

# REPORT

DRAFT

*Approved* *[Signature]*

## *Site Investigation Report*

**Niagara Mohawk Power Corporation  
Oneida (141 Cedar Street) Former MGP Site  
Oneida, New York**

**April 2001**

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BLASLAND, BOUCK & LEE, INC.  
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# 1. Introduction

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This report summarizes field activities and analytical results of the Site Investigation (SI) conducted at the Former Manufactured Gas Plant (MGP) site at 141 Cedar Street in Oneida, New York (the site) (Figure 1). The SI and subsequent remedial program are being implemented in accordance with a Voluntary Cleanup Agreement (VCA) (Index Number: D7-0001-99-04) for the site issued by the NYSDEC in February 2000. Blasland, Bouck & Lee, Inc. (BBL) performed the SI on behalf of Niagara Mohawk Power Corporation (NMPC) to supply the information needed to support a remedial program for the site, anticipated to consist of the excavation and off-site treatment and disposal of fill materials. The SI was performed in accordance with:

- A May 24, 2000 letter from NMPC to the New York State Department of Environmental Conservation (NYSDEC) (NMPC, 2000) that provided the Cedar Street Site Investigation Work Plan;
- The *Oneida (Sconondoa Street) Former MGP Site Preliminary Remedial Design Work Plan* (BBL, 2000) that provided field and laboratory protocols for the investigations at the 141 Cedar Street site; and
- Letters dated July 14, 2000 and November 3, 2000 from NMPC to the NYSDEC that modified and increased the scope of the activities outlined in the May 24, 2000 Work Plan (copies of these letters are provided in Attachment 1).

## 2. Site Setting and Background

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### 2.1 Site Description

The site consists of approximately ¼-acre of paved, unoccupied land on the southeast side of Cedar Street in the City of Oneida (the City) (Figure 2). The City owns this property and the adjacent approximately 2½-acre parcel at 129 Cedar Street, as well as the parcels at 147 and 153 Cedar Street (Figure 2). This SI was conducted to meet the requirements of the VCA, and the subsequent remedial program will be conducted to meet the requirements of the VCA. This remedial program will, in turn, assist the City in the redevelopment of these properties as part of a "Brownfields" program. The City has conducted limited demolition activities at the 153 Cedar Street parcel to make the property more desirable for prospective development. As part of the redevelopment program, it is expected that the 141 Cedar Street site will not be readily useable due to the elevation changes between the site and the adjacent properties (i.e., the site is generally higher than the adjacent properties), and accordingly, regrading of the site is anticipated to be required. The site regrading will require excavation and disposal of the materials currently below the site grade.

The site is generally level and is supported by retaining walls of generally poor structural condition along the northeast, southeast, and southwest boundaries. The retaining walls support fill material used to provide a level grade from Cedar Street to the southeastern site boundary. This fill material is approximately 12 feet deep at the southeastern site boundary, tapering to existing grade at Cedar Street, to the northwest. Land on the adjoining properties generally slopes downward from west to east.

In the central-eastern portion of the site, a slight circular depression in the pavement is apparent in the approximate location of a former gasholder. The depression likely resulted from settlement of backfill inside the former holder.

### 2.2 Site History

A summary of the site history was provided in the November 17, 1999 Phase I Environmental Site Assessment (HYGEIA, 1997) of the adjacent 153 Cedar Street parcel, completed for the City by HYGEIA of N.Y. Inc. (HYGEIA). This assessment found evidence that an MGP, identified as the Oneida Gas Works, was present at 141 Cedar Street sometime before 1890 until no later than 1899. This information is consistent with the general site history discussed in the Stage 1A Cultural Resources Assessment completed for this SI and presented as Attachment 2 to this report. The MGP included a coal shed, a retorts building, a purifying room, and an octagonal gasholder. As discussed in the Cultural Resources Assessment (Attachment 2), a small brass and iron foundry also operated on the southwest portion of the site from before 1890 to sometime between 1895 and 1899. By 1899, the property was occupied by the Oneida Rubber Tire Works and, by 1909, the Coles Tool & Machine Co., which used the former gasholder as a cistern having a 63,000-gallon capacity. By 1923, the gasholder structure had apparently been removed, while the other former MGP structures remained. By 1930, the former MGP structures had been removed, and the site was used for used car sales. By 1956, the building formerly housing the brass foundry had also been removed.

### 2.3 Previous Investigations of Adjacent Parcels

BBL completed a VISTA Information Solutions (VISTA) database search to provide information regarding nearby properties with environmental records. The database search identified 19 listed sites within ¼-mile and a total of 24 sites within a 1-mile radius. Based on the information presented in the database search, none of these sites had

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an apparent environmental impact on the 141 Cedar Street site. The results of this search and a copy of the VISTA report were presented in the Work Plan (NMPC, 2000).

Before BBL's investigation of the site, no site-specific investigations had been completed. Harza Engineering Company (Harza) had completed site investigations of the adjacent City-owned 129 and 153 Cedar Street parcels (Harza, 2000a; 2000b). Harza found no apparent effects of MGP-related constituents in the subsurface or groundwater at the 129 and 153 Cedar Street parcels. Harza also reported that groundwater flow beneath the 129 and 153 Cedar Street parcels is generally from west to east and that groundwater flow from the site is likely toward Oneida Creek, approximately ¼-mile northeast. Some of the monitoring wells installed by Harza are located generally downgradient of the 141 Cedar Street parcel. As presented in the Harza reports, no volatile organic compounds (VOCs), semivolatile organic compounds (SVOCs), pesticides, or polychlorinated biphenyls (PCBs) were detected in any groundwater samples from the 129 and 153 Cedar Street parcels. Cyanide was detected in two groundwater samples collected from monitoring wells MW-2 and MW-4, at the 129 Cedar Street parcel, at a concentration of 1 part per billion (ppb) in each sample. Cyanide was also detected in three groundwater samples collected from monitoring wells MW-1 and MW-3, at the 153 Cedar Street parcel, and from monitoring well MW-2, located upgradient and off site of the 153 Cedar Street parcel, at a concentration of 1 ppb in each sample. These results are orders of magnitude below the New York State Groundwater Quality Standard of 200 ppb for cyanide.

## 2.4 Site Historical Investigation

BBL researched historical information to further understand historical features and operations in the area of the site to assist in defining the nature and extent of MGP impacts, if any, at and in the vicinity of the site. Based on an 1875 Atlas of Madison County (Beer's, 1875), what appears to be a tailrace was located southeast of the site or along the southeastern portion of the site (Figure D.3 of Attachment 2). Additional investigations, beyond those originally proposed in the Work Plan, were completed along the presumed axis of this apparent tailrace to assess potential impacts of former MGP on the tailrace. The results of these investigations are discussed in the following sections.

### 3. Site Investigation Activities and Results

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To facilitate the City's redevelopment of the Cedar Street parcels and to meet the requirements of the VCA, BBL performed SI activities from July to December 2000. The SI provided data necessary to develop and evaluate appropriate soil excavation and disposition alternatives for the site and to develop a soil removal and construction program (including limits of removal) that will be detailed in a Remedial Work Plan in accordance with Paragraph I.D.2 of the VCA. The SI consisted of the following components:

- Site Survey;
- Cultural Resources Assessment;
- Soil Investigation (consisting of waste characterization sampling, Resource Conservation and Recovery Act [RCRA] characterization, and delineation sampling); and
- Groundwater Investigation.

The SI activities were performed in accordance with the Work Plan and, as generally discussed in Section 1, the following supporting documents:

- The *Field Sampling Plan* included in the *Oneida (Sconondoa Street) Former MGP Site Preliminary RD Work Plan* (Sconondoa Street RD Work Plan) (BBL, 2000), which presents field protocols used during the SI;
- The *Quality Assurance Project Plan* included in the Sconondoa Street RD Work Plan, which provides general and laboratory procedures, including quality assurance/quality control (QA/QC) procedures used during the SI; and
- The *Health and Safety Plan (HASP) Addendum*, included as Attachment 1 to the Work Plan, which augments the HASP provided in the Sconondoa Street RD Work Plan. These documents provide project-specific health and safety procedures followed by field personnel during the SI.

Supporting documents from the Sconondoa Street RD Work Plan were used for this site because the SI was conducted in conjunction with the RD activities at the Oneida (Sconondoa Street) Former MGP Site. The sites are close to one another, and the constituents of potential concern are similar at both sites, since both sites supported MGP facilities.

#### 3.1 Site Survey

Before field-sampling activities were conducted, the site was surveyed to create a site map and to assist in anticipated soil removal activities to be conducted as a remedial program. The result of this site survey was the development of a base map, which is provided as Figure 2.

#### 3.2 Cultural Resources Assessment

Binghamton University, as a subcontractor to BBL, conducted a Stage IA Cultural Resources Assessment of the site, consisting of background research, a walkover, and an assessment of the cultural resources. The walkover of

the site was conducted before intrusive aspects of the field investigation began. The results of the Stage IA Cultural Resources Assessment are summarized in Attachment 2. The assessment found that the site has limited research potential with respect to 19th-20th century commerce and industry in the City, and there appears to be limited potential for prehistoric sites within the area of the site. The assessment report found that, based on the results of the Phase IA survey, a Stage IB archeological survey is not warranted for the 141 Cedar Street parcel.

### 3.3 Soil Investigation

The soil investigation program consisted of three components as outlined below.

- *Waste Characterization* - Site soils, which may be removed for off-site treatment, were characterized using the criteria required by Environmental Soil Management of New York, LLC (ESMI). The ESMI criteria were used because NMPC anticipates that the soil to be excavated from the site may contain MGP residuals. (NMPC has used ESMI for treatment of MGP-impacted soils from other sites.)
- *RCRA Characterization* - Site soils that may be removed were characterized using RCRA characterization analyses for ignitability, corrosivity, reactivity, and full Toxicity Characteristic Leaching Procedure (TCLP) analytes to assess whether the soils would be characteristically hazardous.
- *Delineation Sampling* - Delineation sampling was performed to determine if constituents of concern (both MGP-related and non-MGP-related constituents) are present in soils below those that may be excavated during the remedial program, or if MGP-related constituents have affected the surface or subsurface soils of the City-owned adjoining properties. A secondary objective of this program was to locate the fill or native material (e.g., peat, silt) interface, thus defining the thickness of fill material on and around the site.

The soil investigation included drilling 16 soil borings, excavating 5 test pits, and collecting 27 subsurface soil samples to address the objectives stated above. Figure 2 shows the investigation locations, and Table 1 summarizes the analytical samples collected at each location. Soils encountered in the test pits and soil samples collected from the soil borings were characterized by a geologist and screened with a photoionization detector (PID) for the presence of detectable VOCs, if any. Subsurface conditions encountered at each soil boring and test pit are summarized in Attachment 3.

#### 3.3.1 Waste Characterization Sampling

Seven soil samples were collected from soil borings SB-1 through SB-5 for soil characterization (Figure 2). This number of samples meets ESMI's requirements for disposal of 2,500 cubic yards (cy) of soil, the estimated volume of fill expected to be excavated as calculated from the approximate dimensions of the site. In accordance with the Work Plan, samples were collected as follows:

- One shallow soil boring (SB-1) was advanced in the western portion of the site, where the fill is relatively thin.
- Two borings (SB-2 and SB-3) were advanced in the eastern portion of the site, where the fill thickness is greater. From each of these borings, two samples were collected and divided approximately into equal upper and lower intervals.
- Soil borings SB-4A and SB-5 were advanced just outside the holder's perimeter and in the approximate center of the former gasholder, respectively. A sample was collected from each of these borings to represent the entire thickness of fill.

At the request of NYSDEC, a test pit (TP-1) was excavated within the location of the former gasholder to determine the contents and integrity of the former gasholder, if present. The test pit was installed to the presumed bottom of the former gasholder at a depth of approximately 10 feet below ground surface (bgs). Only apparent demolition debris (bricks, concrete, metal) was observed in the location of the former holder, and no MGP-related material was observed in the test pit; therefore, no samples were collected for waste characterization. The test pit was visually characterized by depth and photographed. A test pit log is provided in Attachment 3. The test pit was subsequently backfilled with the material removed from this test pit. The location of this test pit is shown on Figure 2.

Samples collected for soil characterization were homogenized across the sample interval defined in the Work Plan. A summary of the waste characterization analytical results of samples collected for potential treatment and disposal is presented in Table 2. Laboratory reports are provided in Attachment 4. The locations of the soil borings are shown on Figure 2.

### 3.3.2 RCRA Characterization

In accordance with the Work Plan, three discrete soil samples were collected from soil borings SB-2, SB-3, and SB-5 (Figure 2). The samples were submitted for laboratory analysis for ignitability, corrosivity, reactivity, and full TCLP analytes to determine if the soil exhibits the characteristic(s) of a RCRA hazardous waste for treatment and disposal purposes. Sampling locations were determined based on PID readings, odors, staining, or tars observed in the soil. Samples were generally collected from intervals exhibiting characteristics typical of the materials at the site (i.e., not biased toward the most/least visually or PID-impacted soils). Table 1 summarizes the intervals from which these samples were collected.

Based on the analytical results of these samples, the soil is not characterized as a RCRA hazardous waste. However, a homogenized waste characterization soil sample collected (for potential disposal at ESMI) from soil boring SB-3 (6 to 12 feet bgs) contained lead in the TCLP extract at a concentration of 5.46 parts per million (ppm) (regulatory level in the TCLP extract is 5 ppm). All of the other nine results for TCLP lead extract samples were below 0.78 ppm, with a geometric mean of 0.22 ppm for all 10 lead extract samples, suggesting that lead is not a pervasive or widespread issue at the site. Additionally, the material that comprised the sample from SB-3 was noted to consist entirely of shattered brick over 4 feet of the 6-foot sample interval. While brick was observed elsewhere at and near the site, no other subsurface exploration location, completed as part of this investigation, encountered such extensive thicknesses of brick. A summary of the RCRA hazardous waste characterization analytical results is presented in Table 3 and the laboratory reports are provided in Attachment 4.

### 3.3.3 Delineation Sampling

Fifteen discrete soil samples were collected from soil borings and test pits to determine the soil quality at the base of the soil expected to be excavated during the remedial program (i.e., below the depth of the retaining wall) and to define the horizontal extent of MGP-related materials immediately southeast of the site. Ten of these samples were collected from soil borings and submitted for full Target Compound List/Target Analyte List (TCL/TAL) analysis and total cyanide analysis, and five were collected from four off-site test pits and submitted for analysis of polycyclic aromatic hydrocarbons (PAHs). The installation of these additional test pits was in accordance with a November 3, 2000 letter from NMPC to NYSDEC (Attachment 1). Locations of soil samples were determined based on PID readings, odors, staining, or tars observed in the soil. Analytical samples were generally collected from intervals exhibiting no apparent impacts, thus providing a "clean" sample result that would define the limit of the effected materials. Table 1 summarizes the location and interval for these samples, and Figure 2 shows the

locations of the respective soil borings and test pits. Subsurface logs for these investigations are provided in Attachment 3.

During the initial phase of work (installation of soil borings SB-1 through SB-5), tar was observed in only one soil boring (SB-2) at an approximate depth of 16.5 feet bgs. Due to the presence of tar in this boring, four additional soil borings (SB-6 through SB-9) were installed both on and off the 141 Cedar Street parcel during a second mobilization in August 2000. The specific rationale for installing each of the four additional soil borings was presented in a July 14, 2000 letter from NMPC to NYSDEC (provided in Attachment 1). Tar was observed in one of these off-site soil borings (SB-8) at a depth from 4 to 5.2 feet bgs. To further define the extent of MGP-related impacts observed in SB-8 to the southeast of the site, six additional soil borings (SB-10 through SB-15) were installed on the 153 Cedar Street parcel. Tar was observed on gravel from the 4 to 5 feet bgs interval in boring SB-11 and in soil boring SB-12 from 3 to 3.5 feet bgs. Tar, however, was not observed in soil borings SB-10 and SB-15, installed to define the southeastern extent of tar observed in SB-11 and SB-12. In all such cases where tar was observed, the tar was black and highly viscous. A description of the materials encountered in each of the soil borings is provided in Attachment 3.

In accordance with a November 3, 2000 letter from NMPC to NYSDEC (Attachment 1), four additional test pits (TP-2 through TP-5) (Figure 2) were installed on December 7, 2000 on the 153 Cedar Street parcel to assess the possible migration of potentially MGP-related constituents along a former tailrace channel and to further delineate the presence of tar observed in soil borings southeast of the site. Tar was not observed in any of these test pits. While the November 3, 2000 letter discussed only visual characterization of the subsurface materials from these test pits, analytical samples collected from these test pits were analyzed for PAHs to provide further confirmation as to the lack of MGP-related materials in the areas investigated.

The soil sample collected from SB-2 (10 to 12 feet bgs) was not considered "delineation" samples because this sample was collected from within the material assumed to be excavated during the impending remedial program. Therefore, the analytical result of this sample is not included in the discussion of delineation sampling results below. Table 4 presents the soil analytical results, and Figure 3 presents a distribution of total benzene, toluene, ethylbenzene, and xylene (BTEX), total PAHs, and total cyanide detected in the delineation soil samples. A data usability report of the soil delineation data was prepared by BBL and is provided in Attachment 4.

Where appropriate, analytical results from the site investigations of the 129 and 153 Cedar Street parcels are cited for comparison purposes (Harza, 2000a; 2000b). The visual descriptions of soil and the analytical results of one sample collected from a test pit (129-TP-6, collected at 6 feet bgs) (Figure 2), installed immediately north of the site during the 129 Cedar Street site investigation, were used to help define the limits of potentially MGP-related material to the north.

### 3.3.3.1 BTEX

Total BTEX concentrations ranged from nondetect in soil samples collected from soil borings SB-4A (13 to 15 feet bgs), SB-6 (16 to 18 feet bgs and 18 to 20 feet bgs), SB-13 (4 to 6 feet bgs), and SB-15 (4 to 6 feet bgs) to 170 ppm at soil boring SB-2 (19 to 23 feet bgs). The elevated total BTEX concentration in the soil sample collected from SB-2 (19 to 23 feet bgs) is likely associated with the tar observed in this boring at approximately 16.5 feet bgs. BTEX was also detected at 35 ppm in a soil sample collected from soil boring SB-8 (6.5 to 8 feet bgs). Tar was also observed in this soil boring at 4.5 to 6 feet bgs.

Concentrations of BTEX were not detected in a soil sample collected immediately north of the site from test pit TP-6 (6 feet bgs) installed during the 129 Cedar Street investigation. Tar was also not observed in soils from this test pit.

### 3.3.3.2 PAHs

Total PAH concentrations ranged from nondetect for soil samples collected from soil boring SB-6 (16 to 18 feet bgs and 18 to 20 feet bgs) to 1,699.8 ppm in a sample collected from soil boring SB-2 (19 to 23 feet bgs). As discussed above for the total BTEX concentrations in this sample, the elevated total PAH concentration in the sample collected from SB-2 is likely due to the presence of tar observed in a sampling interval above, at 16.5 feet bgs. The next highest total PAH level (1,097.1 ppm) was detected in soil boring SB-1 at 4 to 6 feet bgs. This elevated PAH concentration may be attributed to the presence of treated timbers observed in soil immediately above the 4 to 6 feet bgs interval. All other delineation samples, both on site and off site, had concentrations of total PAHs less than 500 ppm, and only one of those samples (SB-10, 4 to 7 feet bgs) had a concentration greater than 100 ppm (162 ppm). A soil sample collected immediately north of the site from test pit 129-TP-6 (6 feet bgs) (installed during the 129 Cedar Street investigation) contained 29.8 ppm of total PAHs.

### 3.3.3.3 Cyanide

Total cyanide concentrations ranged from nondetect for samples collected from soil borings SB-1 (4 to 6 feet bgs), SB-2 (10 to 12 feet bgs), SB-4A (13 to 15 feet bgs), SB-8 (6.5 to 8 feet bgs), SB-13 (4 to 6 feet bgs), and SB-14 (5 to 7 feet bgs) to 22.4 ppm in a sample collected from SB-3 (12 to 14 feet bgs). The concentration of total cyanide detected in other soil samples ranged from 0.86 to 14.9 ppm. Cyanide was not detected in test pit 129-TP-6 (6 feet bgs) installed immediately north of the site during the 129 Cedar Street investigation.

### 3.3.3.4 PCBs, Pesticides, and Inorganics

PCBs were not detected in any soil samples collected during the SI field activities (including the waste characterization soil sampling). PCBs also were not detected in test pit 129-TP-6 (6 feet bgs) installed immediately north of the site during the 129 Cedar Street investigation.

Pesticides were detected in three of the 10 delineation samples collected for pesticide analysis. One of these samples, SB-2 (19 to 23 feet bgs), did contain a concentration of a heptachlor epoxide (0.061 ppm) slightly above the Recommended Soil Cleanup Objective (0.02 ppm) specified in NYSDEC's Technical Administrative Guidance Memorandum (TAGM) 4046.

The analytical results indicate that all 10 samples analyzed for inorganics contain concentrations of two or more inorganics above the Recommended Soil Cleanup Objectives specified in TAGM 4046. The inorganics detected above the Recommended Soil Cleanup Objective are summarized in the table below.

| Inorganic   | TAGM 4046 (ppm) | Number of Exceedances | Range of Exceedances (ppm) | Highest Detected at 141 Cedar Street | Range of Concentrations Detected at 129/153 Cedar Street |
|---|-----------------|-----------------------|----------------------------|--------------------------------------|--|
| Arsenic   | 7.5 or SB       | 1                     | 34.7                       | SB-2 (19 to 23 feet bgs)             | 1.0 (129-MW-3[8-12]) to 708 (129-MW-2[8-12])             |
| Beryllium   | SB (0.24)       | 5                     | 0.24 to 0.4                | SB-6 (16 to 18 feet bgs)             | 0.2 (153-SB-3[12-16]) to 1.21 (153-TP-2[2.7])            |
| Cadmium   | 10              | 1                     | 10.6                       | SB-13 (4 to 6 feet bgs)<br>Duplicate | 0.15 (129-MW-3[8-12]) to 2.0 (129-MW-1[12-14])           |
| Copper  | 25 or SB        | 3                     | 26.2 to 77.9               | SB-13 (4 to 6 feet bgs)              | 10.5 (153-SB-3[12-16]) to 85.3 (129-TP-1[5])             |
| Iron  | SB (10,200)     | 6                     | 11,000 to 27,900           | SB-2 (19 to 23 feet bgs)             | 7,930 (153-MW-1[4-6]) to 33,200 (153-TP-2[2.7])          |
| Mercury   | 0.1             | 7                     | 0.13 to 1.5                | SB-14 (5 to 7 feet bgs)              | 0.031 (129-MW-1[12-14]) to 6.88 (129-TP-1[5])            |
| Nickel  | 13 or SB        | 1                     | 18.6                       | SB-2 (19 to 23 feet bgs)             | 8.2 (153-SB-3[12-16]) to 28.6 (153-TP-2[2.7])            |
| Zinc  | SB (28.4)       | 5                     | 78 to 9,040J               | SB-13 (4 to 6 feet bgs)              | 24.1 (153-SB-3[12-16]) to 833 (153-TP-1[0-4])            |
| <b>Notes</b>  |                 |                       |                            |                                      |  |
| SB = Site background. Value given in parentheses is from a background sample (153-MW-2 [8 to 12 feet bgs]) collected during the site investigation for the 153 Cedar Street parcel (Harza, 2000b).    |                 |                       |                            |                                      |  |
| J = Concentrations of zinc in 4 of the 5 samples exceeding the site background concentration were all estimated concentrations, and the laboratory duplicate analyses were not within control limits. |                 |                       |                            |                                      |  |



Although several soil samples exhibited concentrations of inorganics at levels exceeding the TAGM 4046 Recommended Soil Cleanup Objectives, with the exception of cadmium and zinc, the concentrations were within the range of the concentration of inorganics detected in subsurface soil samples collected during the site investigations completed at the adjoining 129 and 153 Cedar Street parcels (Harza, 2000a; 2000b). The one soil sample that exceeded the TAGM Recommended Soil Cleanup Objective level for cadmium was only slightly higher (10.6 ppm compared with 10.0 ppm) than the TAGM level, suggesting that this is not a pervasive or widespread issue at the site. Zinc concentrations, however, are generally much higher than the background concentrations and the concentrations observed elsewhere at the adjacent 129 and 153 Cedar Street parcels. This suggests that, with the exception of zinc, the inorganic constituents detected in soil in and around the site are not attributed to the former site activities, but may be attributed to either background or other potential localized sources. Zinc, however, may be present at the site due to the historical use of the property. As indicated by the Cultural Resources Assessment (Attachment 2), an iron and brass foundry existed at the site from before 1890 to sometime between 1895 and 1899. Zinc is a primary component of brass, suggesting that the zinc may be present as a result of the historical use of the site as a brass foundry and not as a result of the use of the site as an MGP. Furthermore, the concentrations of zinc detected in soil on and around the site were well below the NYSDEC TAGM 3028 Action Level and below the United States Environmental Protection Agency (USEPA) Region III Risk-Based Concentration Residential Screening Level of 23,000 ppm.

### 3.4 Groundwater Investigation

The SI activities also included the installation of a temporary monitoring well (PZ-1) at one of the soil boring locations (SB-9) presumed to be located downgradient from the site. This temporary monitoring well was installed on August 10, 2000, and a filtered and nonfiltered groundwater sample was collected from this well on August 15, 2000 for analysis of TCL/TAL constituents and total cyanide, as detailed in the Work Plan. The temporary well was abandoned on December 7, 2000 by hand-pulling the screen and riser and backfilling the borehole with bentonite chips.

VOCs, SVOCs, PCBs, pesticides, and total cyanide were not detected in the filtered or nonfiltered groundwater sample. A total of four inorganic constituents (aluminum, iron, selenium, and zinc) were detected in one or both of the filtered/unfiltered samples at concentrations above the New York State Ambient Groundwater Standards (Technical and Operational Guidance Series [1.1.1] Memorandum, June 1998). As shown in the table below, with the exception of zinc, the levels of these inorganic constituents are consistent with the levels of inorganics detected in nonfiltered groundwater samples collected from monitoring wells during the 129 and 153 Cedar Street site investigations (Harza, 2000a; 2000b). The elevated concentrations of zinc maybe the result of the elevated concentrations of zinc in soil at the 141 Cedar Street parcel. The elevated zinc concentrations in soil are attributed to the historical use of the site as a brass foundry.

| Inorganic | NYS TOGS Standard (ppb) | PZ-1 (ppb)                        | Range at 129/153 Cedar Street (ppb)        |
|-----------|-------------------------|-----------------------------------|--|
| Aluminum  | 100                     | 210 (total)                       | 164 (129-MW-4) to 25000 (129-MW-1)         |
| Iron      | 300                     | 790 (dissolved) & 1,400 (total)   | 294 (129-MW-4) to 68200 (129-MW-1)         |
| Selenium  | 10                      | 19 (dissolved)                    | Not Detected (129-MW-4) to 97.2 (129-MW-1) |
| Zinc      | 300                     | 1,100 (dissolved) & 1,300 (total) | 92.7 (153-MW-3) to 478 (129-MW-1)          |

The full list of analytical results is presented in Table 5. The groundwater analytical results are provided as a "results only" data package in Attachment 4.

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In addition, a round of water level measurements was collected from this temporary well and the existing monitoring wells on the 129 and 153 Cedar Street properties to further assess the groundwater flow patterns near the 141 Cedar Street property. The water level measurements are summarized in Table 6. As shown on Figure 4, shallow groundwater flows in a southeasterly direction.

## 4. Conclusions and Recommendations

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### 4.1 Summary

This section provides a summary of the site conditions encountered and the analytical results completed during this investigation.

#### 4.1.1 Site Conditions

Fill material at the site ranges in thickness from approximately 4 feet near Cedar Street, at soil boring SB-1, to approximately 19 feet near the southern corner of the site, at soil boring SB-2. The fill material on site is composed primarily of a fine to coarse sand and gravel with varying amounts of demolition debris (e.g., brick, concrete, asphalt, wood, and metal) and potential MGP-related waste materials (e.g., cinders, coal, ash, slag, and minor amounts of tar). A soil boring and test pit completed inside of the former gasholder indicates the floor of the holder is at approximately 10 feet bgs. Immediately southeast of the site retaining wall on the 153 Cedar Street parcel, the fill is approximately 5 to 6 feet thick. The upper 2 to 4 feet of fill in this area consists primarily of a brown fine sand, likely resulting from the recent "Brownfields" remediation activities completed on this parcel. The lower portion of the fill in this off-site area is composed of a fine to coarse sand and gravel with demolition debris (e.g., brick, concrete, and wood) and potential MGP-related waste materials (e.g., cinders, slag, and very locally, tar). Both on site and off site, the fill is underlain by a fine sand and silt, peaty silt, or clay, depending on location.

On site, the groundwater table lies approximately 12 to 14 feet bgs, while immediately southeast of the site on the 153 Cedar Street parcel, the water table was observed at approximately 2.5 feet bgs. The difference in depth to water is a direct result of the ground surface elevation difference between the on-site area and the off-site areas. Shallow groundwater flow is generally from west to east across the site. A groundwater sample collected during the SI, approximately 40 feet downgradient from the site at PZ-1, and previous groundwater analytical results from monitoring well 129-MW-4, also located downgradient of the site, indicate that potential MGP-related constituents detected in soil are not affecting groundwater downgradient of the 141 Cedar Street parcel.

#### 4.1.2 Analytical Results

Three distinct types of soil sampling were completed as part of this investigation: waste characterization, to characterize the soil which may be removed from the site for off-site treatment by ESMI; RCRA characterization, to assess whether the soil which may be removed from the site would be classified as characteristically hazardous; and delineation sampling, to assess the nature of the material which would remain on site following a potential soil removal scenario. Each of these is discussed below.

##### 4.1.2.1 Waste Characterization

Based on a discussion with ESMI, ESMI will need to review the data collected to assess whether the soil, which may be removed from the site, would be acceptable for treatment at the ESMI facility. This assessment will be undertaken as part of the Remediation Work Plan development.

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#### 4.1.2.2 RCRA Characterization

None of the samples collected specifically to assess RCRA characterization exceeded any of the RCRA limits. Based on this information the material at the site would be considered nonhazardous. However, one sample collected as part of the ESMI waste characterization program was slightly over the regulatory level for lead and would be considered characteristically hazardous for lead. All of the other nine RCRA characterization samples collected for lead were well below the lead limit, as is the mean lead concentration for all samples. Additionally, the sample that failed for lead was collected from material that was composed mostly of shattered brick, suggesting that the brick at this location may be responsible for the elevated lead concentration.

#### 4.1.2.3 Delineation Sampling

To assess the nature of the material that would remain on site following a potential soil removal program, the delineation sampling was targeted toward collection of native soils below the fill materials at and near the site. The results indicate that, in general, removal of the fill materials at the site and a limited removal of fill off site would leave soil at or near the ground surface, which is similar to the soil quality in background areas or present at the adjacent 129 and 153 Cedar Street parcels. Localized removal of native materials would likely be required in the area of the boring SB-2 (on site) and at and near SB-8/SB-11 (off site) to remove materials observed to contain "tar" and the adjacent underlying soil containing elevated concentrations of BTEX and/or PAHs. Additional native soil removal would be required in the vicinity of SB-1 where treated timbers and associated underlying PAH/BTEX-containing native soils were encountered. Based on a meeting between NMPC, BBL, NYSDEC, and the City of Oneida at the site on October 19, 2000 to discuss the "Brownfields" program and the 129, 141, and 153 Cedar Street parcels, the meeting attendees discussed the use of a visual MGP-impacted removal criteria combined with a 500 ppm total PAH remediation level, assuming a cover/cap will be placed over the area and used for commercial purposes, as an acceptable remediation level for the 141 Cedar Street site and areas immediately adjacent to the site. This level has been used to implement remedial activities at other MGP sites under the consent of the NYSDEC. With the exception of the soil samples collected from on-site soil borings SB-1 (4 to 6 feet bgs) and SB-2 (19 to 23 feet bgs), the total PAH concentrations detected in non-fill soil samples collected during the SI, both on site and off site, are less than 500 ppm. Because the elevated BTEX concentrations were detected only at the locations where there were elevated total PAH concentrations, removal of the soil with elevated PAH concentrations and tar, would concurrently remove the soil with elevated BTEX concentrations.

The delineation sampling also included the analysis of cyanide (all samples below TAGM 4046 Recommended Soil Cleanup Objectives), PCBs (all nondetect), pesticides (all below TAGM 4046 Objectives, except of one sample which would be removed based on the elevated PAH concentrations) and inorganic compounds. All inorganic compounds with the exception of cadmium and zinc were either below TAGM 4046 Objectives, or below concentrations of these metals on the adjacent 129 and 153 Cedar Street parcels. Cadmium was only slightly above the TAGM objective (10.6 ppm vs. 10.0 ppm). Zinc concentrations were above the background levels; however, a brass foundry (zinc is a primary constituent of brass) was located at the site before 1900, and may be the reason that elevated zinc concentrations are present in site. The concentrations of zinc detected in soil on and around the site were below the TAGM 3028 Action Level and USEPA Region III Risk Based Concentration Residential Screening Level of 23,000 ppm.

#### 4.1.3 Fill Materials

As discussed above, a limited extent of MGP-impacted material was observed in the native materials underlying the site and in the fill/native materials outside the retaining wall along the southeastern site boundary, as indicated

by tar-containing soils in SB-2 on site and at two borings (SB-8 and SB-11) immediately southeast of the 141 Cedar Street retaining wall. Although the tar-containing soils observed in SB-2, SB-8, and SB-11 are believed to be delineated, the extent of impacted soil at SB-1 is not well defined. Elevated PAH and BTEX concentrations were detected in the soil sample from SB-1 (4 to 6 feet bgs); however, this sample was collected immediately below an interval of soil containing treated timbers. The elevated concentration of the constituents detected in soil boring SB-1 (4 to 6 feet bgs) may be attributed to the presence of treated timbers observed immediately above the 4 to 6 feet bgs interval.

## 4.2 Recommendations

To assess the distribution of the PAHs and BTEX detected in SB-1, NMPC recommends limited additional investigation activities in this area. To address this data gap, NMPC will provide the NYSDEC with a letter work plan.

Upon completion of the additional field activities and evaluation of the resulting data, NMPC will provide the NYSDEC with a brief letter report discussing the findings of the additional work. Based on those findings, NMPC will develop a remedial program to remove the material within the retaining wall and MGP-related materials observed outside the retaining wall for subsequent off-site disposal/treatment in accordance with the VCA for this site. The Remediation Work Plan will be developed and submitted to the NYSDEC upon the NYSDEC's acceptance of this summary report.

