.0    | ug/L  | U     |        |
| 000541731  | 1,3-DICHLOROBENZENE         | 4.0    | ug/L  | U     |        |
| 000106467  | 1,4-DICHLOROBENZENE         | 4.0    | ug/L  | U     |        |
| 000095501  | 1,2-DICHLOROBENZENE         | 4.0    | ug/L  | U     |        |
| 000100516  | BENZYL ALCOHOL              | 10     | ug/L  | U     |        |
| 000095487  | 2-METHYLPHENOL              | 4.0    | ug/L  | U     |        |
| 000108601  | BIS(2-CHLOROISOPROPYL)ETHER | 4.0    | ug/L  | U     |        |
| 000106445  | 4-METHYLPHENOL              | 4.0    | ug/L  | U     |        |
| 000621647  | N-NITROSO-DI-N-PROPYLAMINE  | 4.0    | ug/L  | U     |        |
| 000067721  | HEXACHLOROETHANE            | 4.0    | ug/L  | U     |        |
| 000098953  | NITROBENZENE                | 4.0    | ug/L  | U     |        |
| 000078591  | ISOPHORONE                  | 4.0    | ug/L  | U     |        |
| 000088755  | 2-NITROPHENOL               | 4.0    | ug/L  | U     |        |
| 000105679  | 2,4-DIMETHYLPHENOL          | 4.0    | ug/L  | U     |        |
| 000111911  | BIS(2-CHLOROETHOXY)METHANE  | 4.0    | ug/L  | U     |        |
| 000120832  | 2,4-DICHLOROPHENOL          | 4.0    | ug/L  | U     |        |
| 000120821  | 1,2,4-TRICHLOROBENZENE      | 4.0    | ug/L  | U     |        |
| 000065850  | BENZOIC ACID                | 48     | ug/L  | U     |        |
| 000091203  | NAPHTHALENE                 | 4.0    | ug/L  | U     |        |
| 000106478  | 4-CHLOROANILINE             | 48     | ug/L  | U     |        |
| 000087683  | HEXACHLOROBUTADIENE         | 4.0    | ug/L  | U     |        |
| 000059507  | 4-CHLORO-3-METHYLPHENOL     | 4.0    | ug/L  | U     |        |
| 000091576  | 2-METHYL NAPHTHALENE        | 4.0    | ug/L  | U     |        |
| 000077474  | HEXACHLOROCYCLOPENTADIENE   | 4.0    | ug/L  | U     |        |
| 000088062  | 2,4,6-TRICHLOROPHENOL       | 4.0    | ug/L  | U     |        |
| 000095954  | 2,4,5-TRICHLOROPHENOL       | 4.0    | ug/L  | U     |        |
| 000091587  | 2-CHLORONAPHTHALENE         | 4.0    | ug/L  | U     |        |
| 000088744  | 2-NITROANILINE              | 4.0    | ug/L  | U     |        |
| 000131113  | DIMETHYL PHTHALATE          | 4.0    | ug/L  | U     |        |
| 000208968  | ACENAPHTHYLENE              | 4.0    | ug/L  | U     |        |



U.S. EPA Region 2 Laboratory  
Data Report

Survey Name: Lackawanna Foundry

Project Number: 00120005

\*Sorted By Sample

**AB03264**

Field/Station ID: 1002-RB02

Coll. Ending Date/Time: 12/5/00 15:05

Matrix: Aqueous

Collection Begin Date/Time: 12/05/00 15:00

Sample Description:

**Analysis Type: NVOA GCMS AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u>         | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|-----------------------------|---------------|--------------|---------------------|
| 000606202         | 2,6-DINITROTOLUENE          | 4.0           | ug/L         | U                   |
| 000099092         | 3-NITROANILINE              | 48            | ug/L         | U                   |
| 000083329         | ACENAPHTHENE                | 4.0           | ug/L         | U                   |
| 000051285         | 2,4-DINITROPHENOL           | 10            | ug/L         | U                   |
| 000100027         | 4-NITROPHENOL               | 4.0           | ug/L         | U                   |
| 000132649         | DIBENZOFURAN                | 4.0           | ug/L         | U                   |
| 000121142         | 2,4-DINITROTOLUENE          | 4.0           | ug/L         | U                   |
| 000086737         | FLUORENE                    | 4.0           | ug/L         | U                   |
| 000084662         | DIETHYLPHTHALATE            | 4.0           | ug/L         | U                   |
| 007005723         | 4-CHLOROPHENYL PHENYL ETHER | 4.0           | ug/L         | U                   |
| 000100016         | 4-NITROANILINE              | 4.0           | ug/L         | U                   |
| 000534521         | 4,6-DINITRO-2-METHYLPHENOL  | 4.0           | ug/L         | U                   |
| 000086306         | N-NITROSODIPHENYLAMINE      | 4.0           | ug/L         | U                   |
| 000103333         | DIAZENE, DIPHENYL           | 4.0           | ug/L         | U                   |
| 000101553         | 4-BROMOPHENYL PHENYL ETHER  | 4.0           | ug/L         | U                   |
| 000118741         | HEXACHLOROBENZENE           | 4.0           | ug/L         | U                   |
| 000087865         | PENTACHLOROPHENOL           | 10            | ug/L         | U                   |
| 000085018         | PHENANTHRENE                | 4.0           | ug/L         | U                   |
| 000120127         | ANTHRACENE                  | 4.0           | ug/L         | U                   |
| 000084742         | DI-N-BUTYL PHTHALATE        | 4.0           | ug/L         | U                   |
| 000206440         | FLUORANTHENE                | 4.0           | ug/L         | U                   |
| 000129000         | PYRENE                      | 4.0           | ug/L         | U                   |
| 000085687         | BUTYL BENZYL PHTHALATE      | 4.0           | ug/L         | U                   |
| 000056553         | BENZO(A)ANTHRACENE          | 4.0           | ug/L         | U                   |
| 000218019         | CHRYSENE                    | 4.0           | ug/L         | U                   |
| 000117817         | BIS(2-ETHYLHEXYL)PHTHALATE  | 4.0           | ug/L         | U                   |
| 000117840         | DI-N-OCTYL PHTHALATE        | 4.0           | ug/L         | U                   |
| 000205992         | BENZO(B)FLUORANTHENE        | 4.0           | ug/L         | U                   |
| 000207089         | BENZO(K)FLUORANTHENE        | 4.0           | ug/L         | U                   |
| 000050328         | BENZO(A)PYRENE              | 4.0           | ug/L         | U                   |
| 000193395         | INDENO(1,2,3-CD)PYRENE      | 4.0           | ug/L         | U                   |
| 000053703         | DIBENZO(A,H)ANTHRACENE      | 4.0           | ug/L         | U                   |
| 000191242         | BENZO(G,H,I)PERYLENE        | 4.0           | ug/L         | U                   |
|                   | UNKNOWN ,RT =12.24          | 2.0           | ug/L         | I QT                |

**Analysis Type: PESTICIDES TCL GC AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|---------------------|---------------|--------------|---------------------|
| 319846            | ALPHA-BHC           | 0.010         | ug/L         | U                   |
| 58899             | GAMMA-BHC           | 0.010         | ug/L         | U                   |
| 319857            | BETA-BHC            | 0.010         | ug/L         | U                   |
| 319868            | DELTA-BHC           | 0.010         | ug/L         | U                   |
| 76448             | HEPTACHLOR          | 0.010         | ug/L         | U                   |



U.S. EPA Region 2 Laboratory  
Data Report

Survey Name: Lackawanna Foundry

Project Number: 00120005

\*Sorted By Sample ID

AB03264

Field/Station ID: 1002-RB02

Coll. Ending Date/Time: 12/5/00 15:05

Matrix: Aqueous

Collection Begin Date/Time: 12/05/00 15:00

Sample Description:

Analysis Type: PESTICIDES TCL GC AQUEOUS

| CAS Number | Analyte Name       | Result | Units | Codes | Remark |
|------------|--------------------|--------|-------|-------|--------|
| 309002     | ALDRIN             | 0.010  | ug/L  | U     |        |
| 1024573    | HEPTACHLOR EPOXIDE | 0.010  | ug/L  | U     |        |
| 5566347    | GAMMA-CHLORDANE    | 0.010  | ug/L  | U     |        |
| 5103719    | ALPHA-CHLORDANE    | 0.010  | ug/L  | U     |        |
| 115297     | ENDOSULFAN I       | 0.010  | ug/L  | U     |        |
| 72559      | 4,4'-DDE           | 0.010  | ug/L  | U     |        |
| 60571      | DIELDRIN           | 0.010  | ug/L  | U     |        |
| 72208      | ENDRIN             | 0.010  | ug/L  | U     |        |
| 72548      | 4,4'-DDD           | 0.010  | ug/L  | U     |        |
| 33213659   | ENDOSULFAN II      | 0.010  | ug/L  | U     |        |
| 50293      | 4,4'-DDT           | 0.010  | ug/L  | U     |        |
| 7421934    | ENDRIN ALDEHYDE    | 0.010  | ug/L  | U     |        |
| 72435      | METHOXYCHLOR       | 0.050  | ug/L  | U     |        |
| 1031078    | ENDOSULFAN SULFATE | 0.010  | ug/L  | U     |        |
| 53494705   | ENDRIN KETONE      | 0.010  | ug/L  | U     |        |
| 57749      | CHLORDANE          | 0.10   | ug/L  | U     |        |
| 8001352    | TOXAPHENE          | 0.10   | ug/L  | U     |        |