## 5. References

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# ***Tables***

BLASLAND, BOUCK & LEE, INC. *engineers & scientists*

*consultants with focus*

TABLE 1

**NIAGARA MOHAWK POWER CORPORATION  
ONEIDA (141 CEDAR STREET) FORMER MGP SITE  
SITE INVESTIGATION**

**SAMPLE SUMMARY**

| Location ID | Depth      | Date    | Type | Matrix | RCRA Hazardous Waste Sampling | Treatment/Disposal Sampling (ESMI) | Delineation Sampling |
|-------------|------------|---------|------|--------|-------------------------------|------------------------------------|----------------------|
| SB-1        | (0 - 4')   | 7/18/00 | FS   | Soil   |                               | X                                  |                      |
| SB-1        | (4 - 6')   | 7/18/00 | FS   | Soil   |                               |                                    | X                    |
| SB-2        | (0 - 2')   | 7/19/00 | FS   | Soil   | X                             |                                    |                      |
| SB-2        | (0 - 6')   | 7/19/00 | FS   | Soil   |                               | X                                  |                      |
| SB-2        | (6 - 10')  | 7/19/00 | FS   | Soil   |                               | X                                  |                      |
| SB-2        | (10 - 12') | 7/19/00 | FS   | Soil   |                               |                                    | X                    |
| SB-2        | (10 - 12') | 7/19/00 | DUP  | Soil   |                               |                                    | X                    |
| SB-2        | (19 - 23') | 7/19/00 | FS   | Soil   |                               |                                    | X                    |
| SB-3        | (0 - 6')   | 7/19/00 | FS   | Soil   |                               | X                                  |                      |
| SB-3        | (6 - 12')  | 7/19/00 | FS   | Soil   |                               | X                                  |                      |
| SB-3        | (10 - 12') | 7/19/00 | FS   | Soil   | X                             |                                    |                      |
| SB-3        | (12 - 14') | 7/19/00 | FS   | Soil   |                               |                                    | X                    |
| SB-4A       | (7 - 13')  | 7/20/00 | FS   | Soil   |                               | X                                  |                      |
| SB-4A       | (13 - 15') | 7/20/00 | FS   | Soil   |                               |                                    | X                    |
| SB-5        | (0 - 8')   | 7/20/00 | FS   | Soil   |                               | X                                  |                      |
| SB-5        | (8 - 10')  | 7/20/00 | FS   | Soil   | X                             |                                    |                      |
| SB-6        | (16 - 18') | 8/8/00  | FS   | Soil   |                               |                                    | X                    |
| SB-6        | (18 - 20') | 8/8/00  | FS   | Soil   |                               |                                    | X                    |
| SB-8        | (6.5 - 8') | 8/10/00 | FS   | Soil   |                               |                                    | X                    |
| SB-10       | (4 - 7')   | 8/10/00 | FS   | Soil   |                               |                                    | X                    |
| SB-13       | (4 - 6')   | 8/10/00 | FS   | Soil   |                               |                                    | X                    |
| SB-13       | (4 - 6')   | 8/10/00 | DUP  | Soil   |                               |                                    | X                    |
| SB-14       | (5 - 7')   | 8/10/00 | FS   | Soil   |                               |                                    | X                    |
| SB-15       | (4 - 6')   | 8/10/00 | FS   | Soil   |                               |                                    | X                    |
| TP-2        | (3')       | 12/7/00 | FS   | Soil   |                               |                                    | X <sup>1</sup>       |
| TP-2        | (3')       | 12/7/00 | DUP  | Soil   |                               |                                    | X <sup>1</sup>       |
| TP-3        | (6')       | 12/7/00 | FS   | Soil   |                               |                                    | X <sup>1</sup>       |
| TP-4        | (6')       | 12/7/00 | FS   | Soil   |                               |                                    | X <sup>1</sup>       |
| TP-4        | (3')       | 12/7/00 | FS   | Soil   |                               |                                    | X <sup>1</sup>       |
| TP-5        | (5')       | 12/7/00 | FS   | Soil   |                               |                                    | X <sup>1</sup>       |
| PZ-1        |            | 8/15/00 | FS   | Water  |                               |                                    | X                    |

**Notes:**

- <sup>1</sup>Sampled for PAHs only.
- RCRA Hazardous Waste Sampling = TCLP VOCs, TCLP SVOCs, TCLP Pesticides/Herbicides, TCLP metals, Corrosivity, pH, Reactive Cyanide, Ignitability, and Reactive Sulfide. Analyses completed by Galson Laboratories.
- Treatment/Disposal Sampling (ESMI) = SVOCs, TCLP Metals, PCBs, Total Benzene, % Sulfur, Total organic halides (TOX), Total petroleum hydrocarbons (TPH), and Total cyanide. Analyses completed by Phoenix Environmental Laboratories, Inc.
- Delineation Sampling = TCL VOCs (NYSDEC ASP 95-1 Methods), TCL SVOCs and PAHs (NYSDEC ASP 95-2 Methods), TCL Pesticides (NYSDEC ASP 95-3 Methods), TCL PCBs (NYSDEC ASP 95-3 Methods), and TAL Inorganics (Method CLP-M). Analyses completed by Galson Laboratories.



TABLE 2

**NIAGARA MOHAWK POWER CORPORATION  
ONEIDA (141 CEDAR STREET) FORMER MGP SITE  
SITE INVESTIGATION**

**SOIL WASTE CHARACTERIZATION RESULTS - POTENTIAL DISPOSAL/TREATMENT AT ESMI**

| Sample ID                                     | SB-1    | SB-2    | SB-2    | SB-3    | SB-3    | SB-4A   | SB-5    |
|---|---------|---------|---------|---------|---------|---------|---------|
| Date Collected                                | 7/18/00 | 7/19/00 | 7/19/00 | 7/19/00 | 7/19/00 | 7/20/00 | 7/20/00 |
| Sample Depth (ft.)                            | 0-4     | 0-6     | 6-10    | 0-6     | 6-12    | 7-13    | 0-8     |
| <b>Metals by TCLP(mg/L)</b>                   |         |         |         |         |         |         |         |
| Mercury                                       | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U |
| Silver  | 0.01 U  | 0.01 U  | 0.01 U  | 0.01 U  | 0.01 U  | 0.01 U  | 0.01 U  |
| Arsenic                                       | 0.03    | 0.01 U  | 0.01 U  | 0.02    | 0.01 U  | 0.04    | 0.01 U  |
| Barium  | 0.49    | 0.77    | 1.61    | 0.47    | 0.49    | 0.49    | 1.64    |
| Cadmium                                       | 0.005 U | 0.005 U | 0.019   | 0.005 U | 0.008   | 0.005 U | 0.006   |
| Chromium                                      | 0.01 U  | 0.01 U  | 0.01 U  | 0.01 U  | 0.01 U  | 0.01 U  | 0.01 U  |
| Lead  | 0.101   | 0.101   | 0.406   | 0.129   | 5.46    | 0.096   | 0.147   |
| Selenium                                      | 0.05 U  | 0.05 U  | 0.05 U  | 0.05 U  | 0.05 U  | 0.05 U  | 0.05 U  |
| <b>PCBs (ug/Kg)</b>                           |         |         |         |         |         |         |         |
| Aroclor-1016                                  | 400 U   | 400 U   | 400 U   | 400 U   | 400 U   | 400 U   | 400 U   |
| Aroclor-1221                                  | 400 U   | 400 U   | 400 U   | 400 U   | 400 U   | 400 U   | 400 U   |
| Aroclor-1232                                  | 400 U   | 400 U   | 400 U   | 400 U   | 400 U   | 400 U   | 400 U   |
| Aroclor-1242                                  | 400 U   | 400 U   | 400 U   | 400 U   | 400 U   | 400 U   | 400 U   |
| Aroclor-1248                                  | 400 U   | 400 U   | 400 U   | 400 U   | 400 U   | 400 U   | 400 U   |
| Aroclor-1254                                  | 400 U   | 400 U   | 400 U   | 400 U   | 400 U   | 400 U   | 400 U   |
| Aroclor-1260                                  | 400 U   | 400 U   | 400 U   | 400 U   | 400 U   | 400 U   | 400 U   |
| Aroclor-1262                                  | 400 U   | 400 U   | 400 U   | 400 U   | 400 U   | 400 U   | 400 U   |
| Aroclor-1268                                  | 400 U   | 400 U   | 400 U   | 400 U   | 400 U   | 400 U   | 400 U   |
| <b>Total Benzene (ug/Kg)</b>                  | 4400    | 5.0 U   | 500 U   | 5.0 U   | 5.0 U   | 5.0 U   | 5.0 U   |
| <b>Sulfur %</b>                               | 0.2     | 0.065   | 0.056   | 0.087   | 0.082   | 0.02    | 0.08    |
| <b>TOX (mg/Kg)</b>                            | 25 U    | 25 U    | 25 U    | 25 U    | 25 U    | 25 U    | 25 U    |
| <b>TPH (mg/Kg)</b>                            | 240     | 40 U    | 160     | 720     | 40 U    | 23 U    | 540     |
| <b>Total Cyanide (mg/Kg)</b>                  | 2.87    | 0.22 U  | 51.8    | 0.296   | 2.24    | 0.21 U  | 1.12    |
| <b>Semivolatile Organic Compounds (mg/Kg)</b> |         |         |         |         |         |         |         |
| 1,2,4-Trichlorobenzene                        | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| 1,2-Dichlorobenzene                           | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| 1,2-Diphenylhydrazine                         | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| 1,3-Dichlorobenzene                           | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| 1,4-Dichlorobenzene                           | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| 2,4,5-Trichlorophenol                         | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| 2,4,6-Trichlorophenol                         | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| 2,4-Dichlorophenol                            | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| 2,4-Dimethylphenol                            | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| 2,4-Dinitrophenol                             | 160 U   | 1.6 U   | 16 U    | 1.6 U   | 1.6 U   | 1.6 U   | 1.6 U   |
| 2,4-Dinitrotoluene                            | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| 2,6-Dichlorophenol                            | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| 2,6-Dinitrotoluene                            | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |

See Notes on Page 3.

TABLE 2

**NIAGARA MOHAWK POWER CORPORATION  
ONEIDA (141 CEDAR STREET) FORMER MGP SITE  
SITE INVESTIGATION**

**SOIL WASTE CHARACTERIZATION RESULTS - POTENTIAL DISPOSAL/TREATMENT AT ESMI**

| Sample ID   | SB-1    | SB-2    | SB-2    | SB-3    | SB-3    | SB-4A   | SB-5    |
|---|---------|---------|---------|---------|---------|---------|---------|
| Date Collected  | 7/18/00 | 7/19/00 | 7/19/00 | 7/19/00 | 7/19/00 | 7/20/00 | 7/20/00 |
| Sample Depth (ft.)  | 0-4     | 0-6     | 6-10    | 0-6     | 6-12    | 7-13    | 0-8     |
| <b>Semivolatile Organic Compounds (mg/Kg) (continued)</b> |         |         |         |         |         |         |         |
| 2-Chloronaphthalene                                       | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| 2-Chlorophenol  | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| 2-Methylnaphthalene                                       | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| 2-Methylphenol (o-cresol)                                 | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| 2-Nitroaniline  | 160 U   | 1.6 U   | 16 U    | 1.6 U   | 1.6 U   | 1.6 U   | 1.6 U   |
| 2-Nitrophenol   | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| 3,3'-Dichlorobenzidine                                    | 66 U    | 0.66 U  | 6.6 U   | 0.66 U  | 0.66 U  | 0.66 U  | 0.66 U  |
| 3-Nitroaniline  | 160 U   | 1.6 U   | 16 U    | 1.6 U   | 1.6 U   | 1.6 U   | 1.6 U   |
| 4,6-Dinitro-2-methylphenol                                | 160 U   | 1.6 U   | 16 U    | 1.6 U   | 1.6 U   | 1.6 U   | 1.6 U   |
| 4-Bromophenyl phenyl ether                                | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| 4-Chloro-3-methylphenol                                   | 66 U    | 0.66 U  | 6.6 U   | 0.66 U  | 0.66 U  | 0.66 U  | 0.66 U  |
| 4-Chloroaniline   | 66 U    | 0.66 U  | 6.6 U   | 0.66 U  | 0.66 U  | 0.66 U  | 0.66 U  |
| 4-Chlorophenyl phenyl ether                               | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| 4-Methylphenol (o-cresol)                                 | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| 4-Nitroaniline  | 160 U   | 1.6 U   | 16 U    | 1.6 U   | 1.6 U   | 1.6 U   | 1.6 U   |
| 4-Nitrophenol   | 160 U   | 1.6 U   | 16 U    | 1.6 U   | 1.6 U   | 1.6 U   | 1.6 U   |
| Acenaphthene  | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| Acenaphthylene  | 53      | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| Anthracene  | 87      | 0.33 U  | 8.5     | 0.43    | 0.37    | 0.33 U  | 0.33 U  |
| Benzidine   | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| Benzo[a]anthracene  | 160     | 0.33 U  | 12      | 1.5     | 1.7     | 0.33 U  | 1.1     |
| Benzo[a]pyrene  | 110     | 0.33 U  | 7.6     | 2.0     | 1.7     | 0.33 U  | 1.2     |
| Benzo[b]fluoranthene                                      | 160     | 0.33 U  | 9.0     | 2.4     | 2.8     | 0.33 U  | 1.5     |
| Benzo[g,h,i]perylene                                      | 34      | 0.33 U  | 3.3 U   | 0.38    | 0.33 U  | 0.33 U  | 0.33 U  |
| Benzo[k]fluoranthene                                      | 76      | 0.33 U  | 8.4     | 2.4     | 1.8     | 0.33 U  | 2.4     |
| Benzoic acid  | 160 U   | 1.6 U   | 16 U    | 1.6 U   | 1.6 U   | 1.6 U   | 1.6 U   |
| Benzyl alcohol  | 66 U    | 0.66 U  | 6.6 U   | 0.66 U  | 0.66 U  | 0.66 U  | 0.66 U  |
| Butyl benzyl phthalate                                    | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| bis (2-Chloroethoxy) methane                              | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| bis (2-Chloroethyl) ether                                 | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| bis (2-Chloroisopropyl) ether                             | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| bis(2-Ethylhexyl)phthalate                                | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| Chrysene  | 130     | 0.33 U  | 99      | 1.2     | 1.3     | 0.33 U  | 0.65    |
| Di-n-butyl phthalate                                      | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| Di-n-octyl phthalate                                      | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| Dibenz[a,h]anthracene                                     | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| Dibenzofuran  | 37      | 0.33 U  | 4.5     | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |

See Notes on Page 3.

TABLE 2

**NIAGARA MOHAWK POWER CORPORATION  
ONEIDA (141 CEDAR STREET) FORMER MGP SITE  
SITE INVESTIGATION**

**SOIL WASTE CHARACTERIZATION RESULTS - POTENTIAL DISPOSAL/TREATMENT AT ESMI**

| Sample ID   | SB-1    | SB-2    | SB-2    | SB-3    | SB-3    | SB-4A   | SB-5    |
|---|---------|---------|---------|---------|---------|---------|---------|
| Date Collected  | 7/18/00 | 7/19/00 | 7/19/00 | 7/19/00 | 7/19/00 | 7/20/00 | 7/20/00 |
| Sample Depth (ft.)  | 0-4     | 0-6     | 6-10    | 0-6     | 6-12    | 7-13    | 0-8     |
| <b>Semivolatile Organic Compounds (mg/Kg) (continued)</b> |         |         |         |         |         |         |         |
| Diethyl phthalate   | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| Dimethyl phthalate  | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| Fluoranthene  | 330     | 0.36    | 23      | 2.2     | 3.1     | 0.33    | 1.2     |
| Fluorene  | 47      | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| Hexachlorobenzene   | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| Hexachlorobutadiene                                       | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| Hexachlorocyclopentadiene                                 | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| Hexachloroethane  | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| Indeno[1,2,3-cd]pyrene                                    | 38      | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| Isophorone  | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| N-Nitroso-di-n-propylamine                                | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| N-Nitrosodimethylamine                                    | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| N-Nitrosodiphenylamine                                    | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| Naphthalene   | 180     | 0.33 U  | 6.9     | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| Nitrobenzene  | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| Pentachlorophenol   | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| Phenanthrene  | 290     | 0.4     | 25      | 1.6     | 1.4     | 0.33 U  | 0.63    |
| Phenol  | 33 U    | 0.33 U  | 3.3 U   | 0.33 U  | 0.33 U  | 0.33 U  | 0.33 U  |
| Pyrene  | 260     | 0.33 U  | 17      | 1.8     | 2.6     | 0.33 U  | 0.99    |

**Notes:**

- Analytical methods used were: Toxicity Characteristic Leaching Procedure (TCLP) extraction method E1311 for metals by USEPA SW-846 Method 6010B/7470, Total benzene by USEPA SW-846 Method 8021, Sulfur by ASTM Method D129, Total organic halides (TOX) by USEPA SW-846 Method 9092, Total petroleum hydrocarbons (TPH) by USEPA Method 8100, Total cyanide by USEPA SW-846 Method 9010, Polychlorinated biphenyls (PCBs) by USEPA SW-846 Method 8082, Semi-volatile organic compounds (SVOCs) by USEPA SW-846 Method 8270.
- The laboratory analytical results were reported as "results only" data packages.
- U = Compound was not detected at a concentration exceeding the laboratory detection limit.
- mg/L = milligrams per liter.
- ug/Kg = micrograms per kilogram.
- mg/Kg = milligrams per kilogram.
- Samples were homogenized across the entire sample depth interval.

TABLE 3

**NIAGARA MOHAWK POWER CORPORATION  
ONEIDA (141 CEDAR STREET) FORMER MGP SITE  
SITE INVESTIGATION**

**SOIL ANALYTICAL RESULTS - RCRA HAZARDOUS WASTE CHARACTERIZATION**

| Sample ID                                 | SB-2     | SB-3     | SB-5     |
|---|----------|----------|----------|
| Date Collected                            | 7/19/00  | 7/19/00  | 7/20/00  |
| Sample depth (ft.)                        | 0-2      | 10-12    | 8-10     |
| <b>VOCs by TCLP(ug/L)</b>                 |          |          |          |
| Benzene                                   | 50 U     | 50 U     | 50 U     |
| 2-Butanone                                | 100 U    | 100 U    | 110 *    |
| Carbon Tetrachloride                      | 50 U     | 50 U     | 50 U     |
| Chlorobenzene                             | 50 U     | 50 U     | 50 U     |
| Chloroform                                | 50 U     | 50 U     | 50 U     |
| 1,2-Dichloroethane                        | 50 U     | 50 U     | 50 U     |
| 1,1-Dichloroethene                        | 50 U     | 50 U     | 50 U     |
| Tetrachloroethene                         | 50 U     | 50 U     | 50 U     |
| Trichloroethene                           | 50 U     | 50 U     | 50 U     |
| Vinyl Chloride                            | 50 U     | 50 U     | 50 U     |
| <b>SVOCs by TCLP(ug/L)</b>                |          |          |          |
| Pyridine                                  | 100 U    | 100 U    | 100 U    |
| 1,4-Dichlorobenzene                       | 100 U    | 100 U    | 100 U    |
| 2-Methylphenol                            | 100 U    | 18 J     | 100 U    |
| 3 & 4-Methylphenol                        | 200 U    | 200 U    | 200 U    |
| Hexachloroethane                          | 100 U    | 100 U    | 100 U    |
| Nitrobenzene                              | 100 U    | 100 U    | 100 U    |
| Hexachlorobutadiene                       | 100 U    | 100 U    | 100 U    |
| 2,4,6-Trichlorophenol                     | 100 U    | 100 U    | 100 U    |
| 2,4,5-Trichlorophenol                     | 100 U    | 100 U    | 100 U    |
| 2,4-Dinitrotoluene                        | 100 U    | 100 U    | 100 U    |
| Hexachlorobenzene                         | 100 U    | 100 U    | 100 U    |
| Pentachlorophenol                         | 250 U    | 250 U    | 250 U    |
| <b>Pesticide/Herbicides by TCLP(ug/l)</b> |          |          |          |
| Gamma-BHC (lindane)                       | 5.0 U    | 5.0 U    | 5.0 U    |
| Chlordane                                 | 20 U     | 20 U     | 20 U     |
| Endrin                                    | 5.0 U    | 5.0 U    | 5.0 U    |
| Heptachlor                                | 5.0 U    | 5.0 U    | 5.0 U    |
| Heptachlor Epoxide                        | 5.0 U    | 5.0 U    | 5.0 U    |
| Metoxychlor                               | 20 U     | 20 U     | 20 U     |
| Toxaphene                                 | 100 U    | 100 U    | 100 U    |
| 2,4-Dinitrotoluene                        | 50 U     | 50 U     | 50 U     |
| 2,4,5-TP (Silvex)                         | 50 U     | 50 U     | 50 U     |
| <b>Metals by TCLP(mg/L)</b>               |          |          |          |
| Mercury                                   | 0.0003 U | 0.0003 U | 0.0003 U |
| Arsenic                                   | 0.01 U   | 0.016    | 0.01 U   |

See Notes on Page 2.

TABLE 3

**NIAGARA MOHAWK POWER CORPORATION  
ONEIDA (141 CEDAR STREET) FORMER MGP SITE  
SITE INVESTIGATION**

**SOIL ANALYTICAL RESULTS - RCRA HAZARDOUS WASTE CHARACTERIZATION**

| Sample ID                               | SB-2    | SB-3    | SB-5    |
|---|---------|---------|---------|
| Date Collected                          | 7/19/00 | 7/19/00 | 7/20/00 |
| Sample depth (ft.)                      | 0-2     | 10-12   | 8-10    |
| <b>Metals by TCLP(mg/L) (continued)</b> |         |         |         |
| Barium                                  | 1.8     | 1 U     | 1.3     |
| Cadmium                                 | 0.0087  | 0.0061  | 0.037   |
| Chromium                                | 0.01 U  | 0.01 U  | 0.01 U  |
| Lead                                    | 0.14    | 0.074   | 0.78    |
| Selenium                                | 0.02 U  | 0.02 U  | 0.02 U  |
| Silver                                  | 0.01 U  | 0.01 U  | 0.01 U  |
| Corrosivity/pH (SU)                     | 10.3    | 8.0     | 7.7     |
| Reactive Cyanide (mg/kg)                | 100 U   | 100 U   | 100 U   |
| Ignitability                            | NEG     | NEG     | NEG     |
| Reactive Sulfide (mg/kg)                | 100 U   | 100 U   | 100 U   |

**Notes:**

1. Analytical methods used were: Toxicity Characteristic Leaching Procedure (TCLP) extraction method E1311 for volatile organic compounds (VOCs) by United States Environmental Protection Agency (USEPA) SW-846 Method 8260, semi-volatile organic compounds (SVOCs) by USEPA SW-846 Method 8270, Pesticide/Herbicide by USEPA SW-846 Method 8151A, and Metals by USEPA SW-846 Method 6010/7470; and Reactivity by USEPA SW-846 Chapter 7; Corrosivity by USEPA SW-846 Method 9045; and Ignitability by USEPA SW-846 Method 1030.
2. The laboratory analytical results were reported as "results only" data packages.
3. U = Compound was not detected at a concentration exceeding the laboratory detection limit.
4. ug/L = micrograms per liter.
5. mg/L = milligrams per liter.
6. mg/Kg = milligrams per kilogram.
7. NEG = Sample did not ignite or support combustion.
8. SU = Standard Units.
9. \* = This compound was detected in the sample as well as the associated TCLP blank.

TABLE 4

NIAGARA MOHAWK POWER CORPORATION  
 ONEIDA (141 CEDAR STREET) FORMER MGP SITE  
 SITE INVESTIGATION

SOIL ANALYTICAL RESULTS - DELINEATION SAMPLING

| Location ID<br>Depth Range<br>Date Sampled<br>Sample Type | SB-1<br>(4 - 6')<br>7/18/00<br>FS | SB-2<br>(10 - 12')<br>7/19/00<br>FS | SB-2<br>(10 - 12')<br>7/19/00<br>DUP | SB-2<br>(19 - 23')<br>7/19/00<br>FS | SB-3<br>(12 - 14')<br>7/19/00<br>FS | SB-4A<br>(13 - 15')<br>7/20/00<br>FS | SB-6<br>(16 - 18')<br>8/8/00<br>FS | SB-6<br>(18 - 20')<br>8/8/00<br>FS | SB-8<br>(6.5 - 8')<br>8/10/00<br>FS | SB-10<br>(4 - 7')<br>8/10/00<br>FS |
|---|-----------------------------------|-------------------------------------|--------------------------------------|-------------------------------------|-------------------------------------|--------------------------------------|------------------------------------|------------------------------------|-------------------------------------|------------------------------------|
| <b>TCL Volatile Organic Compounds</b>                     |                                   |                                     |                                      |                                     |                                     |                                      |                                    |                                    |                                     |                                    |
| 1,1,1-Trichloroethane                                     | 1.4 U                             | 0.013 U                             | 0.014 U                              | 4.9 U                               | --                                  | 0.012 U                              | 0.012 U                            | 0.012 U                            | 0.1 U                               | 0.012 U                            |
| 1,1,2,2-Tetrachloroethane                                 | 1.4 U                             | 0.013 UJ                            | 0.014 UJ                             | 4.9 U                               | --                                  | 0.012 U                              | 0.012 U                            | 0.012 U                            | 0.1 U                               | 0.012 U                            |
| 1,1,2-Trichloroethane                                     | 1.4 U                             | 0.013 U                             | 0.014 U                              | 4.9 U                               | --                                  | 0.012 U                              | 0.012 U                            | 0.012 U                            | 0.1 U                               | 0.012 U                            |
| 1,1-Dichloroethane  | 1.4 U                             | 0.013 U                             | 0.014 U                              | 4.9 U                               | --                                  | 0.012 U                              | 0.012 U                            | 0.012 U                            | 0.1 U                               | 0.012 U                            |
| 1,1-Dichloroethene  | 1.4 U                             | 0.013 U                             | 0.014 U                              | 4.9 U                               | --                                  | 0.012 U                              | 0.012 U                            | 0.012 U                            | 0.1 U                               | 0.012 U                            |
| 1,2-Dichlorobenzene                                       | 38 U                              | 0.44 U                              | 4.5 U                                | 0.65 U                              | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 U                               | 8.1 U                              |
| 1,2-Dichloroethane  | 1.4 U                             | 0.013 U                             | 0.014 U                              | 4.9 U                               | --                                  | 0.012 U                              | 0.012 U                            | 0.012 U                            | 0.1 U                               | 0.012 U                            |
| 1,2-Dichloroethene, Total                                 | 1.4 U                             | 0.013 U                             | 0.014 U                              | 4.9 U                               | --                                  | 0.012 U                              | 0.012 U                            | 0.012 U                            | 0.1 U                               | 0.012 U                            |
| 1,2-Dichloropropane                                       | 1.4 U                             | 0.013 U                             | 0.014 U                              | 4.9 U                               | --                                  | 0.012 U                              | 0.012 U                            | 0.012 U                            | 0.1 U                               | 0.012 U                            |
| 1,3-Dichlorobenzene                                       | 38 U                              | 0.44 U                              | 4.5 U                                | 0.65 U                              | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 U                               | 8.1 U                              |
| 1,4-Dichlorobenzene                                       | 38 U                              | 0.44 U                              | 4.5 U                                | 0.65 U                              | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 U                               | 8.1 U                              |
| 2-Butanone  | 1.4 U                             | 0.013 U                             | 0.014 U                              | 4.9 U                               | --                                  | 0.012 U                              | 0.012 U                            | 0.012 U                            | 0.25                                | 0.012 U                            |
| 2-Hexanone  | 1.4 U                             | 0.013 UJ                            | 0.014 UJ                             | 4.9 U                               | --                                  | 0.012 U                              | 0.012 U                            | 0.012 U                            | 0.1 U                               | 0.012 U                            |
| 4-Methyl-2-pentanone                                      | 1.4 U                             | 0.013 UJ                            | 0.014 UJ                             | 4.9 U                               | --                                  | 0.012 U                              | 0.012 U                            | 0.012 U                            | 0.1 U                               | 0.012 U                            |
| Acetone   | 1.4 U                             | 0.013 U                             | 0.014 U                              | 4.9 U                               | --                                  | 0.012 U                              | 0.012 U                            | 0.012 U                            | 0.58                                | 0.021                              |
| Benzene   | 0.38 J                            | 0.013 U                             | 0.002 J                              | 27                                  | --                                  | 0.012 U                              | 0.012 U                            | 0.012 U                            | 8.5 D                               | 0.002 J                            |
| Bromodichloromethane                                      | 1.4 U                             | 0.013 U                             | 0.014 U                              | 4.9 U                               | --                                  | 0.012 U                              | 0.012 U                            | 0.012 U                            | 0.1 U                               | 0.012 U                            |
| Bromoform   | 1.4 U                             | 0.013 U                             | 0.014 U                              | 4.9 U                               | --                                  | 0.012 U                              | 0.012 U                            | 0.012 U                            | 0.1 U                               | 0.012 U                            |
| Bromomethane  | 1.4 U                             | 0.013 U                             | 0.014 U                              | 4.9 U                               | --                                  | 0.012 U                              | 0.012 U                            | 0.012 U                            | 0.1 U                               | 0.012 U                            |
| Carbon disulfide  | 1.4 U                             | 0.013 U                             | 0.014 U                              | 4.9 U                               | --                                  | 0.012 U                              | 0.012 U                            | 0.012 U                            | 0.1 U                               | 0.012 U                            |
| Carbon tetrachloride                                      | 1.4 U                             | 0.013 U                             | 0.014 U                              | 4.9 U                               | --                                  | 0.012 U                              | 0.012 U                            | 0.012 U                            | 0.1 U                               | 0.012 U                            |
| Chlorobenzene   | 1.4 U                             | 0.013 UJ                            | 0.014 UJ                             | 4.9 U                               | --                                  | 0.012 U                              | 0.012 U                            | 0.012 U                            | 0.1 U                               | 0.012 U                            |
| Chlorodibromomethane                                      | 1.4 U                             | 0.013 U                             | 0.014 U                              | 4.9 U                               | --                                  | 0.012 U                              | 0.012 U                            | 0.012 U                            | 0.1 U                               | 0.012 U                            |
| Chloroethane  | 1.4 U                             | 0.013 U                             | 0.014 U                              | 4.9 U                               | --                                  | 0.012 U                              | 0.012 U                            | 0.012 U                            | 0.1 U                               | 0.012 U                            |

See Notes on Page 15.

TABLE 4

**NIAGARA MOHAWK POWER CORPORATION  
ONEIDA (141 CEDAR STREET) FORMER MGP SITE  
SITE INVESTIGATION**

**SOIL ANALYTICAL RESULTS - DELINEATION SAMPLING**

| Location ID<br>Depth Range<br>Date Sampled<br>Sample Type | SB-1<br>(4 - 6')<br>7/18/00<br>FS | SB-2<br>(10 - 12')<br>7/19/00<br>FS | SB-2<br>(10 - 12')<br>7/19/00<br>DUP | SB-2<br>(19 - 23')<br>7/19/00<br>FS | SB-3<br>(12 - 14')<br>7/19/00<br>FS | SB-4A<br>(13 - 15')<br>7/20/00<br>FS | SB-6<br>(16 - 18')<br>8/8/00<br>FS | SB-6<br>(18 - 20')<br>8/8/00<br>FS | SB-8<br>(6.5 - 8')<br>8/10/00<br>FS | SB-10<br>(4 - 7')<br>8/10/00<br>FS |
|---|-----------------------------------|-------------------------------------|--------------------------------------|-------------------------------------|-------------------------------------|--------------------------------------|------------------------------------|------------------------------------|-------------------------------------|------------------------------------|
| Chloroform  | 1.4 U                             | 0.013 U                             | 0.014 U                              | 4.9 U                               | --                                  | 0.012 U                              | 0.012 U                            | 0.012 U                            | 0.1 U                               | 0.012 U                            |
| Chloromethane   | 1.4 U                             | 0.013 U                             | 0.014 U                              | 4.9 U                               | --                                  | 0.012 U                              | 0.012 U                            | 0.012 U                            | 0.1 U                               | 0.012 U                            |
| cis-1,3-Dichloropropene                                   | 1.4 U                             | 0.013 U                             | 0.014 U                              | 4.9 U                               | --                                  | 0.012 U                              | 0.012 U                            | 0.012 U                            | 0.1 U                               | 0.012 U                            |
| Ethylbenzene  | <b>0.48 J</b>                     | 0.013 UJ                            | 0.014 UJ                             | <b>18</b>                           | --                                  | 0.012 U                              | 0.012 U                            | 0.012 U                            | <b>9.2 D</b>                        | 0.012 U                            |
| Methylene chloride  | <b>0.46 J</b>                     | 0.013 U                             | 0.014 U                              | <b>0.66 J</b>                       | --                                  | 0.012 UB                             | 0.012 U                            | 0.012 U                            | 0.1 U                               | 0.012 U                            |
| Styrene   | <b>0.95 J</b>                     | 0.013 UJ                            | 0.014 UJ                             | <b>10</b>                           | --                                  | 0.012 U                              | 0.012 U                            | 0.012 U                            | 0.1 U                               | <b>0.001 J</b>                     |
| Tetrachloroethene   | 1.4 U                             | 0.013 UJ                            | 0.014 U                              | 4.9 U                               | --                                  | 0.012 U                              | 0.012 U                            | 0.012 U                            | 0.1 U                               | 0.012 U                            |
| Toluene   | <b>2.8</b>                        | <b>0.018 J</b>                      | <b>0.003 J</b>                       | <b>28</b>                           | --                                  | 0.012 U                              | 0.012 U                            | 0.012 U                            | <b>0.3</b>                          | <b>0.003 J</b>                     |
| trans-1,3-Dichloropropene                                 | 1.4 U                             | 0.013 U                             | 0.014 U                              | 4.9 U                               | --                                  | 0.012 U                              | 0.012 U                            | 0.012 U                            | 0.1 U                               | 0.012 U                            |
| Trichloroethene   | 1.4 U                             | 0.013 U                             | 0.014 U                              | 4.9 U                               | --                                  | 0.012 U                              | 0.012 U                            | 0.012 U                            | 0.1 U                               | 0.012 U                            |
| Vinyl chloride  | 1.4 U                             | 0.013 U                             | 0.014 U                              | 4.9 U                               | --                                  | 0.012 U                              | 0.012 U                            | 0.012 U                            | 0.1 U                               | 0.012 U                            |
| Xylenes, Total  | <b>15</b>                         | 0.013 UJ                            | 0.014 UJ                             | <b>97</b>                           | --                                  | 0.012 U                              | 0.012 U                            | 0.012 U                            | <b>17. D</b>                        | <b>0.004 J</b>                     |
| Total BTEX  | <b>18.66</b>                      | <b>0.018</b>                        | <b>0.005</b>                         | <b>170</b>                          | --                                  | ND                                   | ND                                 | ND                                 | <b>35</b>                           | <b>0.009</b>                       |
| <b>TCL Semivolatile Organic Compounds</b>                 |                                   |                                     |                                      |                                     |                                     |                                      |                                    |                                    |                                     |                                    |
| 1,2,4-Trichlorobenzene                                    | 38 U                              | 0.44 U                              | 4.5 U                                | 0.65 U                              | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 U                               | 8.1 U                              |
| 2,2'-oxybis(dichloropropane)                              | 38 U                              | 0.44 U                              | 4.5 U                                | 0.65 U                              | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 U                               | 8.1 U                              |
| 2,4,5-Trichlorophenol                                     | 94 U                              | 1.1 U                               | 11 U                                 | 1.6 UJ                              | 1 U                                 | 1 U                                  | 1 U                                | 1 U                                | 3.3 U                               | 20 U                               |
| 2,4,6-Trichlorophenol                                     | 38 U                              | 0.44 U                              | 4.5 U                                | 0.65 UJ                             | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 U                               | 8.1 U                              |
| 2,4-Dichlorophenol  | 38 U                              | 0.44 U                              | 4.5 U                                | 0.65 U                              | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 U                               | 8.1 U                              |
| 2,4-Dimethylphenol  | 38 U                              | 0.44 U                              | 4.5 U                                | <b>20. J</b>                        | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | <b>0.88 J</b>                       | 8.1 U                              |
| 2,4-Dinitrophenol   | 94 U                              | 1.1 U                               | 11 U                                 | 1.6 UJ                              | 1 U                                 | 1 U                                  | 1 U                                | 1 U                                | 3.3 U                               | 20 U                               |
| 2,4-Dinitrotoluene  | 38 U                              | 0.44 U                              | 4.5 U                                | 0.65 UJ                             | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 U                               | 8.1 U                              |
| 2,6-Dinitrotoluene  | 38 U                              | 0.44 U                              | 4.5 U                                | 0.65 UJ                             | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 U                               | 8.1 U                              |
| 2-Chloronaphthalene                                       | 38 U                              | 0.44 U                              | 4.5 U                                | 0.65 UJ                             | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 U                               | 8.1 U                              |
| 2-Chlorophenol  | 38 U                              | 0.44 U                              | 4.5 U                                | 0.65 U                              | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 U                               | 8.1 U                              |

See Notes on Page 15.