Analysis Type: PCBS TCL GC AQUEOUS

| CAS Number | Analyte Name | Result | Units | Codes | Remark |
|------------|--------------|--------|-------|-------|--------|
| 12674-11-2 | AROCLOR 1016 | 0.10   | ug/L  | U     |        |
| 011104282  | AROCLOR 1221 | 0.20   | ug/L  | U     |        |
| 011141165  | AROCLOR 1232 | 0.10   | ug/L  | U     |        |
| 53469-21-9 | AROCLOR 1242 | 0.10   | ug/L  | U     |        |
| 012672296  | AROCLOR 1248 | 0.10   | ug/L  | U     |        |
| 11097-69-1 | AROCLOR 1254 | 0.10   | ug/L  | U     |        |
| 11096-82-5 | AROCLOR 1260 | 0.10   | ug/L  | U     |        |

Analysis Type: METALS TAL ICP AQUEOUS

| CAS Number | Analyte Name | Result | Units | Codes | Remark |
|------------|--------------|--------|-------|-------|--------|
| 007440224  | SILVER       | 10     | ug/L  | U     |        |
| 007429905  | ALUMINUM     | 200    | ug/L  | U     |        |
| 007440393  | BARIUM       | 200    | ug/L  | U     |        |
| 007440417  | BERYLLIUM    | 5.0    | ug/L  | U     |        |
| 007440702  | CALCIUM      | 5,000  | ug/L  | U     |        |
| 007440439  | CADMIUM      | 5.0    | ug/L  | U     |        |
| 007440484  | COBALT       | 50     | ug/L  | U     |        |
| 007440473  | CHROMIUM     | 10     | ug/L  | U     |        |
| 007440508  | COPPER       | 25     | ug/L  | U     |        |
| 007439896  | IRON         | 100    | ug/L  | U     |        |
| 007440097  | POTASSIUM    | 5,000  | ug/L  | U     |        |
| 007439954  | MAGNESIUM    | 5,000  | ug/L  | U     |        |
| 007439965  | MANGANESE    | 15     | ug/L  | U     |        |





**Survey Name: Lackawanna Foundry**

Project Number: 00120005

\*Sorted By Sample

**AB03264** Field/Station ID: 1002-RB02  
Matrix: Aqueous  
Sample Description:

Coll. Ending Date/Time: 12/5/00 15:05  
Collection Begin Date/Time: 12/05/00 15:00

**Analysis Type: METALS TAL ICP AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|---------------------|---------------|--------------|---------------------|
| 007440235         | SODIUM              | 5,000         | ug/L         | U                   |
| 007440020         | NICKEL              | 40            | ug/L         | U                   |
| 007440360         | ANTIMONY            | 60            | ug/L         | U                   |
| 007440622         | VANADIUM            | 50            | ug/L         | U                   |
| 007440666         | ZINC                | 20            | ug/L         | U                   |

**Single Component Analyses**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|---------------------|---------------|--------------|---------------------|
| 7440-38-2         | ARSENIC             | 10            | ug/L         | U                   |
| 7439-92-1         | LEAD                | 3.0           | ug/L         | U                   |
| 7782-49-2         | SELENIUM            | 5.0           | ug/L         | U                   |
| 7440-28-0         | THALLIUM            | 10            | ug/L         | U                   |

**Single Component Analyses**

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|---------------------|---------------|--------------|---------------------|
| 007439976         | MERCURY (ESAT)      | 0.20          | ug/L         | U                   |