TABLE 4

**NIAGARA MOHAWK POWER CORPORATION  
ONEIDA (141 CEDAR STREET) FORMER MGP SITE  
SITE INVESTIGATION**

**SOIL ANALYTICAL RESULTS - DELINEATION SAMPLING**

| Location ID<br>Depth Range<br>Date Sampled<br>Sample Type | SB-1<br>(4 - 6')<br>7/18/00<br>FS | SB-2<br>(10 - 12')<br>7/19/00<br>FS | SB-2<br>(10 - 12')<br>7/19/00<br>DUP | SB-2<br>(19 - 23')<br>7/19/00<br>FS | SB-3<br>(12 - 14')<br>7/19/00<br>FS | SB-4A<br>(13 - 15')<br>7/20/00<br>FS | SB-6<br>(16 - 18')<br>8/8/00<br>FS | SB-6<br>(18 - 20')<br>8/8/00<br>FS | SB-8<br>(6.5 - 8')<br>8/10/00<br>FS | SB-10<br>(4 - 7')<br>8/10/00<br>FS |
|---|-----------------------------------|-------------------------------------|--------------------------------------|-------------------------------------|-------------------------------------|--------------------------------------|------------------------------------|------------------------------------|-------------------------------------|------------------------------------|
| 2-Methylnaphthalene                                       | 12. JD                            | 0.26 J                              | 0.74 JD                              | 110 JD                              | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 2.5                                 | 1.7 JD                             |
| 2-Methylphenol  | 38 U                              | 0.44 U                              | 4.5 U                                | 0.6 J                               | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 U                               | 8.1 U                              |
| 2-Nitroaniline  | 94 U                              | 1.1 U                               | 11 U                                 | 1.6 UJ                              | 1 U                                 | 1 U                                  | 1 U                                | 1 U                                | 3.3 U                               | 20 U                               |
| 2-Nitrophenol   | 38 U                              | 0.44 U                              | 4.5 U                                | 0.65 U                              | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 U                               | 8.1 U                              |
| 3,3'-Dichlorobenzidine                                    | 38 U                              | 0.44 U                              | 4.5 U                                | 0.65 UJ                             | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 UJ                              | 8.1 UJ                             |
| 3-Nitroaniline  | 94 U                              | 1.1 U                               | 11 U                                 | 1.6 UJ                              | 1 U                                 | 1 U                                  | 1 U                                | 1 U                                | 3.3 U                               | 20 U                               |
| 4,6-Dinitro-2-methylphenol                                | 94 U                              | 1.1 U                               | 11 U                                 | 1.6 U                               | 1 U                                 | 1 U                                  | 1 U                                | 1 U                                | 3.3 U                               | 20 U                               |
| 4-Bromophenyl phenyl ether                                | 38 U                              | 0.44 U                              | 4.5 U                                | 0.65 U                              | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 U                               | 8.1 U                              |
| 4-Chloro-3-methylphenol                                   | 38 U                              | 0.44 U                              | 4.5 U                                | 0.26 J                              | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 U                               | 8.1 U                              |
| 4-Chloroaniline   | 38 U                              | 0.44 U                              | 4.5 U                                | 0.65 U                              | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 U                               | 8.1 U                              |
| 4-Chlorophenyl phenyl ether                               | 38 U                              | 0.44 U                              | 4.5 U                                | 0.65 UJ                             | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 U                               | 8.1 U                              |
| 4-Methylphenol  | 38 U                              | 0.44 U                              | 4.5 U                                | 1.5                                 | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 0.18 J                              | 8.1 U                              |
| 4-Nitroaniline  | 94 U                              | 1.1 U                               | 11 U                                 | 1.6 UJ                              | 1 U                                 | 1 U                                  | 1 U                                | 1 U                                | 3.3 U                               | 20 U                               |
| 4-Nitrophenol   | 94 U                              | 1.1 U                               | 11 U                                 | 1.6 UJ                              | 1 U                                 | 1 U                                  | 1 U                                | 1 U                                | 3.3 U                               | 20 U                               |
| Acenaphthene  | 5.7 JD                            | 0.33 J                              | 0.63 JD                              | 24. EJ                              | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 19. JD                              | 1.0 JD                             |
| Acenaphthylene  | 24. JD                            | 0.44                                | 0.61 JD                              | 120 JD                              | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.2 J                               | 4.5 JD                             |
| Anthracene  | 39. D                             | 1.1                                 | 3.0 JD                               | 76. EJ                              | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 3.3                                 | 7.4 JD                             |
| Benzo(a)anthracene  | 120 D                             | 1.6                                 | 7.1 D                                | 55. JD                              | 0.25 J                              | 0.4 U                                | 0.4 U                              | 0.41 U                             | 3.3 J                               | 11. JD                             |
| Benzo(a)pyrene  | 76. JD                            | 1                                   | 4.2 JD                               | 40. JD                              | 0.28 J                              | 0.4 U                                | 0.4 U                              | 0.41 U                             | 2.3 J                               | 12. JD                             |
| Benzo(b)fluoranthene                                      | 98. JD                            | 1.6                                 | 7.9 JD                               | 42. JD                              | 0.32 J                              | 0.4 U                                | 0.4 U                              | 0.41 U                             | 2.9 J                               | 14. JD                             |
| Benzo(g,h,i)perylene                                      | 22. JD                            | 0.68                                | 2.2 JD                               | 9.2 EJ                              | 0.14 J                              | 0.4 U                                | 0.4 U                              | 0.41 U                             | 0.9 J                               | 5.8 JD                             |
| Benzo(k)fluoranthene                                      | 37. JD                            | 0.25 J                              | 1.8 JD                               | 5                                   | 0.14 J                              | 0.4 U                                | 0.4 U                              | 0.41 U                             | 0.94 J                              | 5.8 JD                             |
| bis(2-Chloroethoxy)methane                                | 38 U                              | 0.44 U                              | 4.5 U                                | 0.65 U                              | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 U                               | 8.1 U                              |
| bis(2-Chloroethyl)ether                                   | 38 U                              | 0.44 U                              | 4.5 U                                | 0.65 U                              | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 U                               | 8.1 U                              |
| bis(2-Ethylhexyl)phthalate                                | 38 U                              | 0.44 U                              | 4.5 U                                | 0.92 J                              | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 UJ                              | 8.1 UJ                             |

See Notes on Page 15.



TABLE 4

**NIAGARA MOHAWK POWER CORPORATION  
ONEIDA (141 CEDAR STREET) FORMER MGP SITE  
SITE INVESTIGATION**

**SOIL ANALYTICAL RESULTS - DELINEATION SAMPLING**

| Location ID<br>Depth Range<br>Date Sampled<br>Sample Type | SB-1<br>(4 - 6')<br>7/18/00<br>FS | SB-2<br>(10 - 12')<br>7/19/00<br>FS | SB-2<br>(10 - 12')<br>7/19/00<br>DUP | SB-2<br>(19 - 23')<br>7/19/00<br>FS | SB-3<br>(12 - 14')<br>7/19/00<br>FS | SB-4A<br>(13 - 15')<br>7/20/00<br>FS | SB-6<br>(16 - 18')<br>8/8/00<br>FS | SB-6<br>(18 - 20')<br>8/8/00<br>FS | SB-8<br>(6.5 - 8')<br>8/10/00<br>FS | SB-10<br>(4 - 7')<br>8/10/00<br>FS |
|---|-----------------------------------|-------------------------------------|--------------------------------------|-------------------------------------|-------------------------------------|--------------------------------------|------------------------------------|------------------------------------|-------------------------------------|------------------------------------|
| Butyl benzyl phthalate                                    | 38 U                              | 0.44 U                              | 4.5 U                                | 0.65 UJ                             | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 UJ                              | 8.1 UJ                             |
| Carbazole   | 5.0 JD                            | 0.47                                | 1.2 JD                               | 26. EJ                              | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 7.2                                 | 1.8 JD                             |
| Chrysene  | 80. D                             | 1.4                                 | 5.9 D                                | 40. JD                              | 0.2 J                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 2.7 J                               | 9.8 JD                             |
| Di-n-butyl phthalate                                      | 38 U                              | 0.44 U                              | 4.5 U                                | 0.65 U                              | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 U                               | 8.1 U                              |
| Di-n-octyl phthalate                                      | 38 UJ                             | 0.44 U                              | 4.5 UJ                               | 0.65 U                              | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 UJ                              | 8.1 UJ                             |
| Dibenz(a,h)anthracene                                     | 6.4 JD                            | 0.17 J                              | 4.5 UJ                               | 4.9                                 | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 UJ                              | 1.8 JD                             |
| Dibenzofuran  | 15. JD                            | 0.61                                | 1.6 JD                               | 79. EJ                              | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 10                                  | 3.8 JD                             |
| Diethyl phthalate   | 38 U                              | 0.44 U                              | 4.5 U                                | 0.65 UJ                             | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 U                               | 8.1 U                              |
| Dimethyl phthalate  | 38 U                              | 0.44 U                              | 4.5 U                                | 0.65 UJ                             | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 U                               | 8.1 U                              |
| Fluoranthene  | 200 D                             | 5.4 D                               | 15. D                                | 120 JD                              | 0.29 J                              | 0.046 J                              | 0.4 U                              | 0.41 U                             | 6.7                                 | 18. D                              |
| Fluorene  | 26. JD                            | 0.49                                | 0.95 JD                              | 110 EJ                              | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 9.4 JD                              | 4.5 JD                             |
| Hexachlorobenzene   | 38 U                              | 0.44 U                              | 4.5 U                                | 0.65 U                              | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 U                               | 8.1 U                              |
| Hexachlorobutadiene                                       | 38 U                              | 0.44 U                              | 4.5 U                                | 0.65 U                              | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 U                               | 8.1 U                              |
| Hexachlorocyclopentadiene                                 | 38 U                              | 0.44 U                              | 4.5 U                                | 0.65 UJ                             | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 U                               | 8.1 U                              |
| Hexachloroethane  | 38 U                              | 0.44 U                              | 4.5 U                                | 0.65 U                              | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 U                               | 8.1 U                              |
| Indeno(1,2,3-cd)pyrene                                    | 22. JD                            | 0.56                                | 2.0 JD                               | 8.7 EJ                              | 0.13 J                              | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.0 J                               | 6.5 JD                             |
| Isophorone  | 38 U                              | 0.44 U                              | 4.5 U                                | 0.65 U                              | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 U                               | 8.1 U                              |
| N-Nitroso-di-n-propylamine                                | 38 U                              | 0.44 U                              | 4.5 U                                | 0.65 U                              | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 U                               | 8.1 U                              |
| N-Nitrosodiphenylamine                                    | 38 U                              | 0.44 U                              | 4.5 U                                | 0.65 U                              | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 U                               | 8.1 U                              |
| Naphthalene   | 24. JD                            | 0.68                                | 1.4 JD                               | 610 D                               | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 62. D                               | 3.2 JD                             |
| Nitrobenzene  | 38 U                              | 0.44 U                              | 4.5 U                                | 0.65 U                              | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 U                               | 8.1 U                              |
| Pentachlorophenol   | 94 U                              | 1.1 U                               | 11 U                                 | 1.6 U                               | 1 U                                 | 1 U                                  | 1 U                                | 1 U                                | 3.3 U                               | 20 U                               |
| Phenanthrene  | 95. D                             | 6.0 D                               | 16. D                                | 230 D                               | 0.062 J                             | 0.064 J                              | 0.4 U                              | 0.41 U                             | 12. JD                              | 23. D                              |
| Phenol  | 38 U                              | 0.44 U                              | 4.5 U                                | 0.65 U                              | 0.4 U                               | 0.4 U                                | 0.4 U                              | 0.41 U                             | 1.3 U                               | 8.1 U                              |
| Pyrene  | 210 D                             | 2.8                                 | 17. D                                | 95. JD                              | 0.4                                 | 0.049 J                              | 0.4 U                              | 0.41 U                             | 8.5 J                               | 32. JD                             |
| Total PAHs  | 1097.1                            | 24.76                               | 86.43                                | 1699.8                              | 2.212                               | 0.159                                | ND                                 | ND                                 | 98.24                               | 162                                |

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TABLE 4

**NIAGARA MOHAWK POWER CORPORATION  
ONEIDA (141 CEDAR STREET) FORMER MGP SITE  
SITE INVESTIGATION**

**SOIL ANALYTICAL RESULTS - DELINEATION SAMPLING**

| Location ID<br>Depth Range<br>Date Sampled<br>Sample Type | SB-1<br>(4 - 6')<br>7/18/00<br>FS | SB-2<br>(10 - 12')<br>7/19/00<br>FS | SB-2<br>(10 - 12')<br>7/19/00<br>DUP | SB-2<br>(19 - 23')<br>7/19/00<br>FS | SB-3<br>(12 - 14')<br>7/19/00<br>FS | SB-4A<br>(13 - 15')<br>7/20/00<br>FS | SB-6<br>(16 - 18')<br>8/8/00<br>FS | SB-6<br>(18 - 20')<br>8/8/00<br>FS | SB-8<br>(6.5 - 8')<br>8/10/00<br>FS | SB-10<br>(4 - 7')<br>8/10/00<br>FS |
|---|-----------------------------------|-------------------------------------|--------------------------------------|-------------------------------------|-------------------------------------|--------------------------------------|------------------------------------|------------------------------------|-------------------------------------|------------------------------------|
| <b>Polychlorinated Biphenyls</b>                          |                                   |                                     |                                      |                                     |                                     |                                      |                                    |                                    |                                     |                                    |
| Aroclor-1016  | 3.3 UD                            | 0.33 UD                             | 0.33 UD                              | 0.33 UD                             | 0.033 U                             | 0.033 U                              | 0.04 U                             | 0.033 U                            | 0.033 U                             | 0.033 U                            |
| Aroclor-1221  | 6.6 UD                            | 0.66 UD                             | 0.66 UD                              | 0.66 UD                             | 0.066 U                             | 0.066 U                              | 0.08 U                             | 0.067 U                            | 0.066 U                             | 0.066 U                            |
| Aroclor-1232  | 3.3 UD                            | 0.33 UD                             | 0.33 UD                              | 0.33 UD                             | 0.033 U                             | 0.033 U                              | 0.04 U                             | 0.033 U                            | 0.033 U                             | 0.033 U                            |
| Aroclor-1242  | 3.3 UD                            | 0.33 UD                             | 0.33 UD                              | 0.33 UD                             | 0.033 U                             | 0.033 U                              | 0.04 U                             | 0.033 U                            | 0.033 U                             | 0.033 U                            |
| Aroclor-1248  | 3.3 UD                            | 0.33 UD                             | 0.33 UD                              | 0.33 UD                             | 0.033 U                             | 0.033 U                              | 0.04 U                             | 0.033 U                            | 0.033 U                             | 0.033 U                            |
| Aroclor-1254  | 3.3 UD                            | 0.33 UD                             | 0.33 UD                              | 0.33 UD                             | 0.033 U                             | 0.033 U                              | 0.04 U                             | 0.033 U                            | 0.033 U                             | 0.033 U                            |
| Aroclor-1260  | 3.3 UD                            | 0.33 UD                             | 0.33 UD                              | 0.33 UD                             | 0.033 U                             | 0.033 U                              | 0.04 U                             | 0.033 U                            | 0.033 U                             | 0.033 U                            |
| <b>Pesticides</b>   |                                   |                                     |                                      |                                     |                                     |                                      |                                    |                                    |                                     |                                    |
| 4,4'-DDD  | 0.33 UD                           | 0.033 UD                            | 0.033 UD                             | 0.033 UD                            | 0.0033 U                            | 0.0033 U                             | 0.004 U                            | 0.0033 U                           | 0.0033 U                            | 0.0033 U                           |
| 4,4'-DDE  | R                                 | 0.033 UD                            | 0.033 UD                             | 0.033 UJD                           | 0.0033 U                            | 0.0033 U                             | 0.004 U                            | 0.0033 U                           | 0.0033 U                            | <b>0.08 JND</b>                    |
| 4,4'-DDT  | 0.33 UD                           | 0.033 UD                            | 0.033 UD                             | 0.033 UD                            | 0.0033 U                            | 0.0033 U                             | 0.004 U                            | 0.0033 U                           | 0.0033 U                            | 0.0033 U                           |
| Aldrin  | 0.16 UD                           | 0.016 UD                            | 0.017 UD                             | 0.017 UD                            | 0.0016 U                            | 0.0016 U                             | 0.002 U                            | 0.0017 U                           | 0.0016 U                            | 0.0016 U                           |
| alpha-BHC   | 0.16 UD                           | 0.016 UD                            | 0.017 UD                             | 0.017 UD                            | 0.0016 U                            | 0.0016 U                             | 0.002 U                            | 0.0017 U                           | 0.0016 U                            | 0.0016 U                           |
| alpha-Chlordane   | 0.16 UD                           | 0.016 UD                            | 0.017 UD                             | 0.017 UD                            | 0.0016 U                            | 0.0016 U                             | 0.002 U                            | 0.0017 U                           | 0.0016 U                            | 0.0016 U                           |
| beta-BHC  | 0.16 UD                           | 0.016 UD                            | 0.017 UD                             | 0.017 UD                            | 0.0016 U                            | 0.0016 U                             | 0.002 U                            | 0.0017 U                           | 0.0016 U                            | 0.0016 U                           |
| delta-BHC   | 0.16 UD                           | 0.016 UD                            | 0.017 UD                             | 0.017 UD                            | 0.0016 U                            | 0.0016 U                             | 0.002 U                            | 0.0017 U                           | 0.0016 U                            | 0.0016 U                           |
| Dieldrin  | 0.33 UD                           | 0.033 UD                            | 0.033 UD                             | 0.033 UD                            | 0.0033 U                            | 0.0033 U                             | 0.004 U                            | 0.0033 U                           | 0.0033 U                            | 0.0033 U                           |
| Endosulfan I  | 0.16 UD                           | 0.016 UD                            | 0.017 UD                             | 0.017 UD                            | 0.0016 U                            | 0.0016 U                             | 0.002 U                            | 0.0017 U                           | 0.0016 U                            | 0.0016 U                           |
| Endosulfan II   | 0.33 UD                           | 0.033 UD                            | 0.033 UD                             | 0.033 UD                            | 0.0033 U                            | 0.0033 U                             | 0.004 U                            | 0.0033 U                           | 0.0033 U                            | 0.0033 U                           |
| Endosulfan sulfate  | 0.33 UD                           | 0.033 UD                            | 0.033 UD                             | 0.033 UD                            | 0.0033 U                            | 0.0033 U                             | 0.004 U                            | 0.0033 U                           | 0.0033 U                            | 0.0033 U                           |
| Endrin  | 0.33 UD                           | 0.033 UD                            | 0.033 UD                             | 0.033 UD                            | 0.0033 U                            | 0.0033 U                             | 0.004 U                            | 0.0033 U                           | 0.0033 U                            | 0.0033 U                           |
| Endrin aldehyde   | 0.33 UD                           | 0.033 UD                            | 0.033 UD                             | 0.033 UD                            | 0.0033 U                            | 0.0033 U                             | 0.004 U                            | 0.0033 U                           | 0.0033 U                            | 0.0033 U                           |

See Notes on Page 15.

TABLE 4

**NIAGARA MOHAWK POWER CORPORATION  
ONEIDA (141 CEDAR STREET) FORMER MGP SITE  
SITE INVESTIGATION**

**SOIL ANALYTICAL RESULTS - DELINEATION SAMPLING**

| Location ID<br>Depth Range<br>Date Sampled<br>Sample Type | SB-1<br>(4 - 6')<br>7/18/00<br>FS | SB-2<br>(10 - 12')<br>7/19/00<br>FS | SB-2<br>(10 - 12')<br>7/19/00<br>DUP | SB-2<br>(19 - 23')<br>7/19/00<br>FS | SB-3<br>(12 - 14')<br>7/19/00<br>FS | SB-4A<br>(13 - 15')<br>7/20/00<br>FS | SB-6<br>(16 - 18')<br>8/8/00<br>FS | SB-6<br>(18 - 20')<br>8/8/00<br>FS | SB-8<br>(6.5 - 8')<br>8/10/00<br>FS | SB-10<br>(4 - 7')<br>8/10/00<br>FS |
|---|-----------------------------------|-------------------------------------|--------------------------------------|-------------------------------------|-------------------------------------|--------------------------------------|------------------------------------|------------------------------------|-------------------------------------|------------------------------------|
| Endrin ketone   | 0.33 UD                           | 0.033 UD                            | 0.033 UD                             | 0.033 UD                            | 0.0033 U                            | 0.0033 U                             | 0.004 U                            | 0.0033 U                           | 0.0033 U                            | 0.23 JND                           |
| gamma-BHC (Lindane)                                       | 0.15 JD                           | 0.016 UD                            | 0.017 UD                             | 0.017 UD                            | 0.0016 U                            | 0.0016 U                             | 0.002 U                            | 0.0017 U                           | 0.0016 U                            | 0.0075 J                           |
| gamma-Chlordane   | 0.16 UD                           | 0.016 UD                            | 0.017 UD                             | 0.017 UD                            | 0.0016 U                            | 0.0016 U                             | 0.002 U                            | 0.0017 U                           | 0.0016 U                            | 0.0016 U                           |
| Heptachlor  | 0.16 UD                           | 0.016 UD                            | 0.017 UD                             | 0.017 UD                            | 0.0016 U                            | 0.0016 U                             | 0.002 U                            | 0.0017 U                           | 0.0016 U                            | 0.0016 U                           |
| Heptachlor epoxide  | R                                 | 0.016 UD                            | 0.017 UD                             | 0.061 D                             | 0.0016 U                            | 0.0016 U                             | 0.002 U                            | 0.0017 U                           | 0.0016 U                            | 0.0016 U                           |
| Methoxychlor  | 1.6 UD                            | 0.16 UD                             | 0.17 UD                              | R                                   | 0.016 U                             | 0.016 U                              | 0.02 U                             | 0.017 U                            | 0.016 U                             | 0.016 U                            |
| Toxaphene   | 16 UD                             | 1.6 UD                              | 1.7 UD                               | 1.7 UD                              | 0.16 U                              | 0.16 U                               | 0.2 U                              | 0.17 U                             | 0.16 U                              | 0.16 U                             |
| <b>TAL Inorganics</b>                                     |                                   |                                     |                                      |                                     |                                     |                                      |                                    |                                    |                                     |                                    |
| Aluminum  | 5110                              | 4790                                | 4470                                 | 8460                                | 5320                                | 5090                                 | 5120                               | 3380                               | 1210                                | 2500                               |
| Antimony  | 1.1 NUJ                           | 47.3 NJ                             | 70.5 NJ                              | 2.5 BNJ                             | 1.2 NUJ                             | 1.2 NUJ                              | 1.2 NUJ                            | 1.2 NUJ                            | 2 NUJ                               | 0.61 NUJ                           |
| Arsenic   | 3.9                               | 19.5                                | 47.3                                 | 34.7                                | 3.3                                 | 2.0 B                                | 4                                  | 1.6 B                              | 1.6 U                               | 2.5                                |
| Barium  | 20.5 B                            | 688                                 | 469                                  | 91.8                                | 24.5 B                              | 27.1 B                               | 24. B                              | 16.3 B                             | 140                                 | 12.1 B                             |
| Beryllium   | 0.52 U                            | 0.54 U                              | 0.54 U                               | 0.39 U                              | 0.29 U                              | 0.3 U                                | 0.4 B                              | 0.29 B                             | 0.4 U                               | 0.22 B                             |
| Cadmium   | 0.45 U                            | 4.9                                 | 2.6                                  | 0.79 B                              | 0.48 U                              | 0.48 U                               | 0.48 U                             | 0.48 U                             | 0.8 U                               | 0.83                               |
| Calcium   | 761 B*                            | 12700 *                             | 17300 *                              | 82700 *                             | 28600 *                             | 30100 *                              | 51400 *                            | 52700 *                            | 21600 *                             | 12000 *                            |
| Chromium  | 7.5                               | 17.7                                | 19                                   | 14.7                                | 7.5                                 | 7.9                                  | 7.6                                | 5.2                                | 2.9 B                               | 5.1                                |
| Cobalt  | 4.4 B                             | 11.5 B                              | 10.3 B                               | 8.9 B                               | 5.0 B                               | 4.7 B                                | 6.5 B                              | 4.9 B                              | 0.83 B                              | 3.0 B                              |
| Copper  | 27.5                              | 2490                                | 1020                                 | 53.4                                | 23.3                                | 15.8                                 | 21.4 *                             | 16.5 *                             | 14.1 *                              | 26.2 *                             |
| Cyanide, Total  | 0.55 U                            | 0.63 U                              | 114                                  | 14.9                                | 22.4                                | 0.51 U                               | 1.2                                | 2.7                                | 1.9 U                               | 3.6                                |
| Iron  | 12800                             | 62000                               | 79000                                | 27900                               | 11000                               | 11200                                | 13500                              | 10000                              | 18700                               | 7990                               |
| Lead  | 4.7 NJ                            | 1700 NJ                             | 14600 NJ                             | 67.2 NJ                             | 10.3 NJ                             | 6.3 NJ                               | 5.3 *                              | 4.0 *                              | 4.5 *                               | 17. *                              |
| Magnesium   | 2090 *                            | 10300 *                             | 3480 *                               | 24700 *                             | 11300 *                             | 10600 *                              | 20000 *                            | 21500 *                            | 917 B*                              | 4540 *                             |
| Manganese   | 216 *NJ                           | 3410 *NJ                            | 3220 *NJ                             | 1150 *NJ                            | 219 *NJ                             | 287 *NJ                              | 488 N                              | 377 N                              | 573 JN                              | 103 JN                             |
| Mercury   | 0.057 U                           | 0.58                                | 0.55                                 | 0.26                                | 0.068 B                             | 0.06 U                               | 0.06 U*                            | 0.16 *                             | 0.2 B*                              | 0.13 *                             |
| Nickel  | 11                                | 88.3                                | 37.4                                 | 18.6                                | 10                                  | 9.1 B                                | 12                                 | 8.6 B                              | 2.3 B                               | 11.7                               |

See Notes on Page 15.

TABLE 4

**NIAGARA MOHAWK POWER CORPORATION  
ONEIDA (141 CEDAR STREET) FORMER MGP SITE  
SITE INVESTIGATION**

**SOIL ANALYTICAL RESULTS - DELINEATION SAMPLING**

| <b>Location ID</b>  | <b>SB-1</b>     | <b>SB-2</b>       | <b>SB-2</b>       | <b>SB-2</b>       | <b>SB-3</b>       | <b>SB-4A</b>      | <b>SB-6</b>       | <b>SB-6</b>       | <b>SB-8</b>       | <b>SB-10</b>    |
|---------------------|-----------------|-------------------|-------------------|-------------------|-------------------|-------------------|-------------------|-------------------|-------------------|-----------------|
| <b>Depth Range</b>  | <b>(4 - 6')</b> | <b>(10 - 12')</b> | <b>(10 - 12')</b> | <b>(19 - 23')</b> | <b>(12 - 14')</b> | <b>(13 - 15')</b> | <b>(16 - 18')</b> | <b>(18 - 20')</b> | <b>(6.5 - 8')</b> | <b>(4 - 7')</b> |
| <b>Date Sampled</b> | <b>7/18/00</b>  | <b>7/19/00</b>    | <b>7/19/00</b>    | <b>7/19/00</b>    | <b>7/19/00</b>    | <b>7/20/00</b>    | <b>8/8/00</b>     | <b>8/8/00</b>     | <b>8/10/00</b>    | <b>8/10/00</b>  |
| <b>Sample Type</b>  | <b>FS</b>       | <b>FS</b>         | <b>DUP</b>        | <b>FS</b>         | <b>FS</b>         | <b>FS</b>         | <b>FS</b>         | <b>FS</b>         | <b>FS</b>         | <b>FS</b>       |
| Potassium           | <b>892 B</b>    | <b>470 B</b>      | <b>596 B</b>      | <b>1050 B</b>     | <b>1060 B</b>     | <b>898 B</b>      | <b>936 B</b>      | <b>709 B</b>      | 146 U             | <b>487 B</b>    |
| Selenium            | <b>1.1 B</b>    | <b>6.1</b>        | 1.1 U             | <b>3.7</b>        | <b>2.2</b>        | 0.96 U            | 0.97 U            | 0.97 U            | <b>2.6</b>        | 0.49 U          |
| Silver              | 0.45 U          | 0.53 U            | 0.54 U            | 0.78 U            | 0.48 U            | 0.48 U            | 0.48 U            | 0.48 U            | 0.8 U             | 0.24 U          |
| Sodium              | 94.2 U          | <b>409 B</b>      | <b>226 B</b>      | <b>170 B</b>      | 99.9 U            | 100 U             | <b>110 B</b>      | <b>129 B</b>      | 166 U             | <b>268 B</b>    |
| Thallium            | 1.4 U           | <b>7.5</b>        | 1.6 U             | 2.3 U             | 1.4 U             | 1.4 U             | 1.4 U             | 1.5 U             | 2.4 U             | 0.73 U          |
| Vanadium            | <b>9.7 B</b>    | <b>33.5</b>       | <b>24.9</b>       | <b>12.2 B</b>     | <b>9.7 B</b>      | <b>9.2 B</b>      | <b>10. B</b>      | <b>7.8 B</b>      | <b>2.4 B</b>      | <b>4.8 B</b>    |
| Zinc                | <b>21.1</b>     | <b>3430</b>       | <b>1450</b>       | <b>78</b>         | <b>23.9</b>       | <b>20</b>         | <b>25.7 *</b>     | <b>19.8 *</b>     | <b>10.6 *J</b>    | <b>2210 *J</b>  |

See Notes on Page 15.

TABLE 4

**NIAGARA MOHAWK POWER CORPORATION  
ONEIDA (141 CEDAR STREET) FORMER MGP SITE  
SITE INVESTIGATION**

**SOIL ANALYTICAL RESULTS - DELINEATION SAMPLING**

| Location ID<br>Depth Range<br>Date Sampled<br>Sample Type | SB-13<br>(4 - 6')<br>8/10/00<br>FS | SB-13<br>(4 - 6')<br>8/10/00<br>DUP | SB-14<br>(5 - 7')<br>8/10/00<br>FS | SB-15<br>(4 - 6')<br>8/10/00<br>FS | TP-2<br>(3')<br>12/7/00<br>FS | TP-2<br>(3')<br>12/7/00<br>DUP | TP-3<br>(6')<br>12/7/00<br>FS | TP-4<br>(3')<br>12/7/00<br>FS | TP-4<br>(6')<br>12/7/00<br>FS | TP-5<br>(5')<br>12/7/00<br>FS |
|---|------------------------------------|-------------------------------------|------------------------------------|------------------------------------|-------------------------------|--------------------------------|-------------------------------|-------------------------------|-------------------------------|-------------------------------|
| <b>TCL Volatile Organic Compounds</b>                     |                                    |                                     |                                    |                                    |                               |                                |                               |                               |                               |                               |
| 1,1,1-Trichloroethane                                     | 0.014 U                            | 0.013 U                             | 0.014 U                            | 0.014 U                            | --                            | --                             | --                            | --                            | --                            | --                            |
| 1,1,2,2-Tetrachloroethane                                 | 0.014 U                            | 0.013 U                             | 0.014 UJ                           | 0.014 U                            | --                            | --                             | --                            | --                            | --                            | --                            |
| 1,1,2-Trichloroethane                                     | 0.014 U                            | 0.013 U                             | 0.014 U                            | 0.014 U                            | --                            | --                             | --                            | --                            | --                            | --                            |
| 1,1-Dichloroethane  | 0.014 U                            | 0.013 U                             | 0.014 U                            | 0.014 U                            | --                            | --                             | --                            | --                            | --                            | --                            |
| 1,1-Dichloroethene  | 0.014 U                            | 0.013 U                             | 0.014 U                            | 0.014 U                            | --                            | --                             | --                            | --                            | --                            | --                            |
| 1,2-Dichlorobenzene                                       | 0.45 U                             | 4.2 U                               | 0.46 U                             | 0.47 U                             | --                            | --                             | --                            | --                            | --                            | --                            |
| 1,2-Dichloroethane  | 0.014 U                            | 0.013 U                             | 0.014 U                            | 0.014 U                            | --                            | --                             | --                            | --                            | --                            | --                            |
| 1,2-Dichloroethene, Total                                 | 0.014 U                            | 0.013 U                             | 0.014 U                            | 0.014 U                            | --                            | --                             | --                            | --                            | --                            | --                            |
| 1,2-Dichloropropane                                       | 0.014 U                            | 0.013 U                             | 0.014 U                            | 0.014 U                            | --                            | --                             | --                            | --                            | --                            | --                            |
| 1,3-Dichlorobenzene                                       | 0.45 U                             | 4.2 U                               | 0.46 U                             | 0.47 U                             | --                            | --                             | --                            | --                            | --                            | --                            |
| 1,4-Dichlorobenzene                                       | 0.45 U                             | 4.2 U                               | 0.46 U                             | 0.47 U                             | --                            | --                             | --                            | --                            | --                            | --                            |
| 2-Butanone  | 0.014 U                            | 0.013 U                             | <b>0.026</b>                       | 0.014 U                            | --                            | --                             | --                            | --                            | --                            | --                            |
| 2-Hexanone  | 0.014 U                            | 0.013 U                             | 0.014 UJ                           | 0.014 U                            | --                            | --                             | --                            | --                            | --                            | --                            |
| 4-Methyl-2-pentanone                                      | 0.014 U                            | 0.013 U                             | 0.014 UJ                           | 0.014 U                            | --                            | --                             | --                            | --                            | --                            | --                            |
| Acetone   | <b>0.008 J</b>                     | 0.013 U                             | <b>0.067</b>                       | <b>0.013 J</b>                     | --                            | --                             | --                            | --                            | --                            | --                            |
| Benzene   | 0.014 U                            | 0.013 U                             | <b>0.033</b>                       | 0.014 U                            | --                            | --                             | --                            | --                            | --                            | --                            |
| Bromodichloromethane                                      | 0.014 U                            | 0.013 U                             | 0.014 U                            | 0.014 U                            | --                            | --                             | --                            | --                            | --                            | --                            |
| Bromoform   | 0.014 U                            | 0.013 U                             | 0.014 U                            | 0.014 U                            | --                            | --                             | --                            | --                            | --                            | --                            |
| Bromomethane  | 0.014 U                            | 0.013 U                             | 0.014 U                            | 0.014 U                            | --                            | --                             | --                            | --                            | --                            | --                            |
| Carbon disulfide  | 0.014 U                            | 0.013 U                             | 0.014 U                            | 0.014 U                            | --                            | --                             | --                            | --                            | --                            | --                            |
| Carbon tetrachloride                                      | 0.014 U                            | 0.013 U                             | 0.014 U                            | 0.014 U                            | --                            | --                             | --                            | --                            | --                            | --                            |
| Chlorobenzene   | 0.014 U                            | 0.013 U                             | 0.014 UJ                           | 0.014 U                            | --                            | --                             | --                            | --                            | --                            | --                            |
| Chlorodibromomethane                                      | 0.014 U                            | 0.013 U                             | 0.014 U                            | 0.014 U                            | --                            | --                             | --                            | --                            | --                            | --                            |
| Chloroethane  | 0.014 U                            | 0.013 U                             | 0.014 U                            | 0.014 U                            | --                            | --                             | --                            | --                            | --                            | --                            |

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TABLE 4

**NIAGARA MOHAWK POWER CORPORATION  
ONEIDA (141 CEDAR STREET) FORMER MGP SITE  
SITE INVESTIGATION**

**SOIL ANALYTICAL RESULTS - DELINEATION SAMPLING**

| Location ID<br>Depth Range<br>Date Sampled<br>Sample Type | SB-13<br>(4 - 6')<br>8/10/00<br>FS | SB-13<br>(4 - 6')<br>8/10/00<br>DUP | SB-14<br>(5 - 7')<br>8/10/00<br>FS | SB-15<br>(4 - 6')<br>8/10/00<br>FS | TP-2<br>(3')<br>12/7/00<br>FS | TP-2<br>(3')<br>12/7/00<br>DUP | TP-3<br>(6')<br>12/7/00<br>FS | TP-4<br>(3')<br>12/7/00<br>FS | TP-4<br>(6')<br>12/7/00<br>FS | TP-5<br>(5')<br>12/7/00<br>FS |
|---|------------------------------------|-------------------------------------|------------------------------------|------------------------------------|-------------------------------|--------------------------------|-------------------------------|-------------------------------|-------------------------------|-------------------------------|
| Chloroform  | 0.014 U                            | 0.013 U                             | 0.014 U                            | 0.014 U                            | --                            | --                             | --                            | --                            | --                            | --                            |
| Chloromethane   | 0.014 U                            | 0.013 U                             | 0.014 U                            | 0.014 U                            | --                            | --                             | --                            | --                            | --                            | --                            |
| cis-1,3-Dichloropropene                                   | 0.014 U                            | 0.013 U                             | 0.014 U                            | 0.014 U                            | --                            | --                             | --                            | --                            | --                            | --                            |
| Ethylbenzene  | 0.014 U                            | 0.013 U                             | 0.019 J                            | 0.014 U                            | --                            | --                             | --                            | --                            | --                            | --                            |
| Methylene chloride  | 0.014 U                            | 0.013 U                             | 0.014 U                            | 0.014 U                            | --                            | --                             | --                            | --                            | --                            | --                            |
| Styrene   | 0.002 J                            | 0.013 U                             | 0.014 UJ                           | 0.014 U                            | --                            | --                             | --                            | --                            | --                            | --                            |
| Tetrachloroethene   | 0.014 U                            | 0.013 U                             | 0.014 UJ                           | 0.014 U                            | --                            | --                             | --                            | --                            | --                            | --                            |
| Toluene   | 0.014 U                            | 0.013 U                             | 0.004 J                            | 0.014 U                            | --                            | --                             | --                            | --                            | --                            | --                            |
| trans-1,3-Dichloropropene                                 | 0.014 U                            | 0.013 U                             | 0.014 U                            | 0.014 U                            | --                            | --                             | --                            | --                            | --                            | --                            |
| Trichloroethene   | 0.014 U                            | 0.013 U                             | 0.014 U                            | 0.014 U                            | --                            | --                             | --                            | --                            | --                            | --                            |
| Vinyl chloride  | 0.014 U                            | 0.013 U                             | 0.014 U                            | 0.014 U                            | --                            | --                             | --                            | --                            | --                            | --                            |
| Xylenes, Total  | 0.014 U                            | 0.013 U                             | 0.023 J                            | 0.014 U                            | --                            | --                             | --                            | --                            | --                            | --                            |
| Total BTEX  | ND                                 | ND                                  | 0.079                              | ND                                 | --                            | --                             | --                            | --                            | --                            | --                            |
| <b>TCL Semivolatile Organic Compounds</b>                 |                                    |                                     |                                    |                                    |                               |                                |                               |                               |                               |                               |
| 1,2,4-Trichlorobenzene                                    | 0.45 U                             | 4.2 U                               | 0.46 U                             | 0.47 U                             | --                            | --                             | --                            | --                            | --                            | --                            |
| 2,2'-oxybis(dichloropropane)                              | 0.45 U                             | 4.2 U                               | 0.46 U                             | 0.47 U                             | --                            | --                             | --                            | --                            | --                            | --                            |
| 2,4,5-Trichlorophenol                                     | 1.1 U                              | 10 U                                | 1.2 U                              | 1.2 U                              | --                            | --                             | --                            | --                            | --                            | --                            |
| 2,4,6-Trichlorophenol                                     | 0.45 U                             | 4.2 U                               | 0.46 U                             | 0.47 U                             | --                            | --                             | --                            | --                            | --                            | --                            |
| 2,4-Dichlorophenol  | 0.45 U                             | 4.2 U                               | 0.46 U                             | 0.47 U                             | --                            | --                             | --                            | --                            | --                            | --                            |
| 2,4-Dimethylphenol  | 0.45 U                             | 4.2 U                               | 0.46 U                             | 0.47 U                             | --                            | --                             | --                            | --                            | --                            | --                            |
| 2,4-Dinitrophenol   | 1.1 U                              | 10 U                                | 1.2 U                              | 1.2 U                              | --                            | --                             | --                            | --                            | --                            | --                            |
| 2,4-Dinitrotoluene  | 0.45 U                             | 4.2 U                               | 0.46 U                             | 0.47 U                             | --                            | --                             | --                            | --                            | --                            | --                            |
| 2,6-Dinitrotoluene  | 0.45 U                             | 4.2 U                               | 0.46 U                             | 0.47 U                             | --                            | --                             | --                            | --                            | --                            | --                            |
| 2-Chloronaphthalene                                       | 0.45 U                             | 4.2 U                               | 0.46 U                             | 0.47 U                             | --                            | --                             | --                            | --                            | --                            | --                            |
| 2-Chlorophenol  | 0.45 U                             | 4.2 U                               | 0.46 U                             | 0.47 U                             | --                            | --                             | --                            | --                            | --                            | --                            |

See Notes on Page 15.

TABLE 4

**NIAGARA MOHAWK POWER CORPORATION  
ONEIDA (141 CEDAR STREET) FORMER MGP SITE  
SITE INVESTIGATION**

**SOIL ANALYTICAL RESULTS - DELINEATION SAMPLING**

| Location ID                 | SB-13    | SB-13    | SB-14    | SB-15    | TP-2    | TP-2    | TP-3    | TP-4    | TP-4    | TP-5    |
|-----------------------------|----------|----------|----------|----------|---------|---------|---------|---------|---------|---------|
| Depth Range                 | (4 - 6') | (4 - 6') | (5 - 7') | (4 - 6') | (3')    | (3')    | (6')    | (3')    | (6')    | (5')    |
| Date Sampled                | 8/10/00  | 8/10/00  | 8/10/00  | 8/10/00  | 12/7/00 | 12/7/00 | 12/7/00 | 12/7/00 | 12/7/00 | 12/7/00 |
| Sample Type                 | FS       | DUP      | FS       | FS       | FS      | DUP     | FS      | FS      | FS      | FS      |
| 2-Methylnaphthalene         | 0.3 J    | 0.52 JD  | 0.46 U   | 0.16 J   | 0.43 U  | 0.44 U  | 0.52 U  | 0.46 U  | 1.2 U   | 0.47 U  |
| 2-Methylphenol              | 0.45 U   | 4.2 U    | 0.46 U   | 0.47 U   | --      | --      | --      | --      | --      | --      |
| 2-Nitroaniline              | 1.1 U    | 10 U     | 1.2 U    | 1.2 U    | --      | --      | --      | --      | --      | --      |
| 2-Nitrophenol               | 0.45 U   | 4.2 U    | 0.46 U   | 0.47 U   | --      | --      | --      | --      | --      | --      |
| 3,3'-Dichlorobenzidine      | 0.45 UJ  | 4.2 U    | 0.46 UJ  | 0.47 UJ  | --      | --      | --      | --      | --      | --      |
| 3-Nitroaniline              | 1.1 U    | 10 U     | 1.2 U    | 1.2 U    | --      | --      | --      | --      | --      | --      |
| 4,6-Dinitro-2-methylphenol  | 1.1 U    | 10 U     | 1.2 U    | 1.2 U    | --      | --      | --      | --      | --      | --      |
| 4-Bromophenyl phenyl ether  | 0.45 U   | 4.2 U    | 0.46 U   | 0.47 U   | --      | --      | --      | --      | --      | --      |
| 4-Chloro-3-methylphenol     | 0.45 U   | 4.2 U    | 0.46 U   | 0.47 U   | --      | --      | --      | --      | --      | --      |
| 4-Chloroaniline             | 0.45 U   | 4.2 U    | 0.46 U   | 0.47 U   | --      | --      | --      | --      | --      | --      |
| 4-Chlorophenyl phenyl ether | 0.45 U   | 4.2 U    | 0.46 U   | 0.47 U   | --      | --      | --      | --      | --      | --      |
| 4-Methylphenol              | 0.45 U   | 4.2 U    | 0.46 U   | 0.47 U   | --      | --      | --      | --      | --      | --      |
| 4-Nitroaniline              | 1.1 U    | 10 U     | 1.2 U    | 1.2 U    | --      | --      | --      | --      | --      | --      |
| 4-Nitrophenol               | 1.1 U    | 10 U     | 1.2 U    | 1.2 U    | --      | --      | --      | --      | --      | --      |
| Acenaphthene                | 0.21 J   | 4.2 U    | 8.7 D    | 0.26 J   | 0.43 U  | 0.44 U  | 0.19 J  | 0.052 J | 0.15 J  | 0.47 U  |
| Acenaphthylene              | 0.77     | 1.3 JD   | 0.58     | 0.048 J  | 0.32 J  | 0.3 J   | 0.066 J | 0.13 J  | 1.2 U   | 0.47 U  |
| Anthracene                  | 2        | 3.6 JD   | 6.7 D    | 0.084 J  | 0.58    | 0.53    | 0.26 J  | 0.26 J  | 1.2 U   | 0.47 U  |
| Benzo(a)anthracene          | 3.6 JD   | 6.4 D    | 3.5 J    | 0.42 J   | 5.0 JD  | 5.8 JD  | 0.74 J  | 1.2 J   | 1.2 UJ  | 0.47 U  |
| Benzo(a)pyrene              | 3.1 J    | 5.6 D    | 2.4 J    | 0.5 J    | 5.1 JD  | 5.9 JD  | 0.83 J  | 1.2 J   | R       | 0.47 UJ |
| Benzo(b)fluoranthene        | 3.5 JD   | 7.3 D    | 2.6 J    | 0.71 J   | 7.0 JD  | 7.8 JD  | 0.86 J  | 1.5 J   | R       | 0.47 UJ |
| Benzo(g,h,i)perylene        | 1.5 J    | 1.6 JD   | 1.2 J    | 0.26 J   | 3.0 J   | 2.9 J   | 0.4 J   | 0.85 J  | R       | 0.47 UJ |
| Benzo(k)fluoranthene        | 1.4 J    | 2.6 JD   | 1.1 J    | 0.23 J   | 1.4     | 1.4 J   | 0.31 J  | 0.46 J  | R       | 0.47 UJ |
| bis(2-Chloroethoxy)methane  | 0.45 U   | 4.2 U    | 0.46 U   | 0.47 U   | --      | --      | --      | --      | --      | --      |
| bis(2-Chloroethyl)ether     | 0.45 U   | 4.2 U    | 0.46 U   | 0.47 U   | --      | --      | --      | --      | --      | --      |
| bis(2-Ethylhexyl)phthalate  | 0.45 UJ  | 4.2 U    | 0.46 UJ  | 0.47 UJ  | --      | --      | --      | --      | --      | --      |

See Notes on Page 15.

TABLE 4

**NIAGARA MOHAWK POWER CORPORATION  
ONEIDA (141 CEDAR STREET) FORMER MGP SITE  
SITE INVESTIGATION**

**SOIL ANALYTICAL RESULTS - DELINEATION SAMPLING**

| Location ID                | SB-13    | SB-13    | SB-14    | SB-15    | TP-2    | TP-2    | TP-3    | TP-4    | TP-4    | TP-5    |
|----------------------------|----------|----------|----------|----------|---------|---------|---------|---------|---------|---------|
| Depth Range                | (4 - 6') | (4 - 6') | (5 - 7') | (4 - 6') | (3')    | (3')    | (6')    | (3')    | (6')    | (5')    |
| Date Sampled               | 8/10/00  | 8/10/00  | 8/10/00  | 8/10/00  | 12/7/00 | 12/7/00 | 12/7/00 | 12/7/00 | 12/7/00 | 12/7/00 |
| Sample Type                | FS       | DUP      | FS       | FS       | FS      | DUP     | FS      | FS      | FS      | FS      |
| Butyl benzyl phthalate     | 0.45 UJ  | 4.2 U    | 0.46 UJ  | 0.47 UJ  | --      | --      | --      | --      | --      | --      |
| Carbazole                  | 0.43 J   | 0.92 JD  | 2.8      | 0.23 J   | --      | --      | --      | --      | --      | --      |
| Chrysene                   | 3.2 JD   | 5.5 D    | 2.7 J    | 0.48 J   | 4.2 JD  | 5.3 JD  | 0.66 J  | 1.2 J   | 1.2 UJ  | 0.47 U  |
| Di-n-butyl phthalate       | 0.45 U   | 4.2 U    | 0.46 U   | 0.47 U   | --      | --      | --      | --      | --      | --      |
| Di-n-octyl phthalate       | 0.45 UJ  | 4.2 U    | 0.46 UJ  | 0.47 UJ  | --      | --      | --      | --      | --      | --      |
| Dibenz(a,h)anthracene      | 0.51 J   | 0.53 JD  | 0.34 J   | 0.088 J  | 0.38 J  | 0.38 J  | 0.1 J   | 0.17 J  | R       | 0.47 UJ |
| Dibenzofuran               | 0.62     | 1.1 JD   | 5.8 D    | 0.084 J  | --      | --      | --      | --      | --      | --      |
| Diethyl phthalate          | 0.45 U   | 4.2 U    | 0.46 U   | 0.47 U   | --      | --      | --      | --      | --      | --      |
| Dimethyl phthalate         | 0.45 U   | 4.2 U    | 0.46 U   | 0.47 U   | --      | --      | --      | --      | --      | --      |
| Fluoranthene               | 5.6 D    | 12. D    | 10. D    | 0.67     | 8.1 EJ  | 8.3 EJ  | 1       | 1.9     | 1.2 U   | 0.47 U  |
| Fluorene                   | 0.85     | 1.8 JD   | 7.0 D    | 0.15 J   | 0.089 J | 0.092 J | 0.22 J  | 0.071 J | 1.2 U   | 0.47 U  |
| Hexachlorobenzene          | 0.45 U   | 4.2 U    | 0.46 U   | 0.47 U   | --      | --      | --      | --      | --      | --      |
| Hexachlorobutadiene        | 0.45 U   | 4.2 U    | 0.46 U   | 0.47 U   | --      | --      | --      | --      | --      | --      |
| Hexachlorocyclopentadiene  | 0.45 U   | 4.2 U    | 0.46 U   | 0.47 U   | --      | --      | --      | --      | --      | --      |
| Hexachloroethane           | 0.45 U   | 4.2 U    | 0.46 U   | 0.47 U   | --      | --      | --      | --      | --      | --      |
| Indeno(1,2,3-cd)pyrene     | 1.7 J    | 1.7 JD   | 1.4 J    | 0.29 J   | 2.7 J   | 2.6 J   | 0.5 J   | 0.71 J  | R       | 0.47 UJ |
| Isophorone                 | 0.45 U   | 4.2 U    | 0.46 U   | 0.47 U   | --      | --      | --      | --      | --      | --      |
| N-Nitroso-di-n-propylamine | 0.45 U   | 4.2 U    | 0.46 U   | 0.47 U   | --      | --      | --      | --      | --      | --      |
| N-Nitrosodiphenylamine     | 0.45 U   | 4.2 U    | 0.46 U   | 0.47 U   | --      | --      | --      | --      | --      | --      |
| Naphthalene                | 0.3 J    | 0.51 JD  | 0.38 J   | 2.1      | 0.11 J  | 0.089 J | 0.52 U  | 0.046 J | 1.2 U   | 0.47 U  |
| Nitrobenzene               | 0.45 U   | 4.2 U    | 0.46 U   | 0.47 U   | --      | --      | --      | --      | --      | --      |
| Pentachlorophenol          | 1.1 U    | 10 U     | 1.2 U    | 1.2 U    | --      | --      | --      | --      | --      | --      |
| Phenanthrene               | 6.2 D    | 10. D    | 23. D    | 0.37 J   | 3       | 2.8     | 0.38 J  | 1.2     | 1.2 U   | 0.47 U  |
| Phenol                     | 0.45 U   | 4.2 U    | 0.46 U   | 0.47 U   | --      | --      | --      | --      | --      | --      |
| Pyrene                     | 10. JD   | 12. D    | 15. JD   | 1.1 J    | 7.9 D   | 9.7 D   | 1.4     | 2.5     | 1.2 U   | 0.47 U  |
| Total PAHs                 | 46.44    | 72.96    | 86.6     | 7.92     | 48.879  | 53.891  | 7.916   | 13.449  | NA      | ND      |

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TABLE 4

**NIAGARA MOHAWK POWER CORPORATION  
ONEIDA (141 CEDAR STREET) FORMER MGP SITE  
SITE INVESTIGATION**

**SOIL ANALYTICAL RESULTS - DELINEATION SAMPLING**

| Location ID           | SB-13    | SB-13    | SB-14    | SB-15    | TP-2    | TP-2    | TP-3    | TP-4    | TP-4    | TP-5    |
|-----------------------|----------|----------|----------|----------|---------|---------|---------|---------|---------|---------|
| Depth Range           | (4 - 6') | (4 - 6') | (5 - 7') | (4 - 6') | (3')    | (3')    | (6')    | (3')    | (6')    | (5')    |
| Date Sampled          | 8/10/00  | 8/10/00  | 8/10/00  | 8/10/00  | 12/7/00 | 12/7/00 | 12/7/00 | 12/7/00 | 12/7/00 | 12/7/00 |
| Sample Type           | FS       | DUP      | FS       | FS       | FS      | DUP     | FS      | FS      | FS      | FS      |
| Endrin ketone         | 0.033 U  | 0.0033 U | 0.0033 U | 0.0033 U | --      | --      | --      | --      | --      | --      |
| gamma-BHC (Lindane)   | 0.016 U  | 0.0017 U | 0.0017 U | 0.0017 U | --      | --      | --      | --      | --      | --      |
| gamma-Chlordane       | 0.016 U  | 0.0017 U | 0.0017 U | 0.0017 U | --      | --      | --      | --      | --      | --      |
| Heptachlor            | 0.016 U  | 0.0017 U | 0.0017 U | 0.0017 U | --      | --      | --      | --      | --      | --      |
| Heptachlor epoxide    | 0.016 U  | 0.0017 U | 0.0017 U | 0.0017 U | --      | --      | --      | --      | --      | --      |
| Methoxychlor          | 0.16 U   | R        | R        | 0.017 U  | --      | --      | --      | --      | --      | --      |
| Toxaphene             | 1.6 U    | 0.17 U   | 0.17 U   | 0.17 U   | --      | --      | --      | --      | --      | --      |
| <b>TAL Inorganics</b> |          |          |          |          |         |         |         |         |         |         |
| Aluminum              | 3150     | 3400     | 3580     | 3960     | --      | --      | --      | --      | --      | --      |
| Antimony              | 5.5 BJN  | 2.9 BJN  | 0.69 NUJ | 0.71 NUJ | --      | --      | --      | --      | --      | --      |
| Arsenic               | 6.9      | 4        | 1.5      | 2.2      | --      | --      | --      | --      | --      | --      |
| Barium                | 34.6     | 26.9     | 38.8     | 22.6 B   | --      | --      | --      | --      | --      | --      |
| Beryllium             | 0.28 B   | 0.29 B   | 0.24 B   | 0.31 B   | --      | --      | --      | --      | --      | --      |
| Cadmium               | 8.5      | 10.6     | 1.3      | 0.28 U   | --      | --      | --      | --      | --      | --      |
| Calcium               | 9440 *   | 11900 *  | 5760 *   | 5060 *   | --      | --      | --      | --      | --      | --      |
| Chromium              | 17.2     | 14.3     | 4.5      | 5.1      | --      | --      | --      | --      | --      | --      |
| Cobalt                | 3.7 B    | 3.1 B    | 3.3 B    | 3.2 B    | --      | --      | --      | --      | --      | --      |
| Copper                | 74.3 *   | 77.9 *   | 15.3 *   | 21.4 *   | --      | --      | --      | --      | --      | --      |
| Cyanide, Total        | 0.86     | 0.62 U   | 0.68 U   | 1.1      | --      | --      | --      | --      | --      | --      |
| Iron                  | 21000    | 20200    | 7550     | 8410     | --      | --      | --      | --      | --      | --      |
| Lead                  | 374 *    | 219 *    | 168 *    | 28.8 *   | --      | --      | --      | --      | --      | --      |
| Magnesium             | 4060 *   | 6280 *   | 2710 *   | 3090 *   | --      | --      | --      | --      | --      | --      |
| Manganese             | 172 JN   | 159 JN   | 106 JN   | 122 JN   | --      | --      | --      | --      | --      | --      |
| Mercury               | 0.34 *   | 0.23 *   | 1.5 *    | 0.41 *   | --      | --      | --      | --      | --      | --      |
| Nickel                | 10.4     | 8.7      | 9        | 8.6      | --      | --      | --      | --      | --      | --      |

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TABLE 4

**NIAGARA MOHAWK POWER CORPORATION  
ONEIDA (141 CEDAR STREET) FORMER MGP SITE  
SITE INVESTIGATION**

**SOIL ANALYTICAL RESULTS - DELINEATION SAMPLING**

| Location ID                      | SB-13    | SB-13    | SB-14         | SB-15    | TP-2    | TP-2    | TP-3    | TP-4    | TP-4    | TP-5    |
|----------------------------------|----------|----------|---------------|----------|---------|---------|---------|---------|---------|---------|
| Depth Range                      | (4 - 6') | (4 - 6') | (5 - 7')      | (4 - 6') | (3')    | (3')    | (6')    | (3')    | (6')    | (5')    |
| Date Sampled                     | 8/10/00  | 8/10/00  | 8/10/00       | 8/10/00  | 12/7/00 | 12/7/00 | 12/7/00 | 12/7/00 | 12/7/00 | 12/7/00 |
| Sample Type                      | FS       | DUP      | FS            | FS       | FS      | DUP     | FS      | FS      | FS      | FS      |
| <b>Polychlorinated Biphenyls</b> |          |          |               |          |         |         |         |         |         |         |
| Aroclor-1016                     | 0.33 U   | 0.033 U  | 0.033 U       | 0.033 U  | --      | --      | --      | --      | --      | --      |
| Aroclor-1221                     | 0.66 U   | 0.067 U  | 0.066 U       | 0.066 U  | --      | --      | --      | --      | --      | --      |
| Aroclor-1232                     | 0.33 U   | 0.033 U  | 0.033 U       | 0.033 U  | --      | --      | --      | --      | --      | --      |
| Aroclor-1242                     | 0.33 U   | 0.033 U  | 0.033 U       | 0.033 U  | --      | --      | --      | --      | --      | --      |
| Aroclor-1248                     | 0.33 U   | 0.033 U  | 0.033 U       | 0.033 U  | --      | --      | --      | --      | --      | --      |
| Aroclor-1254                     | 0.33 U   | 0.033 U  | 0.033 U       | 0.033 U  | --      | --      | --      | --      | --      | --      |
| Aroclor-1260                     | 0.33 U   | 0.033 U  | 0.033 U       | 0.033 U  | --      | --      | --      | --      | --      | --      |
| <b>Pesticides</b>                |          |          |               |          |         |         |         |         |         |         |
| 4,4'-DDD                         | 0.033 U  | 0.0033 U | 0.0033 U      | 0.0033 U | --      | --      | --      | --      | --      | --      |
| 4,4'-DDE                         | 0.033 U  | 0.0033 U | <b>0.0054</b> | 0.0033 U | --      | --      | --      | --      | --      | --      |
| 4,4'-DDT                         | 0.033 U  | 0.0033 U | 0.0033 U      | 0.0033 U | --      | --      | --      | --      | --      | --      |
| Aldrin                           | 0.016 U  | 0.0017 U | 0.0017 U      | 0.0017 U | --      | --      | --      | --      | --      | --      |
| alpha-BHC                        | 0.016 U  | 0.0017 U | 0.0017 U      | 0.0017 U | --      | --      | --      | --      | --      | --      |
| alpha-Chlordane                  | 0.016 U  | 0.0017 U | 0.0017 U      | 0.0017 U | --      | --      | --      | --      | --      | --      |
| beta-BHC                         | 0.016 U  | 0.0017 U | 0.0017 U      | 0.0017 U | --      | --      | --      | --      | --      | --      |
| delta-BHC                        | 0.016 U  | 0.0017 U | <b>0.0094</b> | 0.0017 U | --      | --      | --      | --      | --      | --      |
| Dieldrin                         | 0.033 U  | 0.0033 U | 0.0033 U      | 0.0033 U | --      | --      | --      | --      | --      | --      |
| Endosulfan I                     | 0.016 U  | 0.0017 U | 0.0017 U      | 0.0017 U | --      | --      | --      | --      | --      | --      |
| Endosulfan II                    | 0.033 U  | 0.0033 U | 0.0033 U      | 0.0033 U | --      | --      | --      | --      | --      | --      |
| Endosulfan sulfate               | 0.033 U  | 0.0033 U | 0.0033 U      | 0.0033 U | --      | --      | --      | --      | --      | --      |
| Endrin                           | 0.033 U  | 0.0033 U | 0.0033 U      | 0.0033 U | --      | --      | --      | --      | --      | --      |
| Endrin aldehyde                  | 0.033 U  | 0.0033 U | 0.0033 U      | 0.0033 U | --      | --      | --      | --      | --      | --      |

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TABLE 4

**NIAGARA MOHAWK POWER CORPORATION  
ONEIDA (141 CEDAR STREET) FORMER MGP SITE  
SITE INVESTIGATION**

**SOIL ANALYTICAL RESULTS - DELINEATION SAMPLING**

| <b>Location ID</b>  | <b>SB-13</b>    | <b>SB-13</b>    | <b>SB-14</b>    | <b>SB-15</b>    | <b>TP-2</b>    | <b>TP-2</b>    | <b>TP-3</b>    | <b>TP-4</b>    | <b>TP-4</b>    | <b>TP-5</b>    |
|---------------------|-----------------|-----------------|-----------------|-----------------|----------------|----------------|----------------|----------------|----------------|----------------|
| <b>Depth Range</b>  | <b>(4 - 6')</b> | <b>(4 - 6')</b> | <b>(5 - 7')</b> | <b>(4 - 6')</b> | <b>(3')</b>    | <b>(3')</b>    | <b>(6')</b>    | <b>(3')</b>    | <b>(6')</b>    | <b>(5')</b>    |
| <b>Date Sampled</b> | <b>8/10/00</b>  | <b>8/10/00</b>  | <b>8/10/00</b>  | <b>8/10/00</b>  | <b>12/7/00</b> | <b>12/7/00</b> | <b>12/7/00</b> | <b>12/7/00</b> | <b>12/7/00</b> | <b>12/7/00</b> |
| <b>Sample Type</b>  | <b>FS</b>       | <b>DUP</b>      | <b>FS</b>       | <b>FS</b>       | <b>FS</b>      | <b>DUP</b>     | <b>FS</b>      | <b>FS</b>      | <b>FS</b>      | <b>FS</b>      |
| Potassium           | 552 B           | 564 B           | 522 B           | 542 B           | --             | --             | --             | --             | --             | --             |
| Selenium            | 1.2             | 1.4             | 0.56 U          | 0.56 U          | --             | --             | --             | --             | --             | --             |
| Silver              | 0.27 U          | 0.25 U          | 0.28 U          | 0.28 U          | --             | --             | --             | --             | --             | --             |
| Sodium              | 859             | 793             | 193 B           | 159 B           | --             | --             | --             | --             | --             | --             |
| Thallium            | 0.81 U          | 0.76 U          | 0.83 U          | 0.85 U          | --             | --             | --             | --             | --             | --             |
| Vanadium            | 8.2             | 8               | 8.5             | 7.0 B           | --             | --             | --             | --             | --             | --             |
| Zinc                | 9040 *J         | 8360 *J         | 1360 *J         | 1900 *J         | --             | --             | --             | --             | --             | --             |

See Notes on Page 15.

TABLE 4

NIAGARA MOHAWK POWER CORPORATION  
ONEIDA (141 CEDAR STREET) FORMER MGP SITE  
SITE INVESTIGATION

SOIL ANALYTICAL RESULTS - DELINEATION SAMPLING

**Notes:**

1. Concentrations given in milligrams per kilogram (mg/Kg); also expressed as parts per million (ppm).
2. Samples analyzed using New York State Department of Environmental Conservation (NYSDEC) Analytical Services Protocol (ASP) 95-1, 95-2, and 95-3 Methods.
3. PAHs = Polycyclic Aromatic Hydrocarbons.
4. Carcinogenic PAHs are italicized.
5. Soil samples collected from test pits TP-2 through TP-5 were analyzed for PAHs only.
6. NA = Not available.
7. ND = Not detected.
8. -- = Not analyzed.
9. Detections are bolded.

**Data Qualifiers:**

- \* = Duplicate analysis not within control limits.
- B = (Inorganic) The reported value was obtained from a reading less than the contract required detection limit (CRDL) but greater than or equal to the instrument detection limit (IDL).
- D = Concentration is based on a diluted sample analysis.
- E = The compound was quantified above the calibration range.
- J = The compound was positively identified; however, the associated numerical value is an estimated concentration only.
- N = (Inorganic) Spiked sample recovery not within control limits.
- R = The sample results are rejected.
- U = The compound was analyzed for but not detected. The associated value is the compound quantitation limit.

TABLE 5

**NIAGARA MOHAWK POWER CORPORATION  
ONEIDA (141 CEDAR STREET) FORMER MGP SITE  
SITE INVESTIGATION**

**GROUNDWATER ANALYTICAL RESULTS**

| Constituent                                      | PZ-1 |
|--|------|
| <b>Semi-Volatile Organic Compounds (cont'd.)</b> |      |
| Nitrobenzene                                     | 10 U |
| Isophorone                                       | 10 U |
| 2-Nitrophenol                                    | 10 U |
| 2,4-Dimethylphenol                               | 10 U |
| bis(2-Chloroethoxy)methane                       | 10 U |
| 2,4-Dichlorophenol                               | 10 U |
| 1,2,4-Trichlorobenzene                           | 10 U |
| Napthalene                                       | 10 U |
| 4-Chloroaniline                                  | 10 U |
| Hexachlorobutadiene                              | 10 U |
| 4-Chloro-3-methylphenol                          | 10 U |
| 2-Methylnaphthalene                              | 10 U |
| Hexachlorocyclopentadiene                        | 10 U |
| 2,4,6-Trichlorophenol                            | 10 U |
| 2,4,5-Trichlorophenol                            | 24 U |
| 2-Chloronaphthalene                              | 10 U |
| 2-Nitroaniline                                   | 24 U |
| Dimethylphthalate                                | 10 U |
| Acenaphthylene                                   | 10 U |
| 2,6-Dinitrotoluene                               | 10 U |
| 3-Nitroaniline                                   | 24 U |
| Acenaphthene                                     | 10 U |
| 2,4-Dinitrophenol                                | 24 U |
| 4-Nitrophenol                                    | 24 U |
| Dibenzofuran                                     | 10 U |
| 2,4-Dinitrotoluene                               | 10 U |
| Diethylphthalate                                 | 10 U |
| Fluorene   | 10 U |
| 4-Chlorophenyl-phenyl ether                      | 10 U |
| 4-Nitroaniline                                   | 24 U |
| 4,6-Dinitro-2-methylphenol                       | 24 U |
| N-Nitrosodiphenylamine                           | 10 U |
| 4-Bromophenyl-phenyl ether                       | 10 U |
| Hexachlorobenzene                                | 10 U |
| Pentachlorophenol                                | 24 U |
| Phenanthrene                                     | 10 U |
| Anthracene                                       | 10 U |
| Carbazole  | 10 U |
| Di-n-butylphthalate                              | 10 U |
| Fluoranthene                                     | 10 U |
| Pyrene   | 10 U |
| Butylbenzylphthalate                             | 10 U |
| Benzo (a)anthracene                              | 10 U |
| 3,3'-Dichlorobenzidine                           | 10 U |
| Chrysene   | 10 U |
| bis (2-Ethylhexyl) phthalate                     | 10 U |

See Notes on Page 4.

TABLE 5

**NIAGARA MOHAWK POWER CORPORATION  
ONEIDA (141 CEDAR STREET) FORMER MGP SITE  
SITE INVESTIGATION**

**GROUNDWATER ANALYTICAL RESULTS**

| Constituent                            | PZ-1 |
|--|------|
| <b>Volatile Organic Compounds</b>      |      |
| Chloromethane                          | 10 U |
| Bromomethane                           | 10 U |
| Vinyl Chloride                         | 10 U |
| Chloroethane                           | 10 U |
| Methylene Chloride                     | 10 U |
| Acetone                                | 10 U |
| Carbon Disulfide                       | 10 U |
| 1,1-Dichloroethene                     | 10 U |
| 1,1-Dichloroethane                     | 10 U |
| 1,2-Dichloroethene (total)             | 10 U |
| Chloroform                             | 10 U |
| 1,2-Dichloroethane                     | 10 U |
| 2-Butanone                             | 10 U |
| 1,1,1-Trichloroethane                  | 10 U |
| Carbon Tetrachloride                   | 10 U |
| Bromodichloromethane                   | 10 U |
| 1,2-Dichloropropane                    | 10 U |
| cis-1,3-Dichloropropene                | 10 U |
| Trichloroethene                        | 10 U |
| Dibromochloromethane                   | 10 U |
| 1,1,2-Trichloroethane                  | 10 U |
| Benzene                                | 10 U |
| trans-1,3-Dichloropropene              | 10 U |
| Bromoform                              | 10 U |
| 4-Methyl-2-Pentanone                   | 10 U |
| 2-Hexanone                             | 10 U |
| Tetrachloroethene                      | 10 U |
| 1,1,2,2-Tetrachloroethane              | 10 U |
| Toluene                                | 10 U |
| Chlorobenzene                          | 10 U |
| Ethylbenzene                           | 10 U |
| Xylene (total)                         | 10 U |
| Styrene                                | 10 U |
| <b>Semi-Volatile Organic Compounds</b> |      |
| Phenol                                 | 10 U |
| bis(2-Chloroethyl) ether               | 10 U |
| 2-Chlorophenol                         | 10 U |
| 1,3-Dichlorobenzene                    | 10 U |
| 1,4-Dichlorobenzene                    | 10 U |
| 1,2-Dichlorobenzene                    | 10 U |
| 2-Methylphenol                         | 10 U |
| 2,2'-oxybis(1-Chloropropane)           | 10 U |
| 4-Methylphenol                         | 10 U |
| N-Nitroso-di-n-propylamine             | 10 U |
| Hexachloroethane                       | 10 U |

See Notes on Page 4.

TABLE 5

**NIAGARA MOHAWK POWER CORPORATION  
ONEIDA (141 CEDAR STREET) FORMER MGP SITE  
SITE INVESTIGATION**

**GROUNDWATER ANALYTICAL RESULTS**

| Constituent                                      | PZ-1   |
|--|--------|
| <b>Semi-Volatile Organic Compounds (cont'd.)</b> |        |
| Di-n-octylphthalate                              | 10 U   |
| Benzo (b) fluoranthene                           | 10 U   |
| Benzo (k) fluoranthene                           | 10 U   |
| Benzo (a) pyrene                                 | 10 U   |
| Indeno (1,2,3-cd) pyrene                         | 10 U   |
| Dibenzo(a,h)anthracene                           | 10 U   |
| Benzo(g,h,i)perylene                             | 10 U   |
| <b>PCBs</b>                                      |        |
| Aroclor-1016                                     | 1.0 U  |
| Aroclor-1221                                     | 2.0 U  |
| Aroclor-1232                                     | 1.0 U  |
| Aroclor-1242                                     | 1.0 U  |
| Aroclor-1248                                     | 1.0 U  |
| Aroclor-1254                                     | 1.0 U  |
| Aroclor-1260                                     | 1.0 U  |
| <b>Pesticides</b>                                |        |
| Aldrin   | 0.05 U |
| alpha-BHC  | 0.05 U |
| beta-BHC   | 0.05 U |
| delta-BHC  | 0.05 U |
| gamma-BHC  | 0.05 U |
| alpha-Chlordane                                  | 0.05 U |
| gamma-Chlordane                                  | 0.05 U |
| 4,4'-DDD   | 0.1 U  |
| 4,4'-DDE   | 0.1 U  |
| 4,4'-DDT   | 0.1 U  |
| Dieldrin   | 0.1 U  |
| Endosulfan I                                     | 0.05 U |
| Endosulfan II                                    | 0.1 U  |
| Endosulfan sulfate                               | 0.1 U  |
| Endrin   | 0.1 U  |
| Endrin aldehyde                                  | 0.1 U  |
| Endrin ketone                                    | 0.1 U  |
| Heptachlor                                       | 0.05 U |
| Heptachlor epoxide                               | 0.05 U |
| Methoxychlor                                     | 0.05 U |
| Toxaphene  | 0.05 U |

See Notes on Page 4.