**AB03265** Field/Station ID: 1002-TB01  
Matrix: Aqueous  
Sample Description:

Coll. Ending Date/Time: 12/4/00 00:00

**Analysis Type: VOA TCL GCMS AQUEOUS**

| <u>CAS Number</u> | <u>Analyte Name</u>      | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|--------------------------|---------------|--------------|---------------------|
| 000074873         | CHLOROMETHANE            | 10            | ug/L         | U                   |
| 000075014         | VINYL CHLORIDE           | 10            | ug/L         | U                   |
| 000074839         | BROMOMETHANE             | 10            | ug/L         | U                   |
| 000075003         | CHLOROETHANE             | 10            | ug/L         | U                   |
| 000075354         | 1,1-DICHLOROETHENE       | 10            | ug/L         | U                   |
| 000075150         | CARBON DISULFIDE         | 10            | ug/L         | U                   |
| 000067641         | ACETONE                  | 10            | ug/L         | U                   |
| 000075092         | METHYLENE CHLORIDE       | 10            | ug/L         | U                   |
| 000156605         | TRANS-1,2-DICHLOROETHENE | 10            | ug/L         | U                   |
| 000156592         | CIS-1,2-DICHLOROETHENE   | 10            | ug/L         | U                   |
| 000075343         | 1,1-DICHLOROETHANE       | 10            | ug/L         | U                   |
| 000078933         | 2-BUTANONE               | 10            | ug/L         | U                   |
| 000067663         | CHLOROFORM               | 10            | ug/L         | U                   |
| 000107062         | 1,2-DICHLOROETHANE       | 10            | ug/L         | U                   |
| 000071556         | 1,1,1-TRICHLOROETHANE    | 10            | ug/L         | U                   |
| 000056235         | CARBON TETRACHLORIDE     | 10            | ug/L         | U                   |
| 000071432         | BENZENE                  | 10            | ug/L         | U                   |
| 025323891         | TRICHLOROETHENE          | 10            | ug/L         | U                   |
| 000078875         | 1,2-DICHLOROPROPANE      | 10            | ug/L         | U                   |
| 000074975         | BROMODICHLOROMETHANE     | 10            | ug/L         | U                   |



U.S. EPA Region 2 Laboratory  
Data Report

Survey Name: Lackawanna Foundry

Project Number: 00120005

\*Sorted By Sample ID

**AB03265**

Field/Station ID: 1002-TB01

Coll. Ending Date/Time: 12/4/00 00:00

Matrix: Aqueous

Sample Description:

Analysis Type: VOA TCL GCMS AQUEOUS

| <u>CAS Number</u> | <u>Analyte Name</u>       | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|---------------------------|---------------|--------------|---------------------|
| 010061015         | CIS-1,3-DICHLOROPROPENE   | 10            | ug/L         | U                   |
| 010061026         | TRANS-1,3-DICHLOROPROPENE | 10            | ug/L         | U                   |
| 000079005         | 1,1,2-TRICHLOROETHANE     | 10            | ug/L         | U                   |
| 000124481         | DIBROMOCHLOROMETHANE      | 10            | ug/L         | U                   |
| 000075252         | BROMOFORM                 | 10            | ug/L         | U                   |
| 000108101         | 4-METHYL-2-PENTANONE      | 10            | ug/L         | U                   |
| 000108883         | TOLUENE                   | 10            | ug/L         | U                   |
| 000127184         | TETRACHLOROETHENE         | 10            | ug/L         | U                   |
| 000591786         | 2-HEXANONE                | 10            | ug/L         | U                   |
| 000108907         | CHLOROBENZENE             | 10            | ug/L         | U                   |
| 000100414         | ETHYLBENZENE              | 10            | ug/L         | U                   |
| 001330207         | M,P-XYLENE                | 10            | ug/L         | U                   |
| 000095476         | O-XYLENE                  | 10            | ug/L         | U                   |
| 000100425         | STYRENE                   | 10            | ug/L         | U                   |
| 000079345         | 1,1,2,2-TETRACHLOROETHANE | 10            | ug/L         | U                   |

Project Approval:

*Kenn W. Kulek*

Date:

*1/26/01*



U.S. Environmental Protection Agency  
Region 2 Laboratory

**Data Report: Lackawanna Foundry**

**Project Number: 00120006**

Program: Y206

Project Leader: Donna Janda

| Codes | Explanation  |
|-------|--|
| B     | RESULTS BASED UPON COLONY COUNTS OUTSIDE ACCEPTABLE RANGE        |
| J     | ESTIMATED VALUE  |
| K     | ACTUAL VALUE KNOWN TO BE LESS THAN VALUE GIVEN                   |
| L     | ACTUAL VALUE KNOWN TO BE GREATER THAN VALUE GIVEN                |
| N     | SEE NARRATIVE COMMENTS   |
| V     | SAMPLE RECEIVED BUT NOT ANALYZED DUE TO LAB ACCIDENT             |
| U     | REPORTING LIMIT  |
| QD    | ACCURACY CHECK SAMPLE ABOVE UPPER ACCEPTANCE LIMIT               |
| QE    | ACCURACY CHECK SAMPLE BELOW LOWER ACCEPTANCE LIMIT               |
| QF    | PRECISION OF CALIBRATION CURVE LESS THAN ACCEPTANCE CRITERIA     |
| QJ    | REPORTING LIMIT ESTIMATED DUE TO INTERFERENCE                    |
| QG    | CONTINUING CALIBRATION CHECK DOES NOT MEET ACCEPTANCE CRITERIA   |
| QS    | SPIKE RECOVERIES ABOVE UPPER ACCEPTANCE LIMIT                    |
| QR    | SPIKE RECOVERIES BELOW LOWER ACCEPTANCE LIMIT                    |
| QP    | SAMPLE REPLICATE PRECISION DOES NOT MEET ACCEPTANCE CRITERIA     |
| QH    | RECOMMENDED HOLDING TIME EXCEEDED                                |
| QT    | TENTATIVELY IDENTIFIED COMPOUND                                  |
| QM    | PRESENCE OF MATERIAL VERIFIED BUT NOT QUANTIFIED                 |
| QB    | BLANK CONTAMINATED WITH ANALYTE IN EXCESS OF ACCEPTANCE CRITERIA |
| QQ    | SAMPLE IMPROPERLY MAINTAINED                                     |



test

Survey Name: Lackawanna Foundry

Project Number: 00120006

\*Sorted By Sample ID

AB03266

Field/Station ID: 1002-GW01

Coll. Ending Date/Time: 12/05/00 14:05

Matrix: Aqueous

Collection Begin Date/Time: 12/05/00 14:00

Sample Description:

Single Component Analyses

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|---------------------|---------------|--------------|---------------------|
| 57-12-5           | CYANIDE TOTAL       | 10            | ug/L         | U                   |

AB03267

Field/Station ID: 1002-GW03

Coll. Ending Date/Time: 12/05/00 09:00

Matrix: Aqueous

Collection Begin Date/Time: 12/05/00 08:55

Sample Description:

Single Component Analyses

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|---------------------|---------------|--------------|---------------------|
| 57-12-5           | CYANIDE TOTAL       | 12            | ug/L         |                     |

AB03268

Field/Station ID: 1002-GW04

Coll. Ending Date/Time: 12/05/00 11:05

Matrix: Aqueous

Collection Begin Date/Time: 12/05/00 11:00

Sample Description:

Single Component Analyses

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|---------------------|---------------|--------------|---------------------|
| 57-12-5           | CYANIDE TOTAL       | 10            | ug/L         | U                   |

AB03269

Field/Station ID: 1002-GW05

Coll. Ending Date/Time: 12/05/00 10:00

Matrix: Aqueous

Collection Begin Date/Time: 12/05/00 09:55

Sample Description:

Single Component Analyses

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|---------------------|---------------|--------------|---------------------|
| 57-12-5           | CYANIDE TOTAL       | 10            | ug/L         | U                   |

AB03270

Field Station ID: 1002-GW06

Coll. Ending Date/Time: 12/05/00 14:05

Matrix: Aqueous

Collection Begin Date/Time: 12/05/00 14:00

Sample Description:

Single Component Analyses

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|---------------------|---------------|--------------|---------------------|
| 57-12-5           | CYANIDE TOTAL       | 10            | ug/L         | U                   |

AB03271

Field/Station ID: 1002-RB02

Coll. Ending Date/Time: 12/05/00 15:05

Matrix: Aqueous

Collection Begin Date/Time: 12/05/00 15:00

Sample Description:

Single Component Analyses

| <u>CAS Number</u> | <u>Analyte Name</u> | <u>Result</u> | <u>Units</u> | <u>Remark Codes</u> |
|-------------------|---------------------|---------------|--------------|---------------------|
| 57-12-5           | CYANIDE TOTAL       | 10            | ug/L         | U                   |

Project Approval: Kenn W. Kulda

Date: 1/24/01