TABLE 5

**NIAGARA MOHAWK POWER CORPORATION  
ONEIDA (141 CEDAR STREET) FORMER MGP SITE  
SITE INVESTIGATION**

**GROUNDWATER ANALYTICAL RESULTS**

| Constituent                       | PZ-1 (total) | PZ-1 (dissolved) |
|-----------------------------------|--------------|------------------|
| <b>TAL Inorganic Constituents</b> |              |                  |
| Aluminum                          | 210          | 100 U            |
| Antimony                          | 10 U         | 10 U             |
| Arsenic                           | 10 U         | 10 U             |
| Barium                            | 36           | 32               |
| Beryllium                         | 5 U          | 5 U              |
| Cadmium                           | 5 U          | 5 U              |
| Calcium                           | 450000       | 440000           |
| Chromium                          | 10 U         | 10 U             |
| Cobalt                            | 10 U         | 10 U             |
| Copper                            | 10 U         | 10 U             |
| Iron                              | 1400         | 790              |
| Lead                              | 9.3          | 3 U              |
| Magnesium                         | 21000        | 20000            |
| Manganese                         | 2000         | 1800             |
| Mercury                           | 0.2 U        | 0.2 U            |
| Nickel                            | 20 U         | 20 U             |
| Potassium                         | 10000        | 12000            |
| Selenium                          | 5 U          | 19               |
| Silver                            | 10 U         | 10 U             |
| Sodium                            | 5600         | 5300             |
| Thallium                          | 10 U         | 10 U             |
| Vanadium                          | 10 U         | 10 U             |
| Zinc                              | 1300         | 1100             |
| Cyanide, total                    | 10 U         | NA               |

**Notes:**

1. All concentrations are reported in micrograms per liter (ug/L), or parts per billion (ppb).
2. Samples analyzed using New York State Department of Environmental Conservation (NYSDEC) Analytical Services Protocol (ASP) 95-1, 95-2, and 95-3 Methods.
3. U = Compound was not detected at a concentration exceeding the laboratory detection limit.
4. NA = Not analyzed
5. The analytical laboratory results were not validated.



TABLE 6

**NIAGARA MOHAWK POWER CORPORATION  
ONEIDA (141 CEDAR STREET) FORMER MGP SITE  
SITE INVESTIGATION**

**SUMMARY OF WATER LEVEL MEASUREMENTS**

| Location ID | Measuring<br>Point<br>Elevation | Ground<br>Surface<br>Elevation | Total<br>Depth | Water Level Measurements<br>8/24/00 |                          |
|-------------|---------------------------------|--------------------------------|----------------|-------------------------------------|--------------------------|
|             |                                 |                                |                | Depth to<br>Water                   | Water Level<br>Elevation |
|             | ft. AMSL                        | ft. AMSL                       | ft. MP         | ft. MP                              | ft. AMSL                 |
| 129-MW-1    | 434.87                          | 435.3                          | 17.2           | 10.34                               | 424.5                    |
| 129-MW-2    | 424.79                          | 422.6                          | 16.2           | 6.94                                | 417.9                    |
| 129-MW-3    | 422.51                          | 420.4                          | 16.3           | 6.30                                | 416.2                    |
| 129-MW-4    | 425.99                          | 423.7                          | 15.4           | 7.19                                | 418.8                    |
| 153-MW-1    | 420.94                          | 419.3                          | 15.2           | 3.71                                | 417.2                    |
| 153-MW-2    | 435.13                          | 435.6                          | 17.3           | 8.51                                | 426.6                    |
| PZ-1        | 425.33                          | 423.1                          | 10.1           | 4.78                                | 420.6                    |

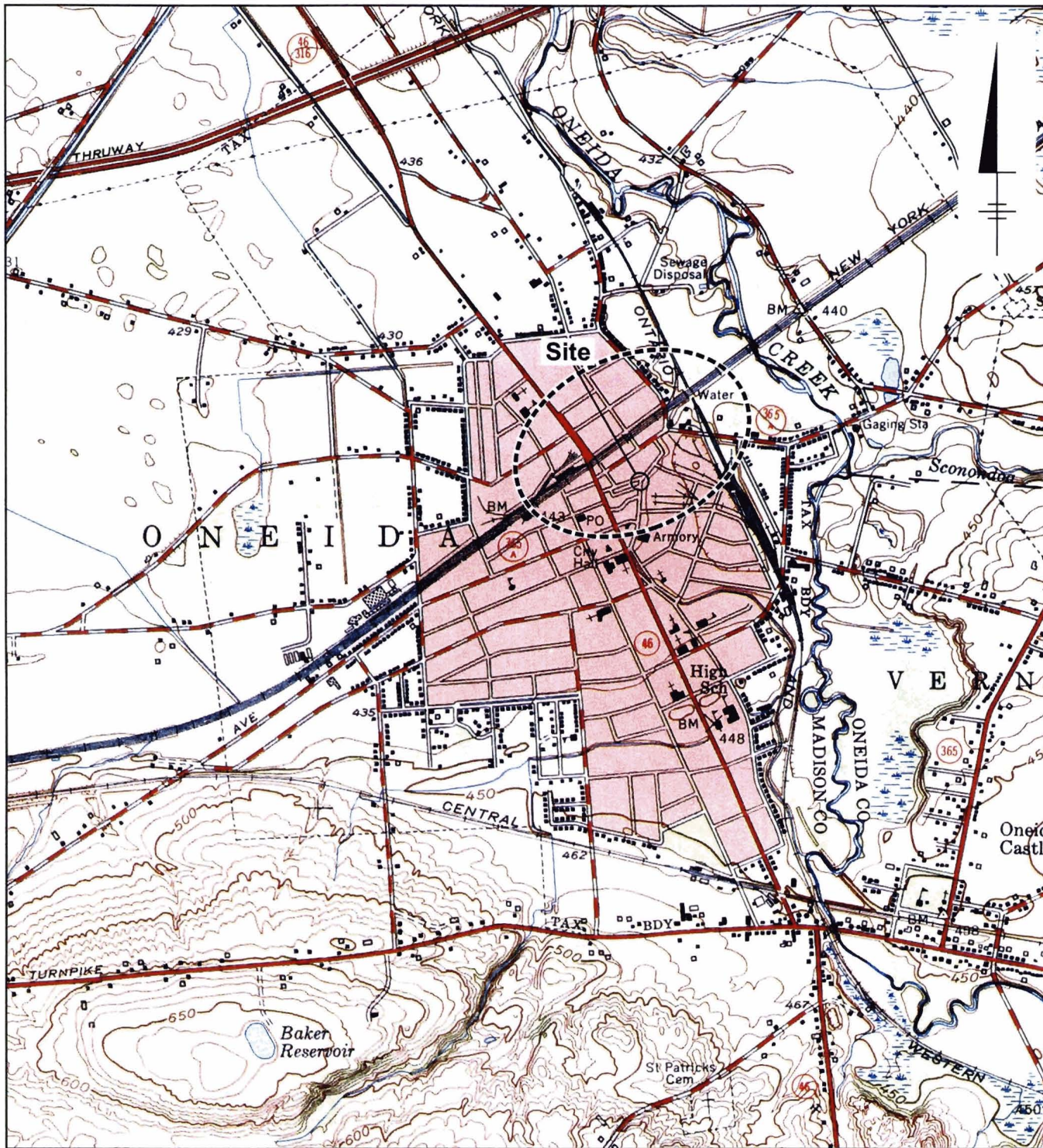
**Notes:**

1. ft. AMSL = feet Above Mean Sea Level.
2. ft. MP = feet below the surveyed measuring point.

## ***Figures***

BLASLAND, BOUCK & LEE, INC. *engineers & scientists*

*consultants with focus*



REFERENCE: BASE MAP SOURCE USGS 7.5 MINUTE QUADS. SERIES ONEIDA, NEW YORK, 1955.

2000' 0 2000'  
Approximate Scale: 1" = 2000'



03/01 SYR-D54-LBR DJH MRC  
36456007/36456n01.cdr

NIAGARA MOHAWK POWER CORPORATION  
ONEIDA (141 CEDAR STREET) FORMER MGP SITE  
SITE INVESTIGATION REPORT

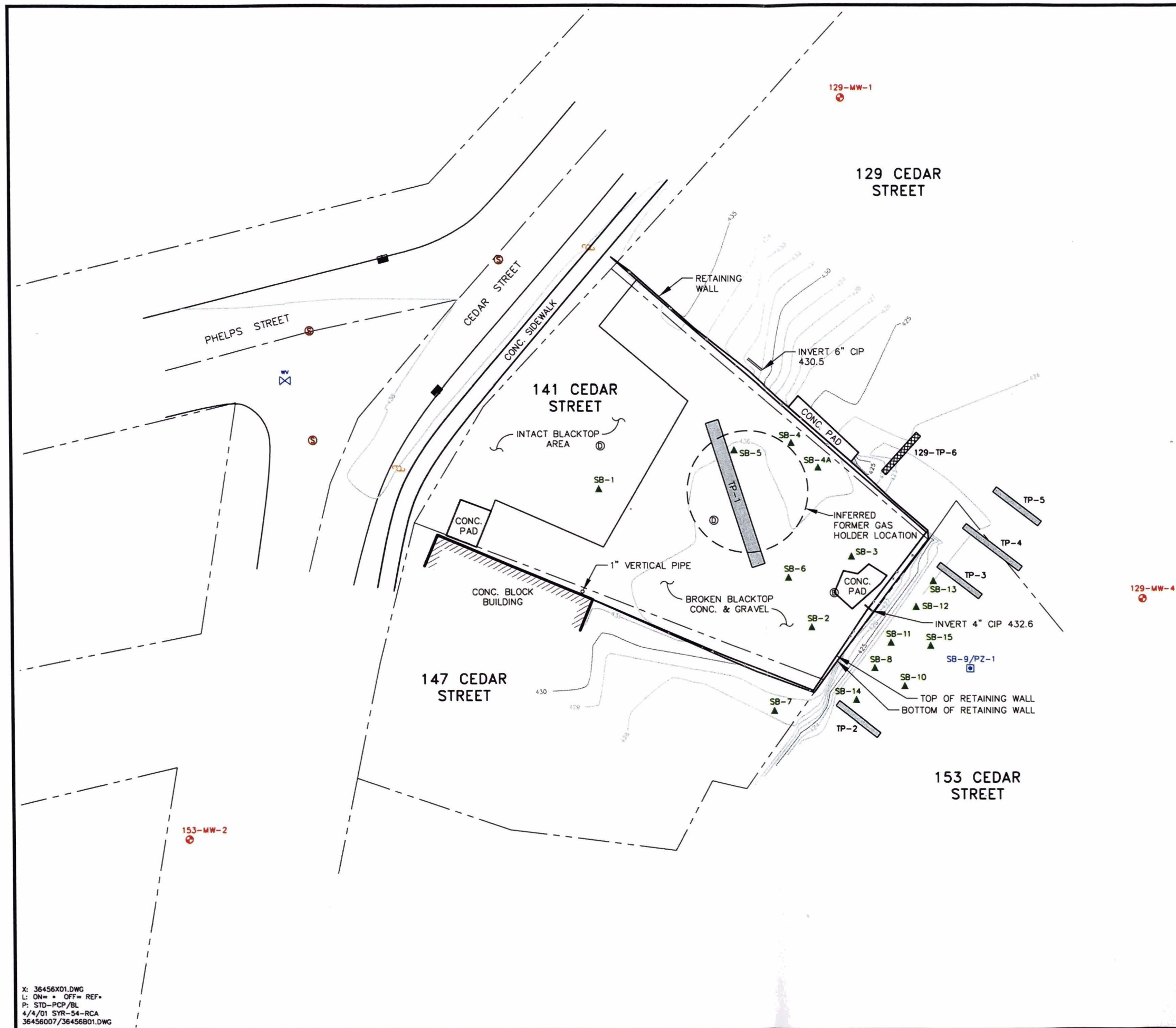
## SITE LOCATION MAP

**BBL**

BLASLAND, BOUCK & LEE, INC.  
engineers & scientists

FIGURE  
1





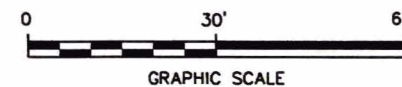
- LEGEND**
- SB-13 ▲ SOIL BORING
  - PZ-1 □ TEMPORARY MONITORING WELL
  - TP-1 ▭ TEST PIT
  - 129-TP-6 ▨ TEST PIT INSTALLED BY HARZA
  - 129-MW-1 ● MONITORING WELL (INSTALLED BY HARZA)
  - UTILITY POLE
  - ⊙ SANITARY MANHOLE
  - ⊞ DROP INLET
  - ⊕ CATCH BASIN
  - ⊗ WATER VALVE
  - CIP CAST IRON PIPE
  - - - APPROXIMATE PROPERTY LINE (SEE NOTE 3)

**NOTES:**

1. NORTH ARROW INDICATES MAGNETIC NORTH AS OBSERVED ON JULY 7, 2000.
2. VERTICAL DATUM IS REFERENCED TO NGVD 1929. HORIZONTAL DATUM IS ASSUMED.
3. NO BOUNDARY SURVEY WAS CONDUCTED, ANY PROPERTY LINES SHOWN ARE DIGITIZED FROM A PHOTOCOPY OF MADISON COUNTY TAX MAP, DATED 1975, AT AN APPROXIMATE SCALE OF 1"= 100', AND ARE APPROXIMATE ONLY.
4. MONITORING WELLS SHOWN ON THIS FIGURE ARE FROM THE FEBRUARY, 2000 SITE INVESTIGATION AND REMEDIAL ACTION REPORT (SI/RAR) FOR THE 129 CEDAR STREET PARCEL (e.g. 129-MW-2) AND THE MAY 2000 DRAFT SI/RAR FOR THE 153 CEDAR STREET PARCEL (e.g. 153-MW-1), COMPLETED BY HARZA ENGINEERING COMPANY.
5. PROPERTY NUMBER PREFIX (i.e. 129- AND 1153-) FOR MONITORING WELL INSTALLED BY HARZA WERE ATTACHED HERE FOR CLARITY.

**REFERENCE DRAWINGS:**

1. MAP SHOWING SITE PLAN - 129 CEDAR STREET, BROWNFIELD INVESTIGATION CITY OF ONEIDA, ONEIDA COUNTY, NEW YORK. DATED 11/15/99 BY HARZA NORTHEAST.



NIAGARA MOHAWK POWER CORPORATION  
ONEIDA (141 CEDAR STREET) FORMER MGP SITE  
**SITE INVESTIGATION REPORT**

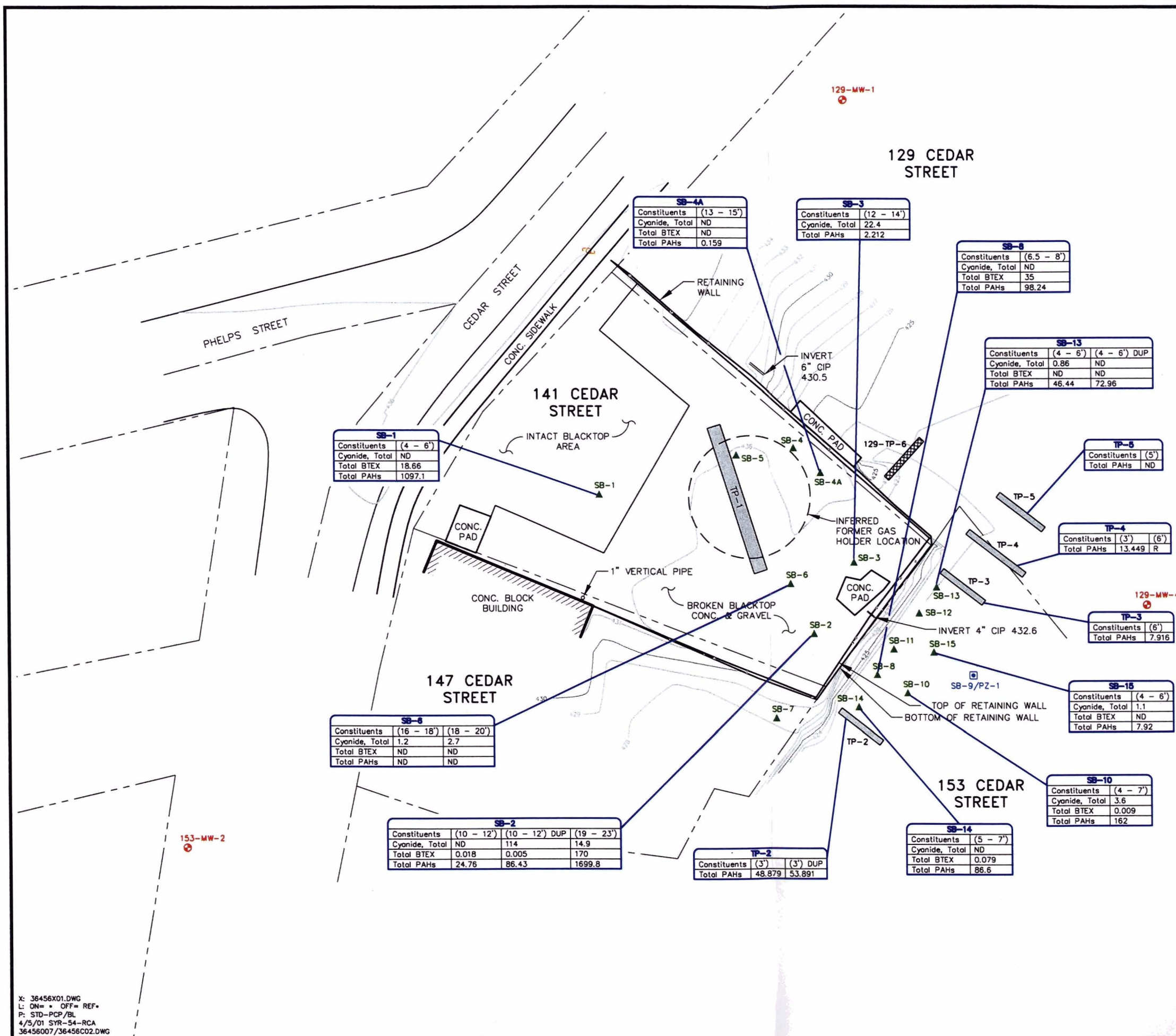
**SITE MAP**

**BBL** BLASLAND, BOUCK & LEE, INC.  
engineers & scientists

FIGURE  
**2**

X: 36456X01.DWG  
L: ON= \* OFF= REF\*  
P: STD-PCP/BL  
4/4/01 SYR-54-RCA  
36456007/36456B01.DWG





**LEGEND**

SB-13 ▲ SOIL BORING  
PZ-1 □ TEMPORARY MONITORING WELL  
TP-1 ▭ TEST PIT  
129-TP-6 ▨ TEST PIT INSTALLED BY HARZA  
129-MW-1 ● MONITORING WELL (INSTALLED BY HARZA)  
○ UTILITY POLE  
CIP CAST IRON PIPE  
--- APPROXIMATE PROPERTY LINE (SEE NOTE 3)

**SAMPLE ID**  
SAMPLE DEPTH, IN FEET BELOW GROUND SURFACE  
DUP = DUPLICATE ANALYSIS  
ND = NOT DETECTED AT OR ABOVE THE QUANTITATION LIMIT  
R = TOTAL BENZENE, TOLUENE, ETHYLBENZENE, XYLENES  
R = TOTAL POLYCYCLIC AROMATIC HYDROCARBONS  
R = ONE OR MORE OF THE RESULTS FOR THE PAH COMPOUNDS WAS REJECTED

ANALYTICAL RESULTS GIVEN IN MILLIGRAMS PER KILOGRAM (mg/Kg), OR PARTS PER MILLION (ppm)

- NOTES:**
1. NORTH ARROW INDICATES MAGNETIC NORTH AS OBSERVED ON JULY 7, 2000.
  2. VERTICAL DATUM IS REFERENCED TO NGVD 1929. HORIZONTAL DATUM IS ASSUMED.
  3. NO BOUNDARY SURVEY WAS CONDUCTED, ANY PROPERTY LINES SHOWN ARE DIGITIZED FROM A PHOTOCOPY OF MADISON COUNTY TAX MAP, DATED 1975, AT AN APPROXIMATE SCALE OF 1"= 100', AND ARE APPROXIMATE ONLY.
  4. MONITORING WELLS SHOWN ON THIS FIGURE ARE FROM THE FEBRUARY, 2000 SITE INVESTIGATION AND REMEDIAL ACTION REPORT (SI/RAR) FOR THE 129 CEDAR STREET PARCEL (e.g. 129-MW-2) AND THE MAY 2000 DRAFT SI/RAR FOR THE 153 CEDAR STREET PARCEL (e.g. 153-MW-1), COMPLETED BY HARZA ENGINEERING COMPANY.
- REFERENCE DRAWINGS:**
1. MAP SHOWING SITE PLAN - 129 CEDAR STREET, BROWNFIELD INVESTIGATION CITY OF ONEIDA, ONEIDA COUNTY, NEW YORK. DATED 11/15/99 BY HARZA NORTHEAST.

0 30' 60'  
GRAPHIC SCALE

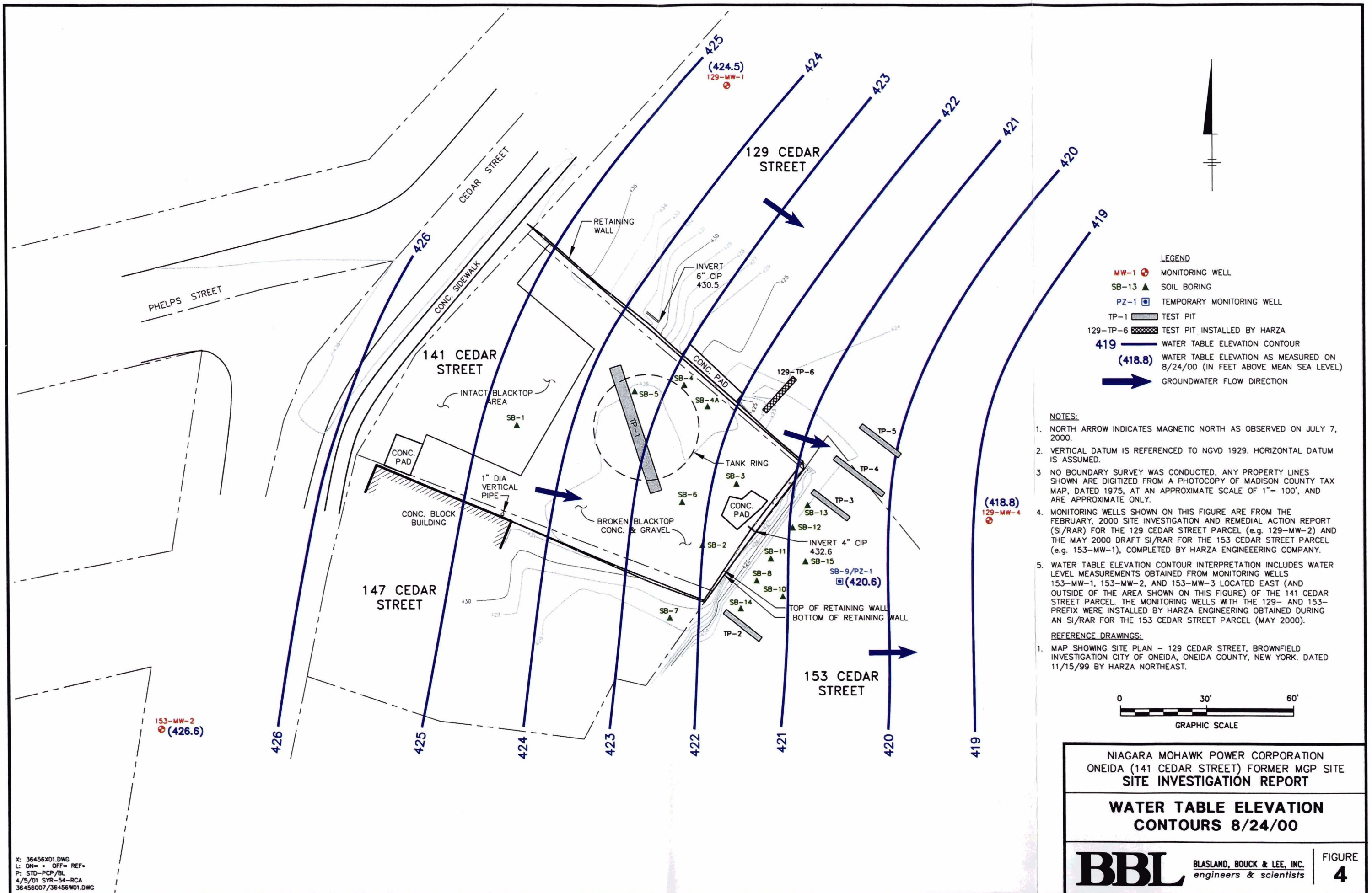
NIAGARA MOHAWK POWER CORPORATION  
ONEIDA (141 CEDAR STREET) FORMER MGP SITE  
**SITE INVESTIGATION REPORT**

**SUMMARY OF TOTAL BTEX, PAHs,  
AND CYANIDE DETECTED IN  
SUBSURFACE SOILS**

**BBL** BLASLAND, BOUCK & LEE, INC.  
engineers & scientists

FIGURE  
**3**





## ***Attachments***

BLASLAND, BOUCK & LEE, INC. *engineers & scientists*

*consultants with focus*

## **Attachment 1**

**July 14, 2000 and November 3, 2000  
Letters from NMPC to NYSDEC**





Steven P. Stucker  
Environmental Analyst

Phone: 315-428-5652  
FAX: 315-460-9670  
E-mail: stuckers@nimo.com

July 14, 2000

Mr. John Helmeset, P.E.  
Bureau of Western Remedial Action  
Division of Environmental Remediation  
New York State Department of Environmental Conservation  
50 Wolf Road  
Albany, New York 12233-7010

Re: Additional Investigations  
Niagara Mohawk Power Corporation  
Oneida (141 Cedar Street) Former MGP Site

Dear Mr. Helmeset:

This letter documents the agreed-upon additional investigation activities to be completed at the Oneida (141 Cedar Street) Former Manufactured Gas Plant (MGP) Site. These additional investigations were agreed to during a conference call between you, Cathy Geraci [Blasland, Bouck & Lee, Inc. (BBL)], George Thomas (BBL), and me on July 24, 2000 and are beyond the scope of work presented in the Supplemental Investigation Work Plan, dated May 24, 2000. The table below summarizes the proposed additional investigation activities and documents the rationale for these activities. A hand-sketched figure of the site and locations of the soil borings installed during the recent investigation activities is attached. This figure was previously provided to you to facilitate the July 24, 2000 conference call.

| Location   | Rationale  |
|--|--|
| <b>SB-6</b><br>Within extent of 141 Cedar Street retaining walls, outside of the former gas holder, located between SB-2 and SB-5. | Assess horizontal extent of presumed MGP-impacted materials observed in SB-2 at 16.5 feet below grade (bg), assess depth to native materials, and provide another boring near the former gas holder. |
| <b>SB-7</b><br>Approximately 10 feet south of the southern retaining wall at 141 Cedar Street, south of SB-2.                      | Assess horizontal extent of presumed MGP-impacted materials observed in SB-2 at 16.5 feet bg and assess depth to native materials.   |
| <b>SB-8</b><br>Approximately 10 feet east of the 141 Cedar Street eastern retaining wall, east of SB-2.                            | Assess horizontal extent of presumed MGP-impacted materials observed in SB-2 at 16.5 feet bg and assess depth to native materials.   |

| Location  | Rationale   |
|---|---|
| <b>SB-9</b><br>Approximately 30 feet to 40 feet east of the eastern retaining wall at 141 Cedar Street. | Provide a soil boring at a presumed downgradient location from 141 Cedar Street. A temporary well will be installed at this location and sampled for Target Compound List/Target Analyte List (TCL/TAL) constituents with a <input type="checkbox"/> results only <input type="checkbox"/> data package if MGP-impacted materials are not identified. Ground-water sampling will follow rudimentary development with a disposable bailer. If feasible, a depth to water and a relative elevation of the water table will also be determined. In addition, a round of water levels from the existing monitoring wells on the 129 and 153 Cedar Street properties will be obtained to further assess the ground-water flow patterns near the 141 Cedar Street property. |

While not discussed in detail during the conference call, NMPC proposes to collect additional soil samples from borings SB-6 through SB-8 for environmental characterization. Approximately one soil sample from each boring will be collected and analyzed for TCL/TAL constituents. NMPC does not propose to collect a soil sample from the SB-9 boring for laboratory analytical analysis because a ground-water sample will be collected from the temporary well installed at this location. If evidence of MGP-impacted materials are identified in any of these borings, additional investigation activities (e.g., borings) will be completed to assess the presence or absence of MGP-impacted materials, as necessary and in accordance with the Voluntary Cleanup Agreement. These additional investigation activities, if any, will be determined with the New York State Department of Environmental Conservation and City of Oneida.

As discussed during the conference call, the soil borings and temporary well installation are proposed to be conducted in the near future, in conjunction with the investigation activities currently underway at the Oneida Sconodoo Street Former MGP site. Once a firm start date is determined for the recommencement of work at the 141 Cedar Street site, NMPC will contact the NYSDEC regarding the proposed schedule.

Please feel free to contact me if you have any questions or require further assistance.

Sincerely,

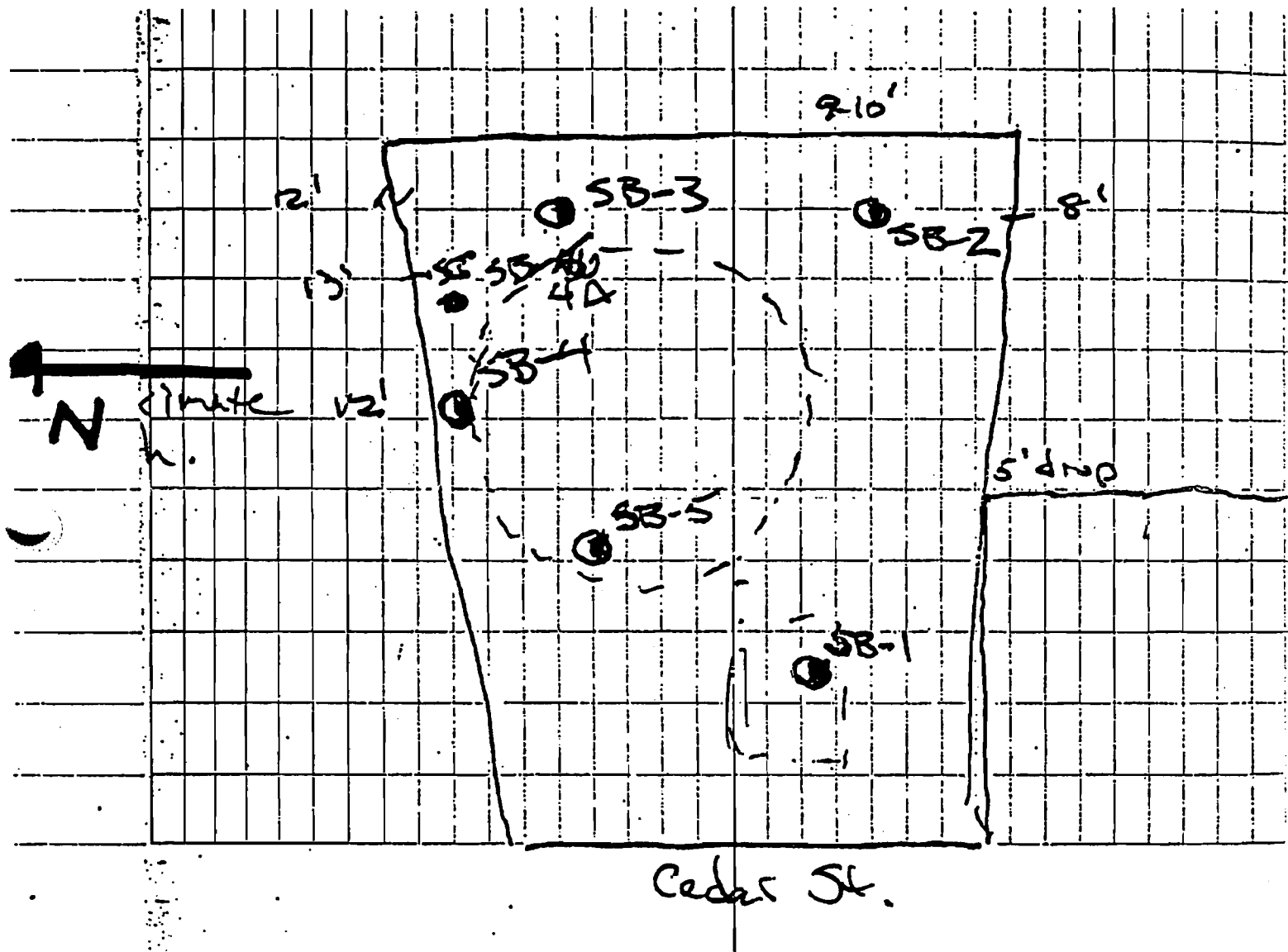


Steven P. Stucker, C.P.G.

Attachment

cc: James W. Bacher, P.E.-City of Oneida  
Michael W. Sherman-Niagara Mohawk Power Corporation  
William C. Weiss-Niagara Mohawk Power Corporation  
M. Cathy Geraci-Blasland, Bouck & Lee, Inc.  
~~George M. Thomas~~-Blasland, Bouck & Lee, Inc.

Oneida (141 Cedar Street) Former MGP Site  
"Working Hand-Sketched Figure"  
NOT-TO-SCALE





*Transmitted Via Facsimile/U.S. Mail*

July 14, 2000

Mr. John A. Helmeset, P.E.  
Environmental Engineer 2  
MGP Section  
Bureau of Western Remedial Action  
Division of Environmental Remediation  
New York State Department of Environmental Conservation  
50 Wolf Road  
Albany, NY 12233-7010

Re: Site Investigation Work Plan  
Niagara Mohawk Power Corporation  
Oneida (141 Cedar Street) Former MGP Site  
Oneida, New York

Dear Mr. Helmeset:

Thank you for your review of the Site Investigation (SI) Work Plan for the Oneida (141 Cedar Street) Former Manufactured Gas Plant (MGP) Site in Oneida, New York. Presented below are Niagara Mohawk Power Corporation's (NMPC's) responses to the New York State Department of Environmental Conservation's (NYSDEC's) comments regarding the SI Work Plan that were presented in your July 7, 2000 letter to NMPC. For ease of presentation, the following presents the NYSDEC's comments followed by our response.

*Comment: Our Department also expects that if contamination is identified at the site that appears to extend beyond the area encompassed by the retaining walls at the site, that additional borings will be completed on the adjoining City owned properties (129 and 153 Cedar Street) as part of this field mobilization to assess the presence or absence of this contamination on those properties.*

**Response:** If evidence of MGP-impacted materials are identified within the area encompassed by the retaining walls at the site, additional investigation activities (e.g., borings) will be completed on the City owned adjoining properties to assess the presence or absence of MGP-impacted materials on those properties, as necessary and in accordance with the Voluntary Cleanup Agreement. These additional investigation activities, if any, will be determined with the NYSDEC and City of Oneida and completed during this field event, if possible, in conjunction with the pre-design characterization activities at the Sconondoa Street Former MGP Site. Drilling and sampling procedures will be conducted in accordance with the Quality Assurance Project Plan (QAPP) and Field Sampling Plan (FSP) provided in the Preliminary Remedial Design Work Plan for the Sconondoa Street Former MGP Site (BBL, May 2000). As stated on the second page of your letter, you requested that if additional site investigation and characterization work is required, it should be conducted in accordance with the procedures in the Remedial Investigation Work Plan for Sconondoa Street. NMPC proposes to use the procedures provided in the Preliminary Remedial Design Work Plan because these procedures are generally consistent

with NYSDEC-approved procedures used for other NMPC sites (e.g., Amsterdam and Watertown).

NMPC understands the NYSDEC's right to require a full investigation as identified in the Voluntary Cleanup Agreement, if warranted based upon site conditions. As detailed in the SI Work Plan, the SI is intended to supply the information needed to support an Interim Remedial Measure (IRM) at the site, anticipated to consist of the excavation and off-site treatment/disposal of approximately 2,500 cubic yards of fill materials.

*Comment: Please add the completion of historical research to the breakdown of work plan tasks for Cedar Street. The general site history contained in this submittal is insufficient to determine the needed level of effort to fully characterize the site. This information includes the SI/RAR report for the adjoining 129 and 153 Cedar Street properties. These reports contain valuable information such as a complete set of analytical data and stratigraphy from the monitoring wells.*

**Response:** As documented in an e-mail from James Bacher, P.E. of the City of Oneida to yourself on July 12, 2000, the SI/RAR report for 153 Cedar Street has not yet been finalized. NMPC will coordinate with the City of Oneida to obtain this report. Pertinent historical information from the SI/RAR reports for the adjoining 129 and 153 Cedar Street properties will be incorporated in NMPC's letter report to the NYSDEC summarizing the SI field activities and analytical results.

*Comment: Wells will need to be installed directly down gradient of the site and any identified areas of contamination to quantify or confirm the lack of groundwater contamination associated with the site. None of the existing wells appears to be directly down gradient of the site and they all appear to be of an extremely shallow depth. Additionally, the provided groundwater contour map appears inaccurate. We strongly suggest that this work be completed as part of the current site work.*

**Response:** As discussed during the June 30, 2000 project meeting at the site, monitoring wells may be installed depending upon the findings of the SI activities. If MGP-impacted materials are not identified during the SI activities, NMPC believes there is no need to install monitoring wells either on-site or downgradient of the site.

NMPC agrees that the existing monitoring wells on the adjacent property do not appear to be directly downgradient from the 141 Cedar Street site, however the analytical results of the investigation activities conducted at 129 Cedar Street do provide relevant information as you indicated on the second page of your letter. The monitoring wells installed as part of the SI activities for 129 Cedar Street are screened from 3 feet to 14 feet below ground surface (bgs), 4 feet to 14 feet bgs, or 8 feet to 18 feet bgs. The ground surface elevation of these monitoring wells are below the surface elevation of the 141 Cedar Street site and some are below the elevation of the toe of the retaining wall. Monitoring wells, if necessary, will be installed in accordance with the FSP provided in the Preliminary Remedial Design Work Plan for the Sconodoa Street Former MGP Site.

*Comment: Additional samples will need to be collected to characterize the soils underlying those bound by the retaining walls. We suggest that a minimum of five additional samples be collected at the currently proposed boring locations. Four at the interface of the soils bound by the retaining wall and those underlying the toe of the retaining wall, and one at the top of the*

*peat layer. This will require the advancement of the proposed borings through the retaining wall soils, deeper than the proposed 15 feet. The depth of advancement of all of the borings beneath the retaining walls should be to the peat layer identified during the 129 Cedar Street brownfield investigation. This depth varies, but should only be approximately 25 feet on the east end of the site which has the greatest amount of fill. We also believe that a test pit should be utilized to characterize the holder contents and integrity instead of a boring.*

*We also request that the boring outside the holder be placed on the north side of the former holder, and that at least one of the remaining two borings be placed adjacent to the eastern retaining wall.*

**Response:** As detailed in the SI Work Plan, discrete soil samples were proposed by NMPC to be collected at the bottom of each of the five soil borings to determine the soil quality at the base of the soil expected to be excavated during the IRM. The intent of this proposed sampling is to characterize the material below the soil expected to be excavated during the IRM (i.e., below the retaining wall). Accordingly and consistent with the NYSDEC's comment, soil samples will be collected below the soils bound by the retaining wall and those underlying the toe of the retaining wall from four of the five soil borings to be installed.

Each of the five soil borings from which these samples are collected will be advanced to the native material (e.g., peat, silt) or to a maximum depth of approximately 30 feet bgs, whichever is shallower. A soil sample will be collected from the fifth soil boring at the top of the native material. Additional soil samples may be collected from the four other borings from the top of the native material based on photoionization detection (PID) readings, odors, staining, or non-aqueous phase liquid (NAPL) observed in the soil, if any. Furthermore, the borings may be advanced further if evidence of MGP-impacted material is identified at the base of the boring as defined above. Borings, however, will not penetrate a confining unit if NAPL or significant indications of NAPL are observed. Also as requested in your July 7, 2000 comment letter, one soil boring will be installed outside and north of the former gas holder and a second boring will be installed adjacent to the eastern retaining wall, as well as within the approximate center of the former gas holder (as detailed in the SI Work Plan).

In addition to the soil boring to be installed in the approximate center of the former gas holder, a test pit will be conducted within the location of the former gas holder to facilitate determining the contents and integrity of the former gas holder, if present. The test pit is anticipated to be at least 6 feet in length and excavated until either significant source materials are encountered, or to within the physical limits of the excavation equipment (backhoe). The installation of the test pit (depth and location) may be limited due to a storm water grate (system) observed to be located in the vicinity of the former gas holder. The test pit will be visually characterized with depth, as well as photographed for future reference. Following visual description, the test pit will be backfilled with the material removed from this test pit. Drilling, test pit excavation, and sampling procedures will be conducted in accordance with the QAPP and FSP for the Sconodoo Street Former MGP Site.

**Comment:** *The soil sample analytical parameters for the samples collected from beneath the soils bounded by the retaining walls should be modified to the full Target Compound and Target Analyte List (TCL/TAL) plus cyanide.*

**Response:** Each of the five discrete soil samples to be collected from beneath the soils bounded by the retaining walls will be analyzed for full TCL/TAL and total cyanide.

**Comment:** *To properly characterize the soil on the site for disposal, they should also be tested for the other Resource Conservation and Recovery Act (RCRA) parameters, i.e. ignitability, and the toxicity characteristic leaching protocol (TCLP) analysis metals.*

**Response:** The SI Work Plan identifies that three discrete soil samples will be collected and submitted for laboratory analysis for the TCLP volatile organic constituents. To address the NYSDEC's comment, each of the three discrete soil samples will also be submitted for laboratory analysis for ignitability, corrosivity, reactivity, and full TCLP to provide information whether or not the soil exhibits a characteristic(s) of a RCRA hazardous waste for treatment/disposal purposes. Soil samples subjected to RCRA characterization analyses will be determined based on PID readings, odors, staining, and NAPL observed, if any. These samples will be representative of the material that will be treated/disposed.

**Comment:** *The proposed plan also calls for the homogenizing of soil boring intervals. We do not find the homogenizing or compositing of samples for semivolatiles and volatiles acceptable to characterize a site or to determine if the material is hazardous waste. The most contaminated interval of the sample core, based on Hnu readings and olfactory observations, should be collected for analysis. Once the material is determined to be hazardous or non hazardous, compositing for characterization for the treatment or disposal facility use would be acceptable to our Department.*

**Response:** As stated in the previous response, three discrete (i.e., not homogenized or composited) soil samples will be collected and submitted for laboratory analysis for ignitability, corrosivity, reactivity, and full TCLP to determine if the soil exhibits a characteristic(s) of a RCRA hazardous waste for treatment/disposal purposes. Soil samples subjected to RCRA hazardous waste characterization will be determined based on PID readings, odors, staining, or NAPL observed, if any. These samples will be representative of the material that will be treated/disposed. Additional soil samples will be homogenized and analyzed for the suite of constituents presented in the SI Work Plan to meet the characterization criteria required for treatment/disposal by Environmental Soil Management of New York, LLC, assuming the soil is not a RCRA hazardous waste.

**Comment:** *The citizen participation plan should be provided to us prior to the commencement of field work at the site.*

**Response:** The Citizen Participation Plan (CPP) will be provided to the NYSDEC prior to commencement of field work at the site. Based on a telephone conversation between myself and Robert W. Schick, P.E. of the NYSDEC on July 12, 2000, the NYSDEC understands the CPP will be provided to the NYSDEC prior to commencement of field work and that NMPC will coordinate with the neighboring properties regarding the upcoming field activities at the site.

**Comment:** *Please provide the qualifications of the University of Binghamton to perform the cultural resource investigation. These qualifications should include the resumes of project staff, resources, jobs completed of similar size and scope, and references.*

**Response:** Please find attached a copy of Binghamton University's qualifications to perform the cultural resource investigation. NMPC has used Binghamton University to perform cultural resource investigations at other NMPC sites and have previously submitted their qualifications to the NYSDEC. As detailed in the attached qualifications, the scope of services to perform the cultural resource investigation complies with Section 106 of the National Historic Preservation Act, the New York State Historic Preservation Act, the New York State Environmental Quality Review Act and the professional standards of the New York Archaeological Council. The Principal Investigator and supervisory staff meet or exceed the professional qualifications presented in 36 CFR 61.

The cultural resource investigation will not be performed prior to conducting the field activities, however, on-site personnel conducting the field activities will use a guide entitled "Major Aboriginal Projectile Points in New York State", developed by Collamor & Associates, Inc. Archaeological Services & Historic Research, to observe cultural resources, if any. If cultural resources are observed during the field activities, NMPC will notify the New York State Historic Preservation Office.

The field walkover portion of the cultural resource investigation is scheduled to be conducted during the week of July 17, 2000. The remainder of the cultural resource investigation will be completed soon after the field walkover.

We hope that this letter satisfactorily addresses the NYSDEC's comments on the SI Work Plan. Please note that the SI field activities at the 141 Cedar Street site are scheduled to start on July 18, 2000. I will call you prior to then to discuss NMPC's responses and confirm this start date.

If you have any questions or require additional information, please do not hesitate to call me at (315) 428-5652.

Sincerely,



Steven P. Stucker, C.P.G.  
Project Manager

TMM/cmd  
Attachment

cc: James Bacher, P.E., City of Oneida  
Michael W. Sherman, Niagara Mohawk Power Corporation  
William C. Weiss, Esq., Niagara Mohawk Power Corporation  
George M. Thomas, P.G., Blasland, Bouck & Lee, Inc.  
M. Cathy Geraci, Blasland, Bouck & Lee, Inc.  
File



**STAGE 1A ARCHAEOLOGICAL SURVEY  
ONEIDA NGP PROJECT  
ONEIDA, NEW YORK**

**I. INTRODUCTION**

The Public Archaeology Facility (PAF) proposes to conduct a Stage 1A Archaeological Survey for the proposed Oneida NGP Project in the City of Oneida, Oneida County, New York.

Our proposed scope of services for this proposed work complies with Section 106 of the National Historic Preservation Act, the New York State Historic Preservation Act, the New York State Environmental Quality Review Act and the professional standards of the New York Archaeological Council (NYAC 1994).

**II. TECHNICAL APPROACH**

The Oneida NGP Project consists of a 1.5 to 2 acre parcel in order for remediation action to occur. The project area would appear to be sensitive for archaeological resources according to the general topography and known sites in the area. Our scope of services will include the following specific tasks:

1. **Literature Review.** Researchers will compile information on the known prehistory, history and environment of the project area to assess the relative sensitivity for finding cultural resources. This research will include site files checks at the New York State Museum (NYSM), Office of Parks, Recreation and Historic Preservation (OPRHP), and Public Archaeology Facility in Binghamton (PAF). Researchers will also check written histories, prehistories, topographic maps, and soil surveys. These combined data will allow an assessment of project sensitivity. The soils information also guides our recommendations for Stage 1B subsurface testing and target depths of hand excavations. In combination, this information will help structure archaeological field testing strategies.
2. **Project Walkover.** Professional archaeologists will visit the project area to determine areas of prior disturbance or places where cultural resources, such as historic foundations and artifact scatters, are visible on the ground surface. This information will contribute to our assessment of project sensitivity.
3. **Summary Report.** At the conclusion of the Stage 1A survey, we will summarize our background and findings in a brief report. We will submit 4 copies of this report to you, all with original photographs. The report will include recommendations regarding whether a Stage 1B survey is warranted for the project and will offer strategies to complete that survey.
4. **Schedule.** We expect that the Stage 1A survey will take two weeks to complete. The final report will be submitted within two weeks of the completion of the background research. This schedule can be modified to meet the specific deadlines of your project.

5. **Professional Experience.** Our facility has conducted professional archaeological surveys since 1972. We are the statewide archaeological consultant to the New York State Museum and Department of Transportation for highway surveys throughout New York State. In addition, we have completed large segments of pipeline surveys for Tennessee Gas Pipeline, as well as smaller DEC and EPA permit projects for sewer lines, gravel/topsoil mines, property developments, and airports in New York and Pennsylvania.
6. **Equipment and Facilities.** The Public Archaeology Facility has the following equipment and facilities to support our cultural resource management projects:
  - 1) Administrative office complex (672 ft<sup>2</sup>) housing supervisory staff;
  - 2) Graphics Lab (741 ft<sup>2</sup>) with a fully-equipped darkroom, drafting tables, blue-line reproduction machine and map storage.
  - 3) Artifact processing lab (475 ft<sup>2</sup>) with sinks, drying and cataloging tables and temporary artifact storage shelves
  - 4) Analysis Lab (696 ft<sup>2</sup>) with floatation set-up, lab tables and below table storage drawers;
  - 5) Permanent artifact curation rooms (400 ft<sup>2</sup> and 300 ft<sup>2</sup>);
  - 6) Equipment storage area (100 ft<sup>2</sup>);
  - 7) Access to four field vehicles (vans, wagons, suburban);
  - 8) Small and large field and lab equipment;
  - 9) Soil floatation apparatus;
  - 10) Six computers and two printers;
  - 11) Access to the Archaeological Analytical Research Facility, specializing in faunal analysis of Northeastern US collections.
7. **Personnel.** I will serve as the Principal Investigator for this project (credentials attached). The Project Director and crew will be assigned when the project is scheduled. All supervisory staff meet or exceed the professional qualifications in 36 CFR 61.

## QUALIFICATIONS OF THE PRINCIPAL INVESTIGATOR

Dr. Nina M. Versaggi  
Director and Principal Investigator, Public Archaeology Facility

Versaggi received her doctorate in Anthropology from SUNY-Binghamton in 1988, her MA from SUNY University at Binghamton in 1976 and her BA from Rutgers University in 1974. She has been active in professional archaeology since 1972. Professional positions held include Director of the Public Archaeology Facility since 1988, Partner in Compliance Survey Associates for 6 years, Guest Curator at the Roberson Museum and Science Center, and Post-doctoral Fellow at the Hartwick College Museums. She serves as principal investigator for all current and past projects of the Public Archaeology Facility whose recent projects include the Rainbow Plaza Data Recovery in Niagara Falls and the state-wide highway subcontract with the New York State Museum and NYSDOT. She has authored "Hunter to Farmer: 10,000 Years of Susquehanna Valley Prehistory," "Prehistoric Hunter-Gatherer Settlement Models: Interpreting the Upper Susquehanna Valley," and "Upland Foraging Sites in the Northeast: Engendering Prehistory," which are based on NYSDOT and pipeline prehistoric data. She is a member of the board for the Preservation Association of the Southern Tier, and for the New York Archaeological Council she chairs the Professional Survey and Report Standards Committee. She serves as an Adjunct Associate Professor at Binghamton University.

## **RESUMES OF KEY PERSONNEL ONEIDA REMEDIATION PROJECT**

**Nina M. Versaggi, PhD**

**Director and Principal Investigator, Public Archaeology Facility**

Versaggi received her doctorate in Anthropology from SUNY-Binghamton in 1988, her MA from SUNY University at Binghamton in 1976 and her BA from Rutgers University in 1974. She has been active in professional archaeology since 1972. Professional positions held include Director of the Public Archaeology Facility since 1988, Partner in Compliance Survey Associates for 6 years, Guest Curator at the Roberson Museum and Science Center, and Post-doctoral Fellow at the Hartwick College Museums. She serves as principal investigator for all current and past projects of the Public Archaeology Facility whose recent projects include the Rainbow Plaza Data Recovery in Niagara Falls and the state-wide highway subcontract with the New York State Museum and NYSDOT. She has authored "Hunter to Farmer: 10,000 Years of Susquehanna Valley Prehistory," "Prehistoric Hunter-Gatherer Settlement Models: Interpreting the Upper Susquehanna Valley," and "Upland Foraging Sites in the Northeast: Engendering Prehistory," which are based on NYSDOT and pipeline prehistoric data. She is a member of the board for the Preservation Association of the Southern Tier, and for the New York Archaeological Council she chairs the Professional Survey and Report Standards Committee. She serves as an Adjunct Associate Professor at Binghamton University.

**Christopher D. Hohman, RPA**

**Assistant to the Director and Project Director, Public Archaeology Facility**

Hohman received his MA in Anthropology from the University of Connecticut in 1986 and his BA in Anthropology from the University of Rhode Island in 1983. He has worked in professional archaeology since 1983 and in 1987 joined the staff of the Public Archaeology Facility. From 1990 to 1991, Hohman served as one of the Coordinators at Garrow and Associates for the 300 + mile Iroquois Gas pipeline project in New York and Connecticut. He has served as project director on a wide range of site examinations and surveys including the Whirlpool Rapids Reconnaissance, and the Rainbow Plaza Reconnaissance and Data Recovery in Niagara Falls. His research interests include Northeast prehistory and 19<sup>th</sup> century rural and urban settlements. The author of numerous cultural resource management reports, he currently directs projects for PAF's statewide highway contract with the State Museum and other projects in New York State and Pennsylvania. Member of ROPA (Register of Professional Archaeologists), the Council for Northeast Historic Archaeology and the Archaeological Society of Connecticut.

**Barbara Ross**

**Assistant Director for Research, Public Archaeology Facility**

Ross received her MS in Public Archaeology from Rensselaer Polytechnic Institute in 1981 and her BA in Anthropology from SUNY at Albany in 1978. She has been active in professional archaeology since 1979 as a principal investigator with various NYS private firms and the NYS Museum's Cultural Resource Survey Program (NYSM CRSP). She served as Assistant Director of the the NYSM CRSP from 1983 - 1995. This position included three years as interim director as well as liaison with NYS client agencies and SUNY subcontractors. Her primary area of interest is late eighteenth and nineteenth century rural residential/agricultural archaeology and architecture/landscape preservation. She joined the staff of the Public Archaeology Facility in 1996 to assist the program director with administration of its statewide highway contract with the NYSM.

**Brian Russell Grills**  
Researcher, Public Archaeology Facility

Grills received his B.A. in History (with a minor in Anthropology) from Ohio Wesleyan University in 1994. He is currently enrolled in the M.A. program in Anthropology at Binghamton University (SUNY), specializing in Northeastern US archaeology. Since June 1995 he has been employed with the Public Archaeology Facility at Binghamton University, Binghamton, NY. As a field technician, he worked on a wide range of reconnaissance, site exams, and data recoveries. His archaeological experience also included working in the artifact processing laboratory assisting with cleaning and cataloging prehistoric and historic artifacts, prehistoric lithic analysis, and soil flotation. Since 1997, he has also participated in PAF projects as a researcher, crew chief and field director. Experience with other groups included working at the Erie Canal Boat Landing Museum in Madison County, NY, student in the SUNY-Cortland fieldschool along the Otselic River, and excavator at the Tarbat Ness Archaeological Research Project, UK.

**Maria Pezzuti**  
Administrative Assistant, Public Archaeology Facility

Pezzuti has administered contracts and grants since 1984, where she was on the staff of the Research Foundation of SUNY, Office of Contract and Grant Administration in Albany, NY. There she monitored the NYS Education Department Highway Salvage contract for all participating SUNY campuses. In 1986 she moved to the Office of Sponsored Programs at Binghamton University where she was responsible for preparing research proposals for submittal. She joined the Public Archaeology Facility in 1989 and since 1992 she has held a concurrent half time positions in the Department of Biological Sciences as Project Coordinator of educational grants from the Howard Hughes Medical Institute. In addition, she has also worked as curator of the Vestal Museum, a small historical museum, and held the positions of secretary and President for the Vestal Historical Society.

**Mary Lou Supa**  
Drafter, Public Archaeology Facility

Supa received her BA and completed graduate course work in Art History. Her work experience ranges from the museum field (collection care and preservation, educational programming and curator) to graphic design and desktop publishing. Supa serves as the drafting director for Public Archaeology Facility providing project maps, site maps and soil profiles for all projects.

## REFERENCES FOR THE PUBLIC ARCHAEOLOGY FACILITY

Stage 1, 2, and 3 Cultural Resources Management Surveys, Statewide subcontract with the New York State Museum and New York Department of Transportation. Completed between 30-60 projects per year, including some in the Niagara region of western New York.

Prepared for: The New York State Museum  
State Education Department  
Albany, New York 12230

Contact: Dr. John P. Hart 518-474-3895

Stage 1 and 2 Cultural Resource Surveys for various airport expansions, road reconstructions, and school expansions.

Prepared for: McFarland-Johnson Engineers, Inc.  
49 Court Street, Metrocenter  
Binghamton, NY 13902-1980

Contact: David Bors 607-723-9421

Stage 1 and 2 Cultural Resource Surveys for the USDA Natural Resources Conservation Service (NRCS). PAF completes between 8 and 12 wetland restoration and barnyard improvement projects per year on tight schedules.

Prepared for: US Department of Agriculture  
Natural Resources Conservation Service  
The Galleries of Syracuse  
441 South Salina Street, Suite 354  
Syracuse, NY 13202-2450

Contact: Florence Swartz 315-477-6533



*Transmitted Via U.S. Mail*

November 3, 2000

Mr. John A. Helmeset, P.E.  
Environmental Engineer 2  
MGP Section  
Bureau of Western Remedial Action  
Division of Environmental Remediation  
New York State Department of Environmental Conservation  
50 Wolf Road  
Albany, New York 12233-7010

Re: Test Pits  
Niagara Mohawk Power Corporation  
Oneida (141 Cedar Street) Former MGP Site  
Oneida, New York

Dear Mr. Helmeset:

As a follow up to the discussions we had following the working meeting at the Oneida City Hall on October 19, 2000 regarding the status of the investigation work at the Oneida (141 Cedar Street) Former Manufactured Gas Plant (MGP) Site, this letter outlines recommendations for limited additional subsurface investigations at the site. As we discussed, the limited investigations are proposed to be completed to evaluate potential migration along a former surface water drainage channel located to the southeast of the 141 Cedar Street site. Evidence of the presence of this surface water drainage channel is based on a 1875 Atlas of Madison County (Beers, 1875) showing this area of the site. A copy of this map is attached.

As outlined during the October 19, 2000 meeting, visual indications (e.g., tar or oily sheens) of potentially MGP-related material were observed in three (SB-8, SB-11, and SB-12) of the eight soil probes completed south and southeast of the 141 Cedar Street site. A summary table of the subsurface investigation findings is attached along with a figure showing the locations of the soil probes. A temporary monitoring well, PZ-1, was installed at one of the soil probe locations (SB-9) and ground water was sampled from this well for Target Compound List/Target Analyte List constituents and total cyanide. Volatile organic compounds, semi-volatile organic compounds, polychlorinated biphenyls, pesticides, and total cyanide were not detected in this ground-water sample. Based on the findings of these investigations, visual indications (e.g., tar or oily sheens) were not observed except at the three soil probe locations directly southeast of the retaining wall; however, to assess the possible migration of potentially MGP-related constituents along the former surface water drainage channel, Niagara Mohawk Power Corporation (NMPC) is recommending excavating several test pits approximately 15 feet apart along the inferred location of the former surface water drainage channel. We currently anticipate that these test pits will be excavated in the assumed downstream direction from soil probe SB-13 and one in the upstream direction from SB-14. Additional test pits may be excavated based on the observations made in the field (e.g., if tars or oily sheens are observed in the outermost test pits).

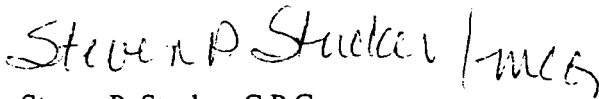
The test pits will be excavated using a decontaminated, rubber-tired backhoe. The actual location of the test pits will be determined in the field. Continuous air monitoring will be conducted as

indicated in the *Preliminary Remedial Design Work Plan*, dated May 2000, for the Sconondoa Street Former MGP Site. Test pit materials will be visually observed and described with respect to the depth. Photographs of the subsurface soil will be taken for future reference. Photoionization detector measurements will be taken on soil samples from the test pit to provide an indication of the relative concentration of total organic volatile compounds. Materials removed from the test pits during excavation will be placed on polyethylene sheeting. Upon completion, the materials from the test pit will be placed back into the excavation. The installation of these test pits are proposed to be conducted in the near future. NMPC will contact the New York State Department of Environmental Conservation (NYSDEC) once a date has been determined to conduct the test pitting activities.

A letter summarizing the field activities and analytical results of the Supplemental Investigation (SI) conducted at the site will include the findings of these test pits. As discussed following the October 19, 2000 meeting, the SI summary letter report will be submitted to the NYSDEC in early December 2000.

Please feel free to contact me if you have any questions.

Sincerely,

A handwritten signature in black ink that reads "Steven P. Stucker / mcs".

Steven P. Stucker, C.P.G.  
Project Manager

GMT/cmd  
Attachments  
93301750.doc

cc: James Bacher, P.E., City of Oneida  
Michael W. Sherman, Niagara Mohawk Power Corporation  
William C. Weiss, Esq., Niagara Mohawk Power Corporation  
George M. Thomas, P.G., Blasland, Bouck & Lee, Inc.  
M. Cathy Geraci, Blasland, Bouck & Lee, Inc.  
File

**Niagara Mohawk Power Corporation  
Oneida (141 Cedar Street) Former MGP Site  
Site Investigation Activities**

**Offsite Boring Summary**

| Boring ID | PID / (Sample Interval) | Sample Description   |
|-----------|-------------------------|--|
| SB-7      | 0.0 (0-4')              | 0.0-0.7' - Tan brown fine SAND, little medium to coarse Sand, fine Gravel, loose, organics (earth worm).<br>0.7-1.2' - Medium to coarse SAND and fine GRAVEL, [FILL], brick fragments, loose, glass fragments.   |
|           | NA (4-8')               | 4.0-8.0' - No recovery.  |
| SB-8      | 0.0 (0-4')              | 0.0-1.8' - Light brown fine to medium SAND and roots.  |
|           | 206 (4-6')              | 4.0-5.2' - Light brown fine to coarse SAND, some Gravel and organics; black tar throughout sample.   |
|           | 25.2 (6-8')             | 5.2-8.0' - Olive-brown to dark brown PEAT, little black staining.  |
| SB-9      | 0.9 (0-2')              | 0.0-1.0' - Medium reddish-brown fine to medium SAND, little coarse Sand, fine Gravel.  |
|           | 2.1 (2-4')              | 1.0-2.1' - Loose fine to coarse SAND, fine Gravel, black, gray, yellow, and red Fill, trace Slag and Brick fragments.<br>2.1-3.5' - Medium brown fine SAND, some Silt, little medium to coarse Sand, fine Gravel, Brick, and Cinders [FILL], moist to wet. |
|           | 0.0 (4-6')              | 4.0-5.0' - Medium to light brown fine SAND and SILT, little Clay, trace coarse Sand, stiff, wet.   |
|           | 1.0 (6-8')              | 5.0'-7.0' - Black Peaty SILT, grading to dark brown-black PEAT.  |
|           |                         |  |
| SB-10     | 2.5 (0-4')              | 0.0-4.0' - Red-brown fine to coarse SAND, organics, and broken concrete [FILL], black staining.  |
|           | 0.0 (4-7')              | 4.0-5.4' - red-brown to tan fine to medium SAND, trace to little Silt and Gravel.  |
|           | 0.0 (7-8')              | 5.4-8.0' - Olive-gray to dark brown PEAT, black staining.  |
| SB-11     | 0.0 (0-4')              | 0.0-4.0' - Red-brown fine to coarse SAND, organics, and slag [FILL].   |
|           | NA (4-8')               | 4.0-8.0' - No recovery in sleeve; black tar on gravel found in shoe.   |
| SB-12     | 0.0 (0-3')              | 0.0-3.5' - Red-brown fine to medium SAND, organics, and little Gravel; trace oily sheen; black staining, and odor.   |
|           | 38.7 (3-3.5')           | Refusal at 3.5' bgs.   |
| SB-13     | 0.4 (0-3')              | 0.0-4.0' - Brown fine to medium SAND, organics, slag, and brick [FILL]; yellow staining.   |
|           | 0.0 (3-4')              |  |
|           | 0.0 (4-6')              | 4.0-7.2' - Olive/yellow-brown to red FILL, fine to coarse Sand; black staining.  |
|           | 0.0 (6-8')              | 7.2-8.0' - Dark brown to black PEAT.   |
| SB-14     | 0.8 (0-2')              | 0.0-1.8' - Medium brown fine to medium SAND, trace coarse Sand and fine Gravel, loose [FILL].  |
|           | 0.0 (2-4')              | 1.8-2.5' - Coarse SAND and fine GRAVEL, [FILL], brick fragments, slag, wet, yellowish brown, red brick.  |
|           | 4.3 (4-6')              | 4.0-5.8' - Medium brown fine SAND, some Silt, wet, soft, trace medium to coarse Sand, possible brick fragments.  |
|           | 28.2 (6-8')             | 5.8-6.2' - Dark brown-black Peaty SILT, little black wood fragments, possible faint odor.<br>6.2-7.5' - Dark brown-black PEAT, low density, moist.   |
|           |                         |  |
| SB-15     | 0.0 (0-2')              | 0.0-1.3' - Medium brown fine to medium SAND, trace coarse Sand, loose, moist.  |
|           | 0.0 (2-4')              | 1.3-2.2' - Fine to coarse SAND, trace Gravel, red Brick, black-gray Fill, possible Slag, broken rock, trace organics, wet.   |
|           | 5.5 (4-6')              | 4.0-5.3' - Medium brown fine SAND, little Silt and medium Sand, trace coarse Sand and fine Gravel, possible brick fragments, wet.  |
|           | 0.0 (6-8')              | 5.3-7.6' - Medium brown-black PEAT, low density, moist.  |

**Notes:**

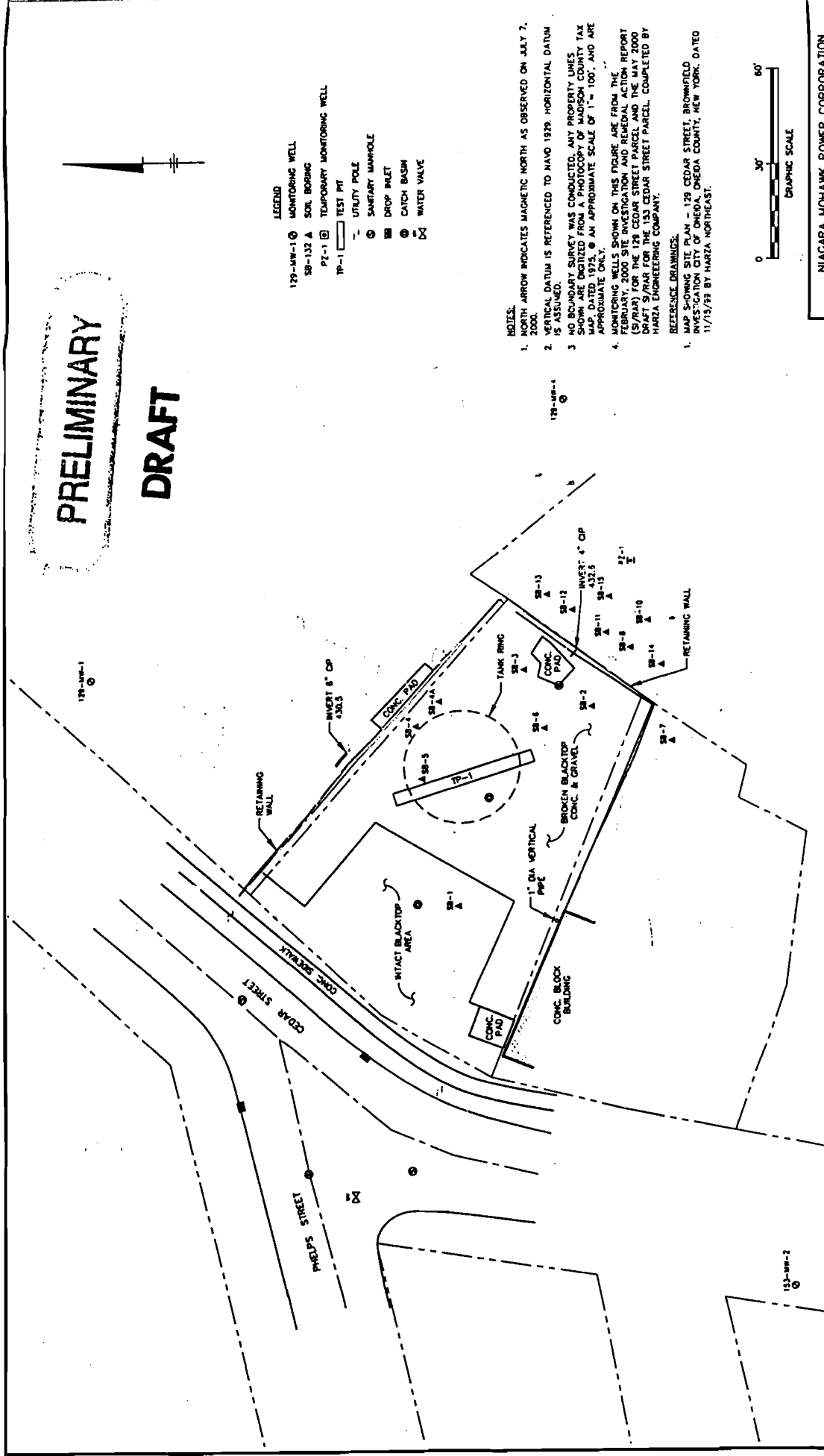
PID - Headspace screening measurements of the relative concentration of volatile organic compounds in soil (in parts per million) obtained with a photoionization detector.

Sample Interval - Indicates the depth interval (in feet below ground surface) penetrated by the direct push sampling device. The intervals indicated in the soil description represent actual soil recovery.



# PRELIMINARY

# DRAFT





## **Attachment 2**

### **Stage 1A Cultural Resources Assessment**



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# **PUBLIC ARCHAEOLOGY FACILITY REPORT**

Department of Anthropology  
State University of New York at Binghamton  
Binghamton, New York

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**CULTURAL RESOURCE MANAGEMENT SURVEY  
STAGE 1A CULTURAL RESOURCE ASSESSMENT**

**ONEIDA MGP PROJECT  
CITY OF ONEIDA  
MADISON COUNTY, NEW YORK  
MCD 05340**

**PREPARED BY:**

**CHRISTOPHER D. HOHMAN**

**PREPARED FOR:**

**BLASLAND, BOUCK AND LEE, INC.  
6723 TOWPATH ROAD  
SYRACUSE, NY 13214**

**AUGUST 2000**

**REPORT OF FIELD RECONNAISSANCE  
STAGE 1A CULTURAL RESOURCE ASSESSMENT  
ONEIDA MGP PROJECT**

**Permit Applicant:** Blasland, Bouck and Lee, Inc.

**Permit Number:**

**Location:** City of Oneida, Madison County, NY (MCD 05340)

**Report prepared by:** Christopher D. Hohman

**Date:** August 2000

**Affiliation:** Public Archaeology Facility  
Binghamton University  
Binghamton, New York 13902  
(607) 777-4786

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**PROJECT SUMMARY**

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A Stage 1A cultural resource assessment was requested for the Oneida MGP project in the City of Oneida, Madison County, New York. The project consists of a parcel at 143 Cedar Street that is approximately .22 acres and will impact between .6 to 6 m (2-19 ft) below the surface on the property. Because the project is located on the site of a former gas works, the project may involve the remediation of the property. The project is situated at an elevation of 136 m (445 ft) ASL. Background research on the prehistory of the surrounding area indicated that the project is situated in a zone of moderate prehistoric sensitivity. Eight known prehistoric/contact period sites are located within 3.2 km (2 mi) of the parcel; four of the eight sites are noted as Contact period villages or cemeteries. Background research on the history of the surrounding area indicated that the parcel is in a zone of high historic sensitivity. There are at least two map documented structures (MDSs) from the 19<sup>th</sup> century located within the project area.

A walkover of the project area revealed that the project is set on top of fill ranging from 2.1-4 m (7-13 ft) in height. The fill has been enclosed and sealed by fieldstone and cement walls on three sides with the fill bordering the sidewalk on the west side of the parcel. A storm drain has been placed within the central portion of the parcel (Attachment F, Photos 1-16).

Since the footprint of the former structures covered approximately 95% of the project area and because of the amount of fill that was placed within the project area in the mid 20<sup>th</sup> century, the project area has limited research potential in regards to 19th-20th century commerce and industry in the City of Oneida. Because the project area is located away from any major water resources (600 m [2000 ft] to Oneida Creek) and ground disturbance has occurred throughout most of the project area, there appears to be limited potential for prehistoric sites within the project area. Therefore, it is recommended that a Stage 1B archaeological survey is not needed for this parcel.

## Part 1: DOCUMENTARY RESEARCH ADDENDUM SITE IDENTIFICATION

### A. Documentary Research Addendum (if needed)

1. ..X.. Local site inventory checked (specify)  
Public Archaeology Facility, Binghamton University  
  
..X.. Division for Historic Preservation/New York State Museum
2. .... Informants interviewed (name, address, specialty)
3. ..X.. Other sources checked (specify)

Abbott, Arthur J. (1943) *Oneida in Madison County today*.

Beauchamp, William M. (1900) *Aboriginal Occupation of New York*. New York State Museum Bulletin 32: 7, Albany.

Beers, D.G. (1875) *Map of Madison County, New York*. Pennsylvania: Pomeroy, Whitman and Co.

Evans, Gordon (1853) *Topographical Map of Madison County, New York*. Philadelphia: Anthony Byles.

French, J.H. (1859) *Gillette's Map of Madison County*. Philadelphia: Jonathan E. Gillette.

Funk, Robert E. (1993) *Archaeological Investigations in the Upper Susquehanna Valley, New York State*. Buffalo: Persimmon Press.

Lehman, Karl H, editor. (1943) *Madison County today*. Oneida Castle, NY.

Parker, Arthur C. (1920) *The Archaeological History of New York*. New York State Museum Bulletins 235-238. Albany: New York State Museum.

Pratt, Peter P. (1976) *Archaeology of the Oneida Iroquois, Volume 1*. Occasional Publications in Northeastern Anthropology, No. 1. Greenfield, MA: Minott Printin and Binding Co.

Ritchie, W. (1980) *The Archaeology of New York State*. Harrison, NY: Harbor Hill Books.

Ritchie, William and Robert Funk (1973) *Aboriginal Settlement Patterns in the Northeast*. Albany: New York State Museum, Memoir 20.

Sanborn Map Company (1890, 1895, 1899, 1904, 1909, 1914, 1923, 1930 and 1956) *Fire Insurance Maps of Oneida*. New York: Sanborn Map Company.

Smith, John E. (1899) *Our county and its people; a description and biographical record of Madison County, New York*. Boston: Boston history company.

Snow, Dean R. (1995) *Mohawk Valley Archaeology: The Sites*. Albany: The Institute for Archaeological Studies.

Tuck, James A. (1971) *Onondaga Iroquois Prehistory: A Study in Settlement Archaeology*. Syracuse: Syracuse University Press.

United States Department of Agriculture (1981) *Soil Survey, Madison County, New York*. Washington, D.C.: USDA Soil Conservation Service with Cornell University Agricultural Experiment Station.

United States Geological Survey (1955) *Oneida, 7.5 minute quadrangle*.

Versaggi, Nina (1986) *Hunter to Farmer: 10,000 Years of Susquehanna Valley Prehistory*. Binghamton, NY: Roberson Center for Arts and Sciences.

Versaggi, Nina (1987) *Hunter-Gatherer Settlement Models and the Archaeological Record: A Test Case for the Upper Susquehanna Valley of New York*. PhD Dissertation. Binghamton, NY.

Whitnall (1943) Geography of Madison County in *Madison County today*.

Whitney, Luna M. Hammond (1872) *History of Madison County, New York*. Syracuse: Truair, Smith and Co.

### **Results of Documentary Research: ENVIRONMENT AND SOILS**

The project area is located at the northern end of Madison County. Madison County is comprised of three main geographical forms: the glacial plain at the northern end; the Allegheny Plateau which covers the rest of the county and the watersheds which divide the drainage into three systems. Oneida is located on the west side of Oneida Creek, which flows into Oneida Lake approximately 10 km (6 mi) to the northwest of the project area. Oneida Creek is located approximately 600 m (2000 ft) east of the project area, although small waterways are noted adjacent to the project area on 19<sup>th</sup> century historic maps (Attachments D.1-D.3). As noted by Abbott (1943), much of Oneida was located on a swamp adjacent to Oneida Creek, with portions being filled in from the mid 19<sup>th</sup> century through the early 20<sup>th</sup> century. The project area itself is located at an elevation of approximately 136 m (445 ft), 2.1-4 m (7-14 ft) above the surrounding surface to the east of the project limits. The setting of the project area consists of an asphalt covered lot with fill enclosed by fieldstone and cement retaining walls (Photos 1-16, Attachment F).

Although the soil survey map of the project area suggests that Teel (Te) silt loam is located within the project area, visual observation suggests that this natural soil horizon may be located adjacent to the project boundaries or under the fill within the project limits. Teel silt loam is formed in silty alluvium that was washed from upland glacial drift. They are nearly level, along streams and subject to flooding. Within an agricultural field, Teel silt loam commonly has a profile of an Ap horizon (12 cm [5 in] of dark brown silt loam), a B21 horizon (28 cm [11 in] of dark reddish gray silt loam), a B22 horizon (55 cm [22 in] of reddish brown silt loam) and a C horizon (50 cm [20 in] of dark reddish gray silt loam that grades to very fine sandy loam (USDA, 1981) (Attachment B.2).

Based on the soil descriptions and the environmental setting, the probability is moderate that prehistoric sites exist in the project area. The project area is located approximately 600 m (2000 ft) west of Oneida Creek, although other water resources may have been available adjacent to the project area. The project area itself is comprised of fill in the vicinity of 19<sup>th</sup> century structures, suggesting that impacts have been made to the original surface underneath these structures.

### **Results of Documentary Research: PREHISTORY**

During the Paleo-Indian period (10,000-8,000 B.C.) people used the rich resources of low-lying streams and riverbanks for the mainstay of their existence, while using more elevated flood plains and glacial kames for their temporary residences (Ritchie 1980). Paleo-Indian people led a highly mobile life, following megafauna (e.g., caribou, mastodon and moose), fishing, and collecting available plant resources.

People followed a seasonally mobile lifestyle during the Archaic period (8,000-1,500 B.C.). As megafauna became less available, deer, bear, elk, and a wide variety of smaller mammals and birds were substituted (Ritchie 1980). Settlement types for the Early and Middle Archaic periods include large camps and smaller, more temporary camps. Late Archaic sites (large central base camps and temporary camps) are generally located near major waterways, with many special purpose camps and general resource processing areas occurring on a variety of different landforms.

The Transitional period (1,500-1,000 B.C.) is characterized by steatite vessels and broad spear points and a subsistence practice focused on hunting, gathering and fishing. Ceramic vessels were manufactured during this period. Small, temporary camps, often oriented toward river or coastal areas, typify settlement patterns during this period (Ritchie and Funk 1973).

The use of ceramic vessels became more widespread during the Early Woodland period (1,000-200 B.C.). In addition, long-distance contact with cultural groups in the Great Lakes region developed. This contact continued into the subsequent Middle Woodland period (200 B.C. - A.D. 800). During the Early and Middle Woodland periods, subsistence was distinguished by a greater reliance on native plant species (Chenopodium, sunflower, tobacco) (Funk 1993; Ritchie 1980). Settlement types for these periods include large and small camps. It was not until the Late Woodland period (A.D. 800-1300) and continuing into the Iroquois period (A.D. 1300-1550) that permanent villages, along with camps, became the dominant settlement pattern. In addition, at this time, maize, beans and squash horticulture, supplemented by the native cultigens, was added to the previous subsistence practices of hunting, fishing and gathering.

The major published sources on New York prehistory and the Oneida and Onondaga Iroquois (Beauchamp 1900, Parker 1920, Ritchie 1980, Ritchie and Funk 1973, Pratt 1976, Tuck 1971 and Snow 1995) do not list any sites within, or near, the project area. However site distributions within surrounding areas in Madison and Oneida Counties suggest fairly dense occupations adjacent to the Oneida River and the waterways that flow through the counties, including Oneida Creek.

A site files search at the New York State Museum/Office of Parks, Recreation, and Historic Preservation (OPR&HP) supplemented the background research conducted at the Public Archaeology Facility (Attachment C). The site files recorded eight prehistoric/contact period sites within 3.2 km (2 mi) of the project area. The sites range in size from traces of occupation to Oneida villages, with occupations being known from the Woodland through Contact periods. However, the sites are located mainly along Oneida Creek, Cowaselon Creek and small tributaries of Oneida Creek. The project area, itself, is situated approximately 600 m (2000 ft) west of Oneida Creek and no other water resources are located in close proximity to the project. The project is also located within the boundaries of the original Oneida Indian Reservation (Lehman 1943).

Considering what is known of the topographic and physiographic location of the Oneida MGP project and the known site types and settings, we can expect that small, dispersed resource processing locations associated with villages along Oneida Creek may occur within the project area. However, 19th and 20th century land modifications within the project area may have erased these early traces in that portion of the project.

## **Results of Documentary Research: HISTORY**

The City of Oneida was originally part of the Town of Sullivan, and was part of property that became the Town of Lenox in 1809. In 1896, the Town of Lenox was split into three towns: Lenox, Lincoln and Oneida (Smith 1899). Settlement began in the Town of Sullivan/Lenox in 1792, with the arrival of settlers, the Klock family and others, at the site of Clockville, approximately 10 km (6 mi) to the southwest of the project area. Much of the location of the City of Oneida was originally a vast hemlock swamp (Abbott 1943). In 1817, the Erie Canal was built and the Oneida feeder took water from Oneida creek to maintain the water level of the canal. The canal and its feeder, as well as the Cowaselon feeder, aided in draining the swamp.



In 1829 and 1830, Sands Higginbotham, a merchant from the Town of Vernon, purchased a considerable amount of property from individuals and from the state. In 1834, he built a residence and moved to Oneida. A few log cabins and homes were built there, mainly by German and Irish immigrants (Abbott 1943). In 1837, the Syracuse and Utica Railroad Company located their railroad across his farm and made a station there - "Oneida Depot". Higginbotham gave his land to the railroad in return for their agreement to stop every train there for meals. The railroad used wood for fuel and the wood in the swamp was cut and stacked for their use, thus clearing the land in the vicinity of the region. In 1839, Higginbotham constructed a hotel there, the Railroad house. He then proceeded to sell lands to settlers on such terms to attract residents and develop the place. The railroad provided faster transportation for products (especially agricultural products and lumber). In addition, the canal feeder was plied, thus allowing cheap transportation to Albany and other points.

The canals and railroad not only brought in laborers, but factories and centers of commerce also developed because of these transportation facilities (Whitnall 1943). Through the mid 19th century, Oneida was still considered a village and consisted primarily of small industries (centered around lumbering and sash/blind factories) (Whitney 1872). However, by the late 19th century, manufacturing became prevalent within the community. Important industries such as the National Casket Company, the Oneida Iron Works, the Westcott Chuck (wedges) Company, the Oneida National Chuck Company, and the Oneida Silver Manufacturing Company and many smaller establishments contributed materially to the growth and activity of the city (Smith 1899).

The City of Oneida has continued to grow, although industries have faded since the mid 20th century. Much of Oneida now centers around smaller industries and commercial businesses within the community.

The historic maps document continual use of the property from as early as the mid 19th century through the mid 20th century. In 1853 (Attachment D.1), there was a shop (possibly a sawmill shop) (MDS A) within the project boundaries. Between 1859 and 1875, two other structures were constructed on the property (MDS B). MDS A was not labeled in 1859 or 1875 as to its use, while the two other structures (MDS B) were noted as Gas Works (Attachment D.3). One of the Gas Works structures was a commercial building, while the second structure was circular, suggesting that it was used for gas storage. The 1890 Sanborn fire insurance map (Attachment D.4) is much more detailed and reveals that the 1853 structure (2 story brick) was being used in 1890 as a blacksmith shop, upholstery business and contains a one story structure to its rear with a brass and iron foundry. The gas works business, known in 1890 as the Oneida Gas Works, was comprised of four different structures, with three of the structures being attached: a 1 story coal shed, a one story office and a one story building; the fourth structure was the circular gasometer building to the east of the business office. In addition, a 1 ½ story storehouse was located to the east of the Upson and Holden Carriage and Sleigh Factory and may be located within the boundaries of the current project area.

Between 1890 and 1914, the gas works buildings served a variety of commercial uses: Oneida gas works, Oneida Rubber Tire Works, and the Coles Tool and Machine Company. Although the buildings continued to be used for blacksmith and wheelright supplies and auto sales and service through 1930, the gas storage structure was removed between 1914 and 1923. Between 1895 and 1899, the 1-story foundry structure was removed and a new 1 story structure was built in its place. The 1853 structure continued to be used for a variety of commercial activities: tinshop, carriage shop, painting, the Ryan Burkhart printing, the W.S. Ryan Printing Company, a machine shop, and a bicycle shop. The structure was removed between 1930 and 1956.

Between 1930 and 1956, following the removal of the gas works buildings and the commercial building, the project area was used as a used car sales lot, with one story overhangs along the north and south ends of the lot, and a small office constructed in the middle of the lot.

The historic maps document continual use of the project area from the mid 19<sup>th</sup> century through the mid 20<sup>th</sup> century. The footprint of the structures for the gas works, as well as the southern commercial structure suggests that although there may be a high sensitivity for historic resources (i.e., foundation walls), there is a low sensitivity for resources with research potential (i.e., sheet midden, shaft features) within the project boundaries.

## **Documentary Summary**

..... no sites reported

...X... sites reported (describe briefly)

Eight prehistoric/contact period sites are located within 3.2 km (2 mi) of the project area. The sites suggest occupation of the region from at least the Woodland to Contact periods. The sites are generally located along Oneida Creek and tributaries that flow into Oneida Creek. No known historic sites, beyond those identified as Native American contact sites, were identified within 3.2 km (2 mi) of the project area. Two MDSs, A and B, are located within the project boundaries. MDS A was a commercial building used for a wide variety of businesses and MDS B was a building complex used for a variety of commercial and industrial purposes, including a gasworks business. The site files also note that the project area may be within the viewshed of a National Register Historic District located along Main, Broad, Grove, Wilbur and Walnut Streets.

## **B. Field Investigation**

### **1. Methodology**

#### **a. Description of structure for survey team (number, organization).**

Dr. Nina M. Versaggi, Director of PAF and Christopher D. Hohman, Project Director with PAF, supervised this project and authored this report. Project maps were drawn by James Levandowski and all administrative work was carried out by Maria Pezutti and Annie Pisani. All field and research personnel meet or exceed the minimum qualifications for professionals in archaeology.

#### **b. Date of survey and soil description of general and subsurface conditions (including season, ground visibility, and relative wetness of soil).**

Walkover of the project area was conducted on July 17, 2000. The surface was covered by asphalt, with retaining walls being visible on the edge of the lot.

#### **c. Outline of walkover strategy.**

The project area was visually inspected by Hohman for an assessment of research potential within the project boundaries. This was accomplished by walking within and adjacent to the project boundaries.

#### **d. Description of general project characteristics.**

The nature of the topography within and adjacent to the project area, as well as the retaining walls on the north, east and south sides suggests that soils within the project area consist of fill constructed on top of a base of solid subsoil. In addition, a water drain has been constructed through the middle of the lot, suggesting that the interior of the lot has been impacted since being filled.

- e. **Description of intensity of coverage and rationale for excluding areas from survey. Attach a map with location and type of each excavation unit, and areas surface inspected. Any areas not surveyed should be clearly delineated.**

All of the project area was observed for an assessment of research potential. The original surface underneath the fill could not be observed, although the construction of all of the 19<sup>th</sup> century structures within the project area suggest that there is little potential for remaining topsoil.

- f. **Description of problems encountered during survey which may have influenced results.**

None

### **Results of Field Investigation**

- ...X... **no sites identified**  
..... **site(s) identified**

Walkover did not identify sites within the project limits for the Oneida MGP Project.

### **2. Recommendations**

- ...X... **no additional work**  
..... **additional investigation**  
..... **project modification to avoid sites**

Because the size of the structures' footprints within the project boundaries covered approximately 95% or more of the lot and the fact that the structures impacted the original surface, we recommend that there is little potential for any prehistoric resources remaining within the limited amount of area not covered by structures. In addition, the commercial use of much of the property and the extent of ground disturbance suggest negligible potential for historic resources with research value.

### **3. Rationale**

- a. **Evaluate the effect of the proposed undertaking on identified cultural resources.**

N/A

- b. **Describe possible precautions, protective measures or project modifications that would avoid or alleviate these impacts.**

N/A

**c. Identify sites and/or areas which require additional study.**

None

**d. Outline the nature and extent of additional investigation(s).**

None

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**END PART 1**

**If site evaluation is not completed at this time, proceed to Part 3.**

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### PART 3: SUPPORTIVE DATA

Reports should include the items listed below. Bracketed information is optional. Put a check next to each item appended.

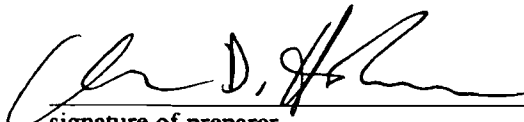
**PLEASE NOTE:** Most attachments listed below often provide precise locational and compositional data on archaeological sites. This information is confidential to protect the resource from vandalism. All attachments with site specific information should be omitted from report copies which will be available to the general public.

- ..A..      qualifications of the principal investigator(s)
- ..B..      topographic and soils maps with project area noted
- ..C..      site files checks (Confidential: Not for Public Release)
- ..D..      copies of relevant historic maps
- ..E..      map(s) of test locations, field inspection, and areas of cultural material; maps must have title, legend, bar scale, and directional arrow.
- .....      record of soil stratigraphy in each test unit.
- ....      artifact catalog
- ..F..      photographs of the project area
- .....      OPR&HP Prehistoric Site form

---

**Certification:** I certify that I directed the cultural resource investigation reported here, that my observations and methods are fully reported, and that this report is complete and accurate to the best of my knowledge.

8/1/00  
\_\_\_\_\_  
date

  
\_\_\_\_\_  
signature of preparer

**ATTACHMENT A:**  
**Qualifications of the Principal Investigators**

**Dr. Nina M. Versaggi**  
**Director and Principal Investigator, Public Archaeology Facility**

Versaggi received her doctorate in Anthropology from SUNY-Binghamton in 1988, her MA from SUNY University at Binghamton in 1976 and her BA from Rutgers University in 1974. She has been active in professional archaeology since 1972. Professional positions held include Director of the Public Archaeology Facility since 1988, Partner in Compliance Survey Associates for 6 years, Guest Curator at the Roberson Museum and Science Center, and Post-doctoral Fellow at the Hartwick College Museums. She serves as principal investigator for all current and past projects of the Public Archaeology Facility whose recent major projects include the Rainbow Plaza Data Recovery in Niagara Falls and the state-wide highway subcontract with the New York State Museum and NYSDOT. She has authored "Hunter to Farmer: 10,000 Years of Susquehanna Valley Prehistory," "Prehistoric Hunter-Gatherer Settlement Models: Interpreting the Upper Susquehanna Valley," and "Upland Foraging Sites in the Northeast: Engendering Prehistory," which are based on NYSDOT and pipeline prehistoric data. She is a member of the board for the Preservation Association of the Southern Tier, and for the New York Archaeological Council she chairs the Professional Survey and Report Standards Committee. She serves as an Adjunct Associate Professor at Binghamton University.

**Christopher D. Hohman, RPA**  
**Project Director and Assistant to the Director, Public Archaeology Facility**

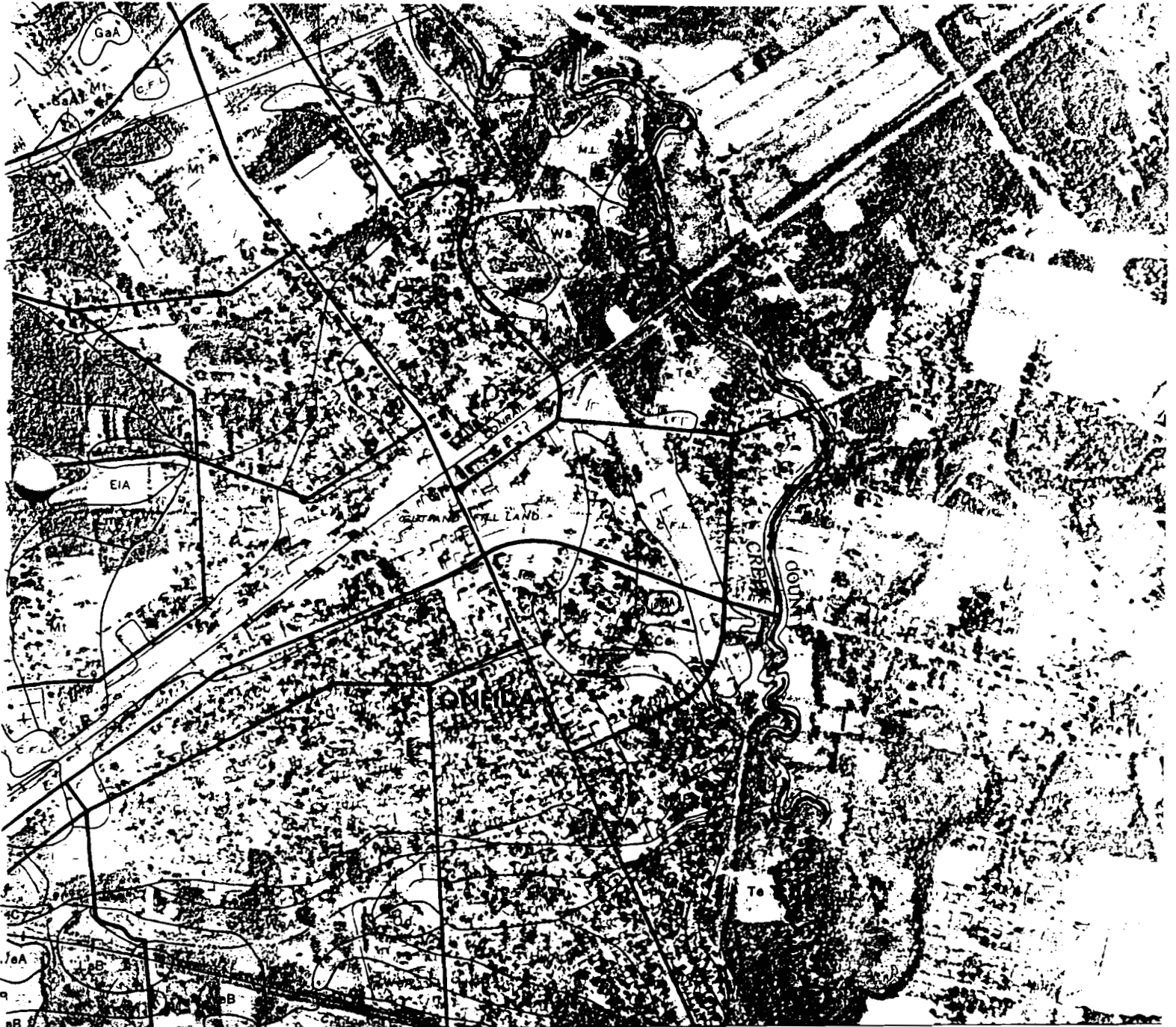
Hohman received his MA in Anthropology from the University of Connecticut in 1986 and his BA in Anthropology from the University of Rhode Island in 1983. He has worked in professional archaeology since 1983 and in 1987 joined the staff of the Public Archaeology Facility. From 1990 to 1991, Hohman served as one of the Coordinators at Garrow and Associates for the 300+ mile Iroquois Gas pipeline project in New York and Connecticut. He has served as project director on a wide range of site examinations and surveys and the Rainbow Plaza Data Recovery in Niagara Falls. His research interests include Northeast prehistory and 19th century rural and urban settlements. The author of numerous cultural resource management reports, he currently directs projects for PAF's statewide highway contract with the State Museum and other projects in New York State and Pennsylvania. Member of the Registry of Professional Archaeologists, Council for Northeast Historic Archaeology and the Archaeological Society of Connecticut.

ATTACHMENT B.1:  
Topographic Map with Project Area Noted

1955 Oneida Quadrangle 7.5' Series



**ATTACHMENT B.2:**  
**Soil Map of the Project Area**





**ATTACHMENT C:**  
**Site Files Checks**  
**(Confidential: Not for Public Release)**

**CONFIDENTIAL; Not for Public Release**  
**OPRHP/NYSM Site Files**  
7/6/00

em 1. Archaeological Sites. Niagara Mohawk Remediation Project, Oneida, Madison County. Oneida Quadrangle.

| SITE NUMBER                               | SITE NAME                                  | DIST. From Project/Dist. From water   | Elevation/Slope                                       | Cultural Affiliation/Site Type  | Testing/Type/Interval Artifacts | Report      |
|---|--|---|---|---|---------------------------------|-------------|
| 1) historic Native American<br>NYSM 3801  | ACP MDSN-19                                | General area 3.2 km (2 mi) W / vicinity of Cowaselon Creek                  | 131 m (430 ft) amsl; flat to gentle slope             | Historic Oneida village   | No information                  | Parker 1920 |
| 2) prehistoric<br>NYSM 4115               | ACP ONID-7                                 | General area 2.4 km (1.5 mi) N / vicinity of Oneida Creek                   | "   | Unidentified prehistoric Village  | "                               | "           |
| 3) historic Native American?<br>NYSM 4118 | ACP ONID-10                                | General area 0.5 km (0.3 mi) S / vicinity of Oneida Creek                   | 134 m (440 ft) amsl; flat to gentle slope             | Historic Native American? Cemetery  | "                               | "           |
| 4) prehistoric<br>NYSM 4116               | ACP ONID-8                                 | General area 1.9 km (1.2 mi) SE onto Vernon Quad / vicinity of Oneida Creek | 137 m (450 ft) amsl; flat to gentle slope             | Unidentified prehistoric Burial   | "                               | "           |
| 5) historic & historic NA<br>NYSM 4117    | ACP ONID-9A, B, C & D; Indian Church Site  | General area 2.4 km (1.5 mi) SE / vicinity of Oneida Creek                  | "   | Fort and block house (probably built under Wm. Johnson's direction, 1756) and village, church and cemetery (Native Americans) | "                               | "           |
| 6) historic NA<br>NYSM 3802               | Canowaraghere; Kanoalohaleo; Oneida Castle | General area 2.3 km (1.4 mi) SE / vicinity Oneida Creek, swamp              | "   | Historic Native American, 1762 Oneida village   | "                               | "           |
| 7) prehistoric<br>NYSM 7466               | ACP MDSN                                   | Lg area begins 2.7 km (1.7 mi) S / vicinity of trib of Oneida Creek         | 152-183 m (500-600 ft) amsl; gentle to moderate slope | Unidentified prehistoric Traces of occupation   | "                               | "           |
| 8) prehistoic<br>NYSM 3831                | ACP MDSN                                   | Lg area begins 2.4 km (1.5 mi) S / vicinity of tributary of Oneida Creek    | 152-213 m (500-700 ft) amsl; gentle to moderate slope | "   | "                               | "           |

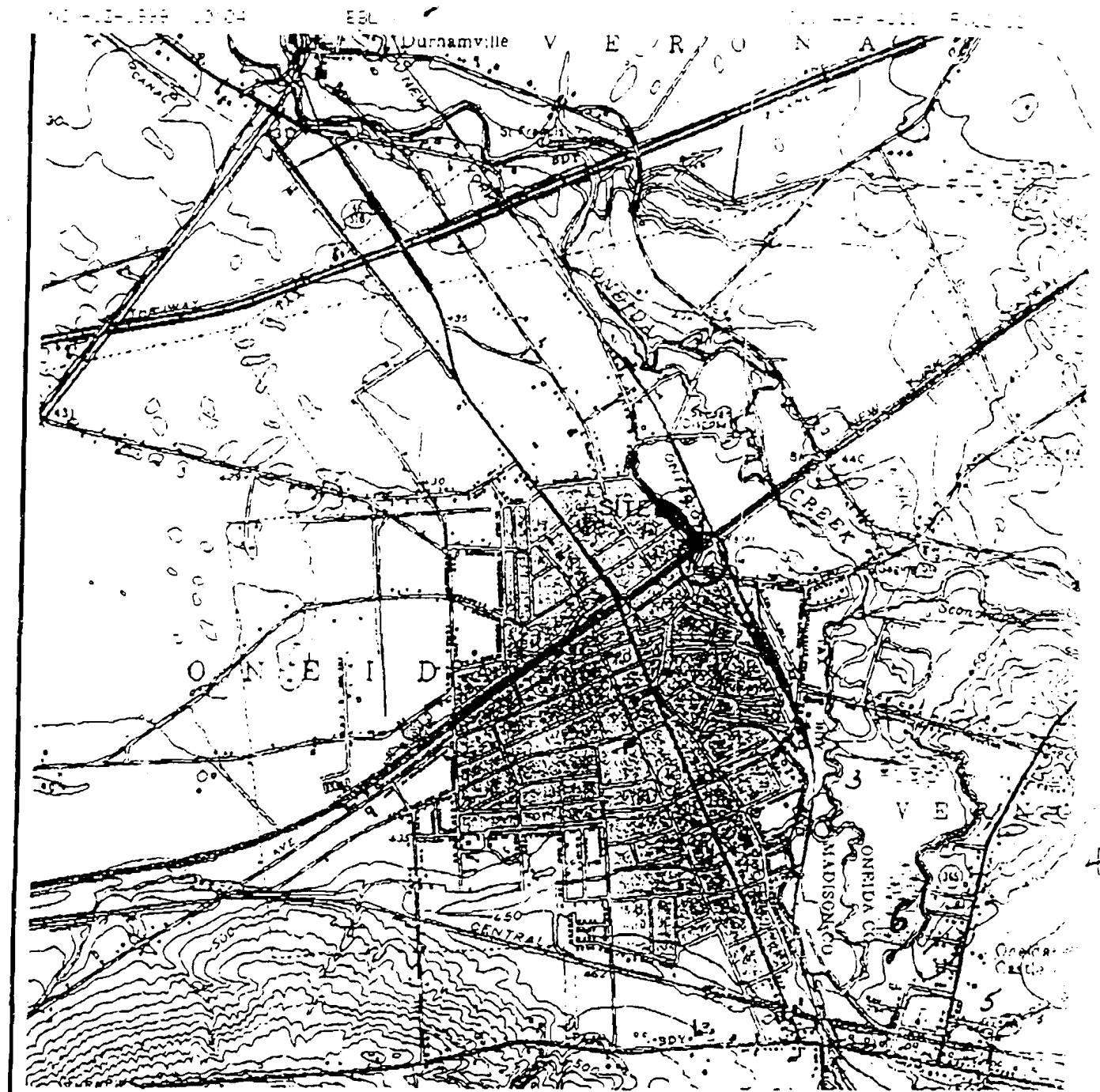
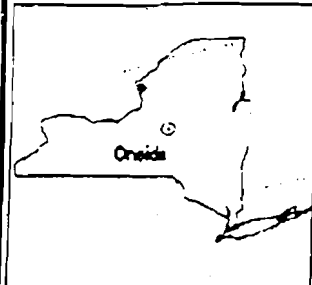


FIGURE 1.1

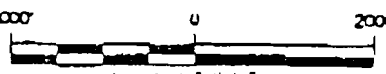


NEW YORK  
QUADRANGLE LOCATION

LATITUDE: N 43° 06' 27"  
LONGITUDE: W 75° 39' 08"



SOURCE: U.S.G.S. 7.5 MINUTE SERIES  
TOPOGRAPHIC MAP,  
ONEIDA, NEW YORK (1955)  
SW/4 ONEIDA 15 QUADRANGLE



Approximate Scale in Feet

**NIAGARA MOHAWK**

NIAGARA MOHAWK POWER  
CORPORATION  
SYRACUSE, NEW YORK

### SITE LOCATION MAP

ONEIDA SCENONDOA STREET  
143 CEDAR STREET, City Oneida, Madison  
0553

PARSONS ENGINEERING SCIENCE, INC.

DESIGN • RESEARCH • PLANNING  
ON EARTH'S SURFACE • WATER • LAND • AIR • AND  
SPACE • IN PRINCIPAL OFFICE

*Oneida  
0553*

**CONFIDENTIAL; Not for Public Release**  
**OFFICE OF PARKS, RECREATION AND HISTORIC PRESERVATION**  
**Field Services Bureau Files Search**

DATE: 7/6/00  
CONDUCTED BY: B. Ross

Project: Niagara Mohawk Remediation, 143 Cedar Street  
Minor Civil Division: city of Oneida (05340)  
County: Madison  
USGS Quadrangle: Oneida

1. Archaeological Sites (within 3.2 km/ 2 mi radius):

Refer to attached table.

2. Surveys and Reports within immediate or adjacent MCDs: (all; with possible exception of Stage IA):

OPR Report #3. *Stage IB Cultural Resource Survey for Proposed Oneida Wastewater Facilities, City of Oneida and Town of Lenox, Madison County, C-36-655, Pratt & Pratt, 5/76 for EPA. One prehistoric site; A05309.000005 and three with historic and prehistoric components; A05309.000003, 4 & 6 (all beyond 3.2 km from current project area). And a Stage IB Addendum, 1/77; no sites.*

OPR Report #10. *Cultural Resource Literature Search and Site Inventory for Oswego Basin, (multiple counties), Deborah Swartz, 10/80 for CORPS.*

OPR Report #38, 90PR1490. *Stage IA/B Cultural Resource Survey for Oneida City Hospital Nursing Home, Pratt and Pratt, 10/90 for HUD. No sites.*

OPR Report #41, 92PR2253. *Stage IA/B Archaeological Survey for Oneida Indian Nation Housing Phase I, City of Oneida, PAF, 11/92 for HUD. Two sites within 20 acres. Historic; A05340.000388 and prehistoric; A05340.000587 – both beyond 3.2 km from current project area. AND Stage II for SUBI-1405, A05340.000587, 3/93; 51 sq ft.*

OPR Report #43, 93PR1307. *Stage IA/B Cultural Resource Survey for Oneida Nation Sewer Project, Part II, Town of Oneida, PAF, 6/93 for EPA. No sites within two acres surveyed.*

OPR Report #47, 95PR0427. *Stage IA/B Cultural Resource Survey for Retail Establishment Site, City of Oneida, Pratt and Pratt, 1/95 for SEQRA. No sites within 33 acres surveyed.*

OPR Report #53, 97PR2466. *Cultural Resource Survey for Hartman Enterprises, Inc., 455 Elizabeth Street, City of Oneida, Hartgen Archaeological Services, 4/98 for SEQRA. No sites within 2.1 acres surveyed.*

3. NR Listed and NR Eligible Properties within, adjacent or within view shed of project area:

No NR eligible or listed properties within, adjacent or within view shed. There is a NRL historic district including many buildings on Main, Broad, Grove Streets as well as Wilbur and Walnut and several other individually eligible residences and commercial structures.

4. Inventoried Structures within, adjacent or within view shed of project area:

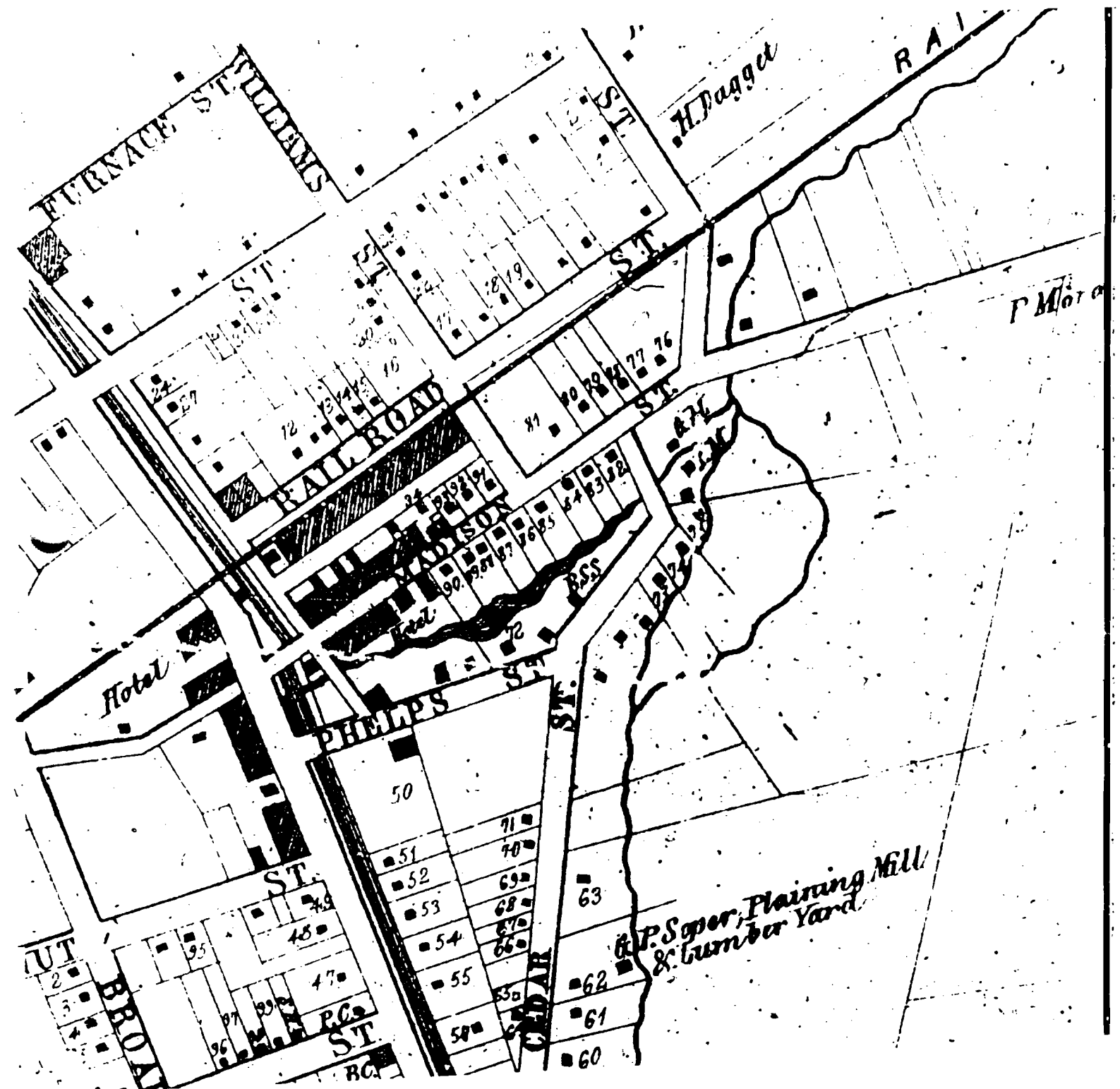
05340.000372, 232 Cedar Street; Cedar House-Fiore Building Apartments

**ATTACHMENT D: Historic Maps for the Project Area.**

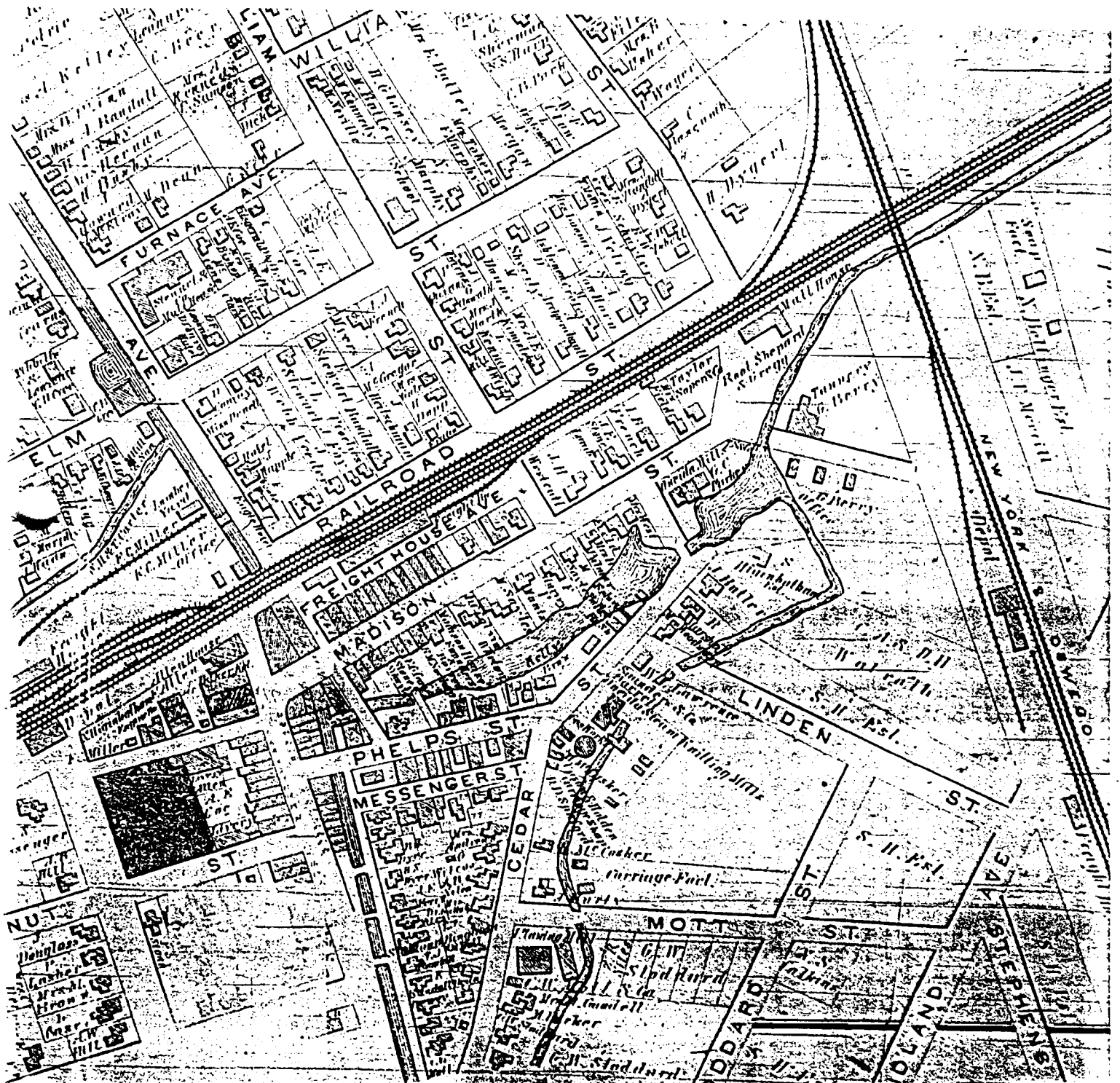
- D.1: 1853 Evans**
- D.2: 1859 French**
- D.3: 1875 Beers**
- D.4: 1890 Sanborn**
- D.5: 1895 Sanborn**
- D.6: 1899 Sanborn**
- D.7: 1904 Sanborn**
- D.8: 1909 Sanborn**
- D.9: 1914 Sanborn**
- D.10: 1923 Sanborn**
- D.11: 1930 Sanborn**
- D.12: 1930/1956 Sanborn**



D.2: 1859 French Map, with project area highlighted

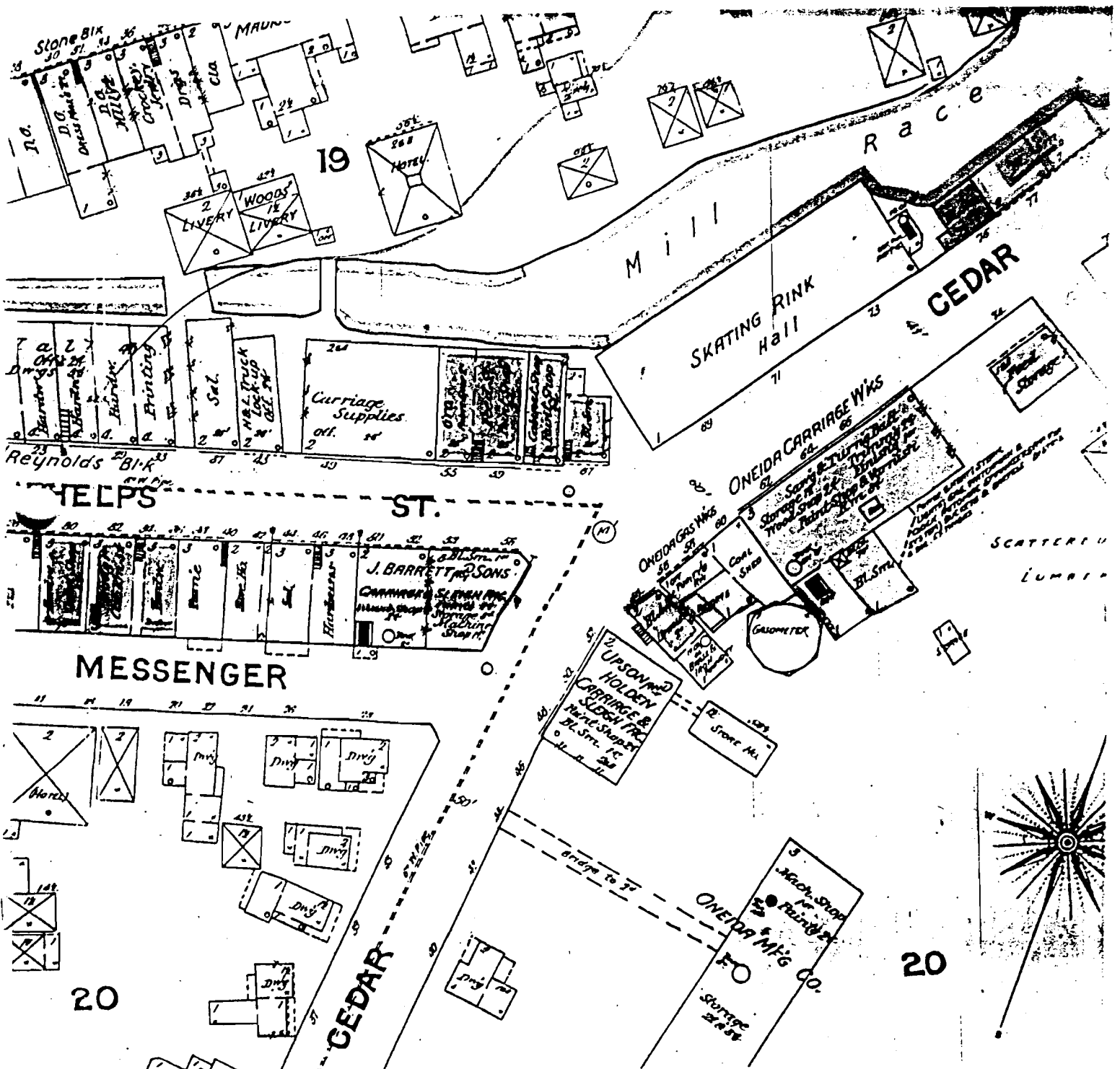


D.3: 1875 Beers Map, with project area highlighted

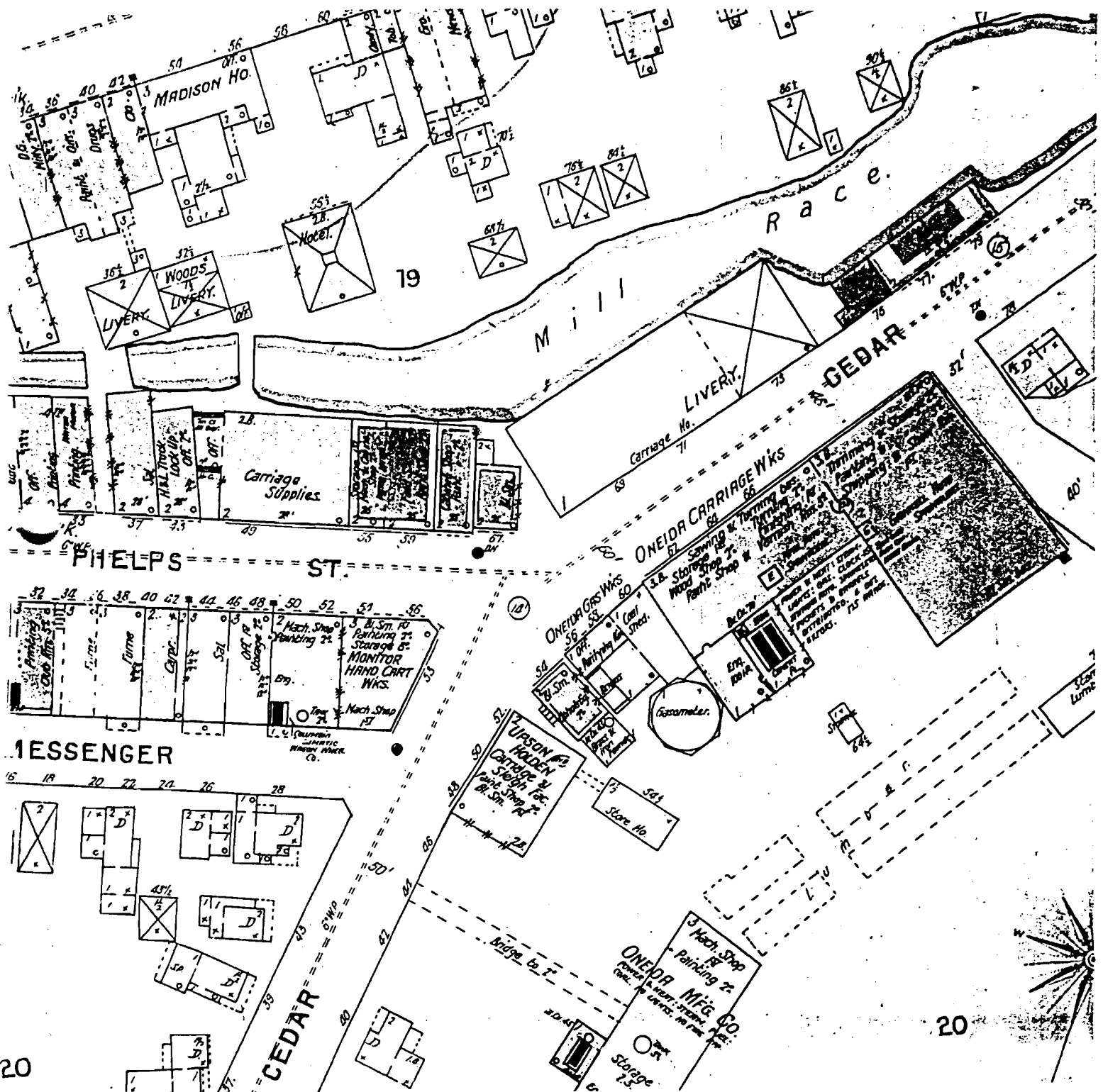




D.4: 1890 Sanborn map, with project area highlighted



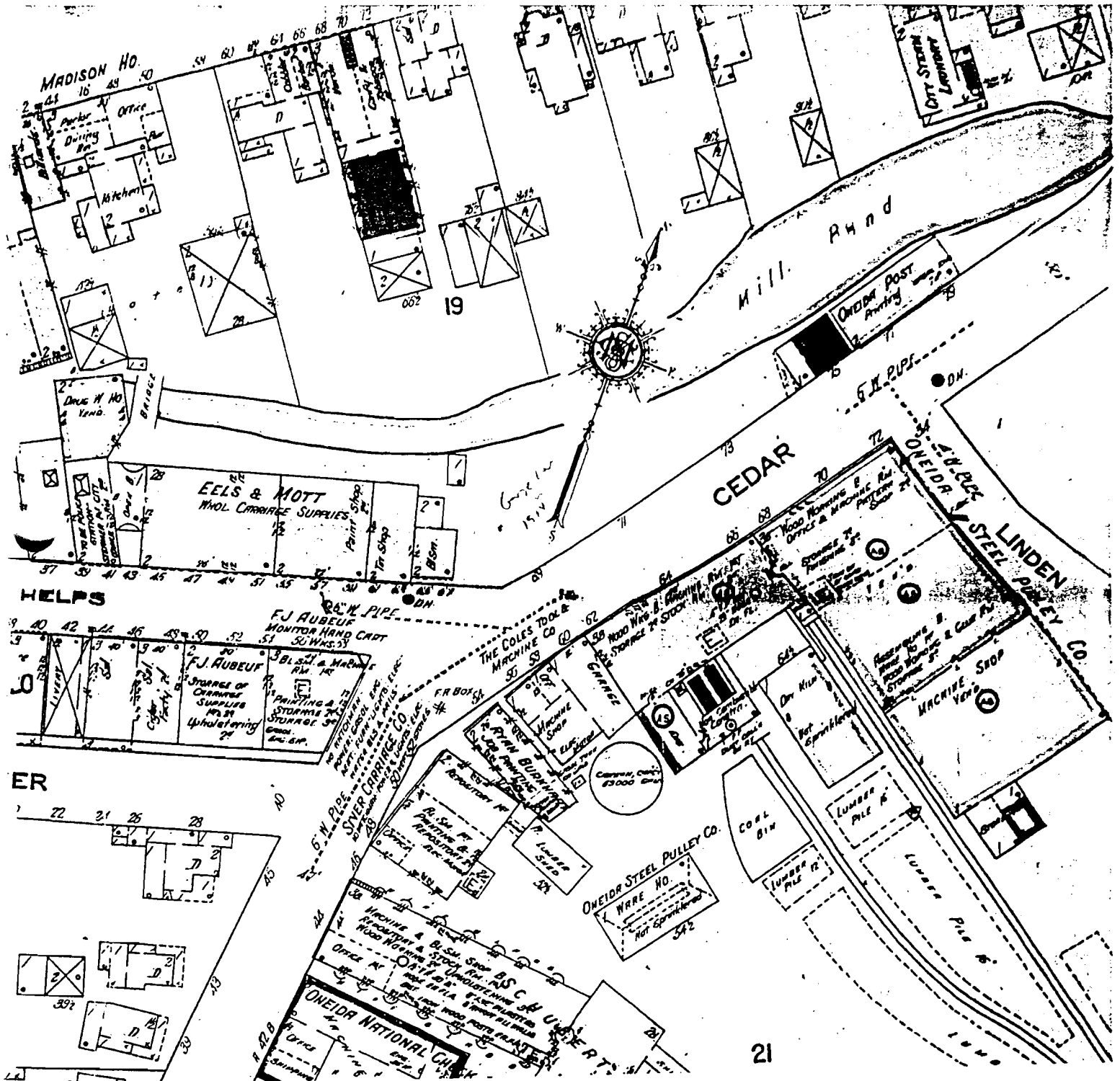
D.5: 1895 Sanborn map, with project area highlighted



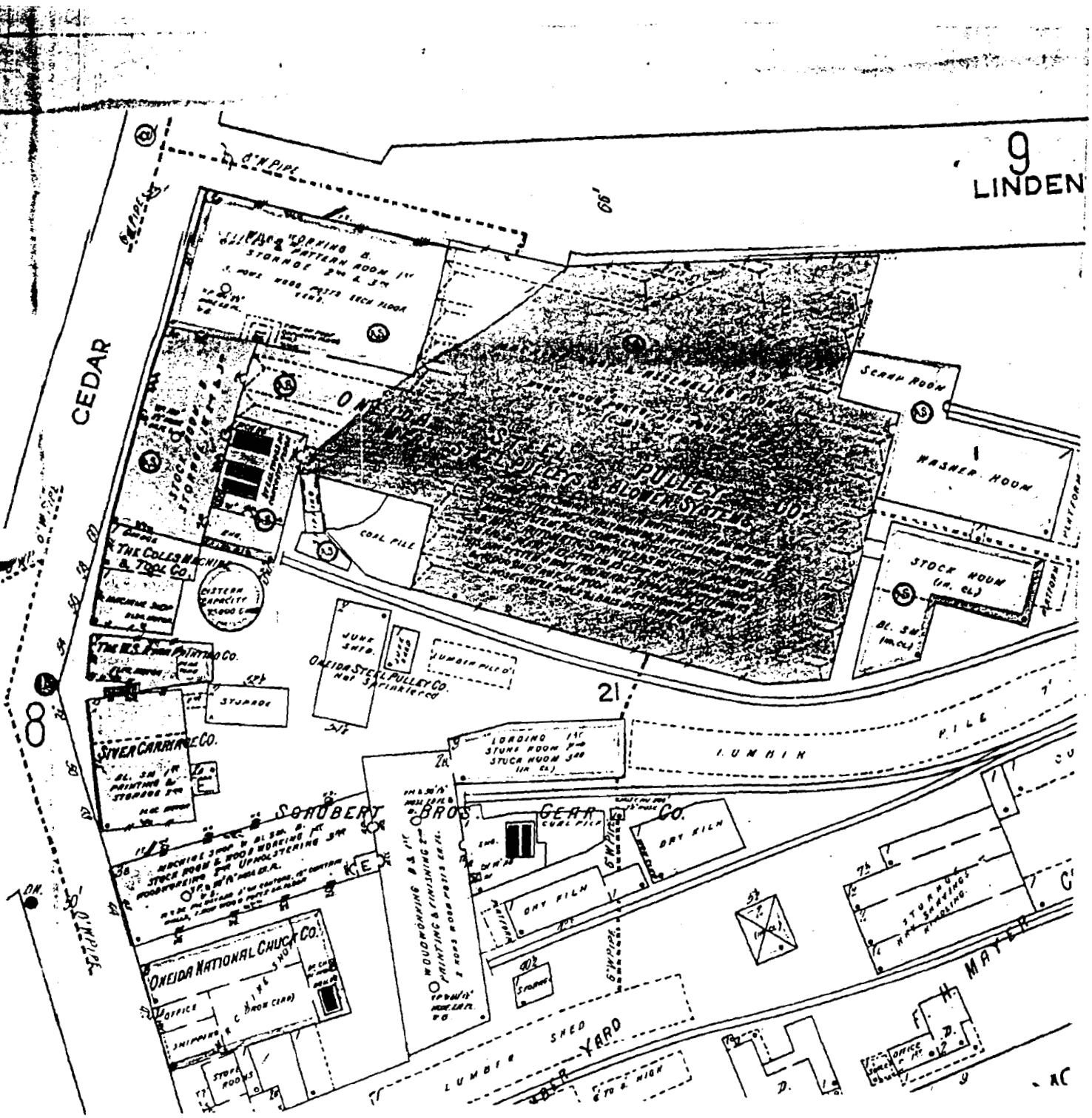


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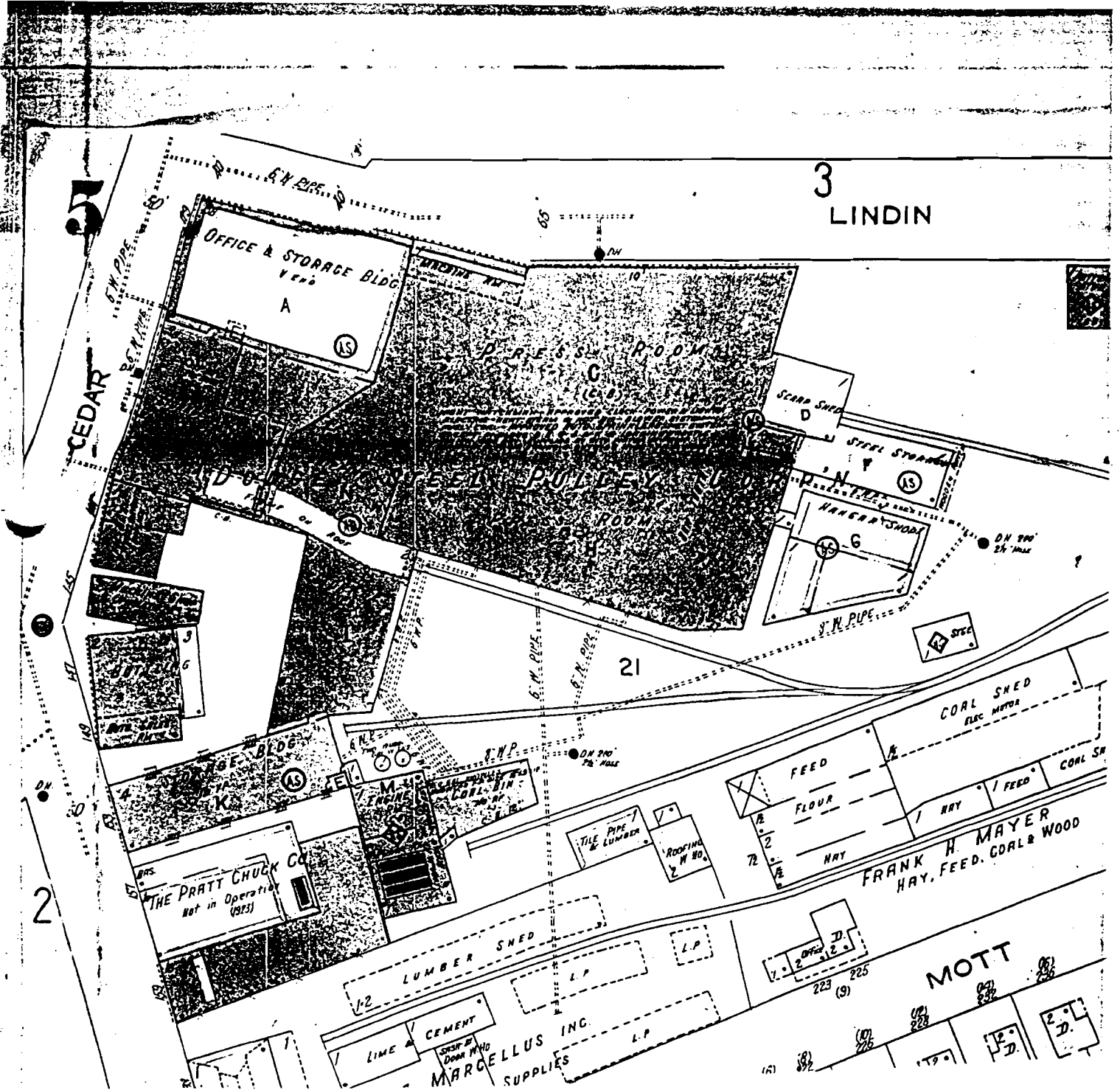
D.8: 1909 Sanborn map, with project area highlighted



D.9: 1914 Sanborn map, with project area highlighted

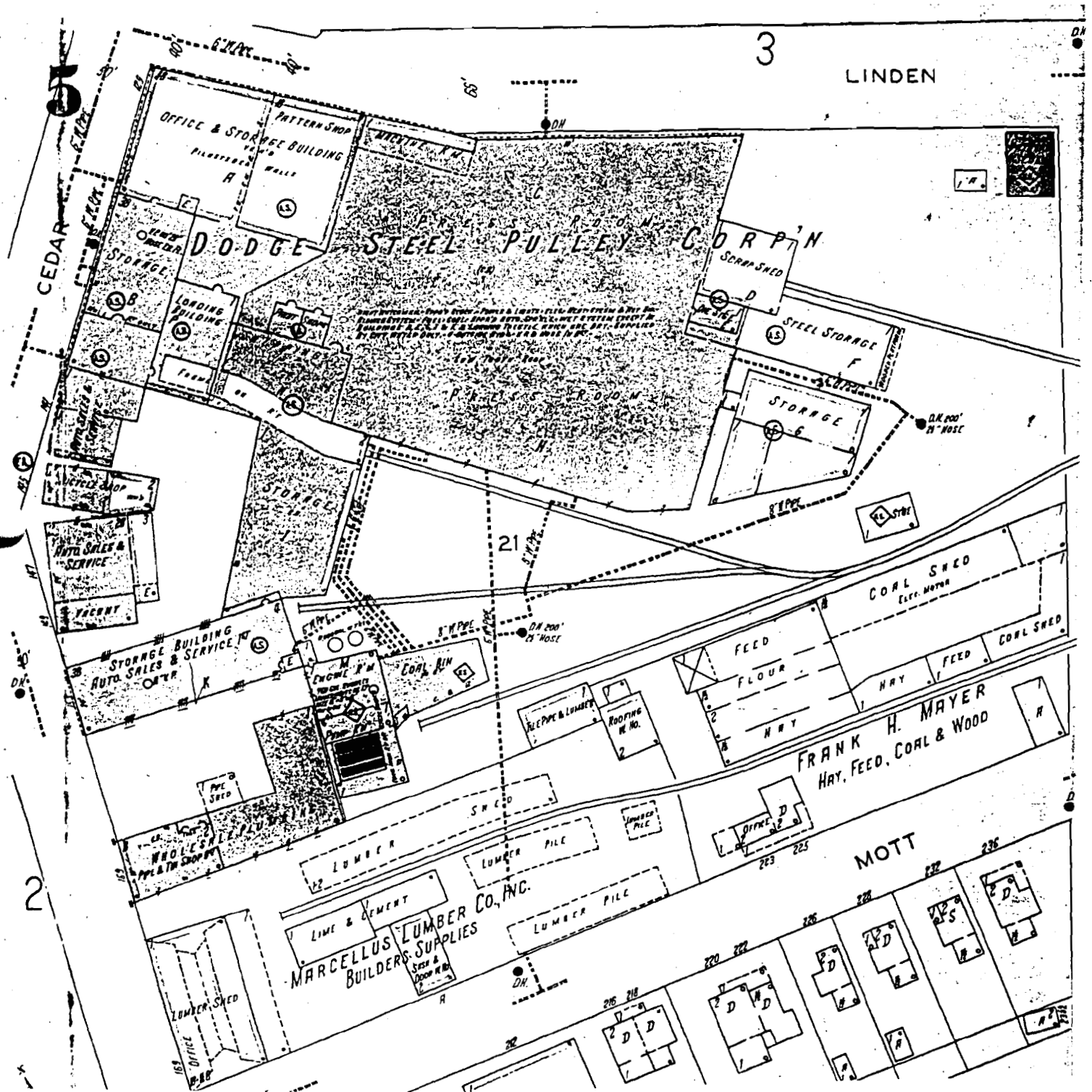


D.10: 1923 Sanborn map, with project area highlighted



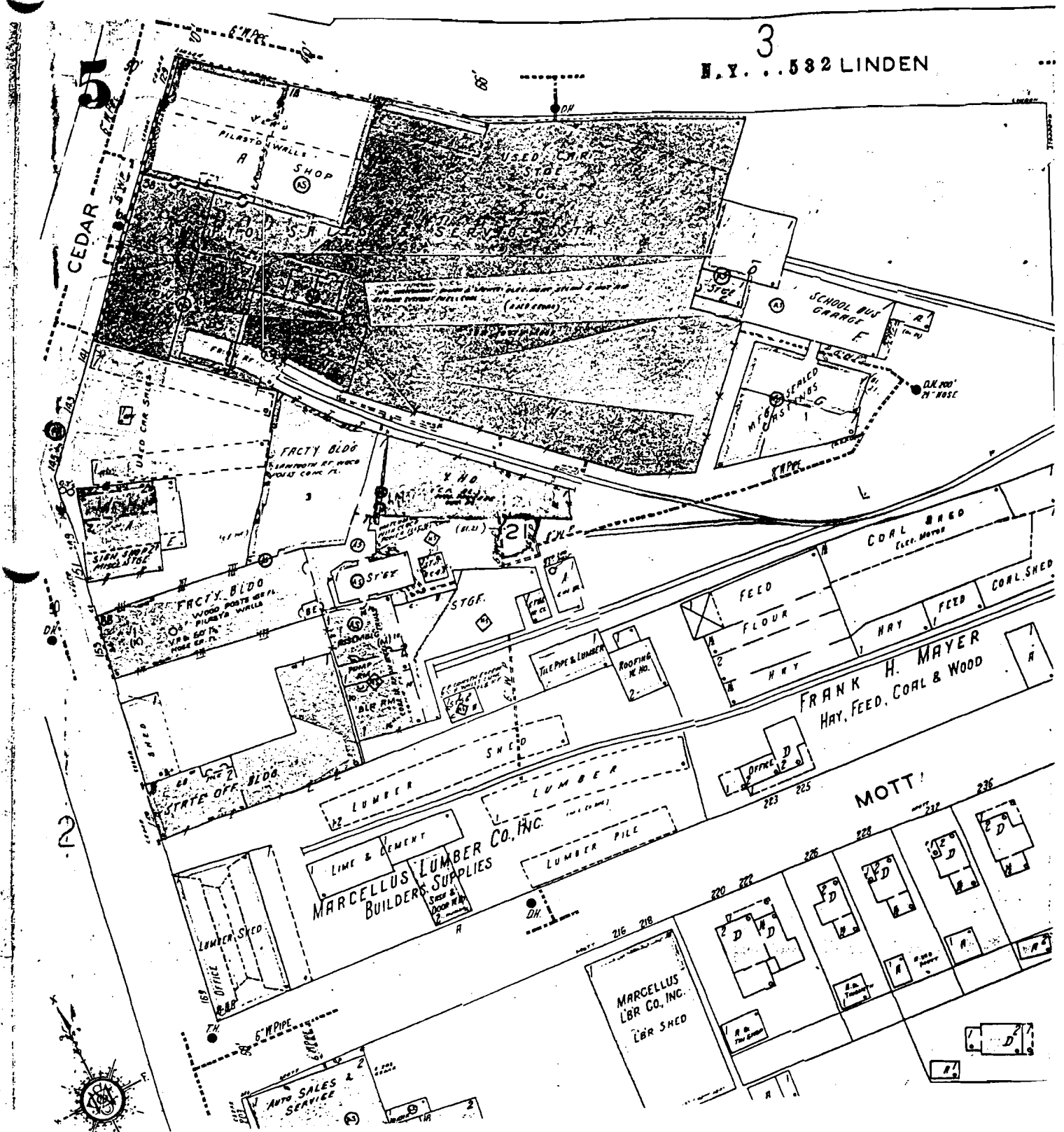


D.11: 1930 Sanborn map, with project area highlighted

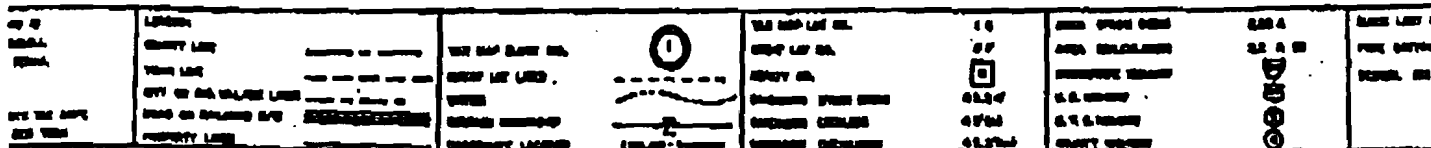




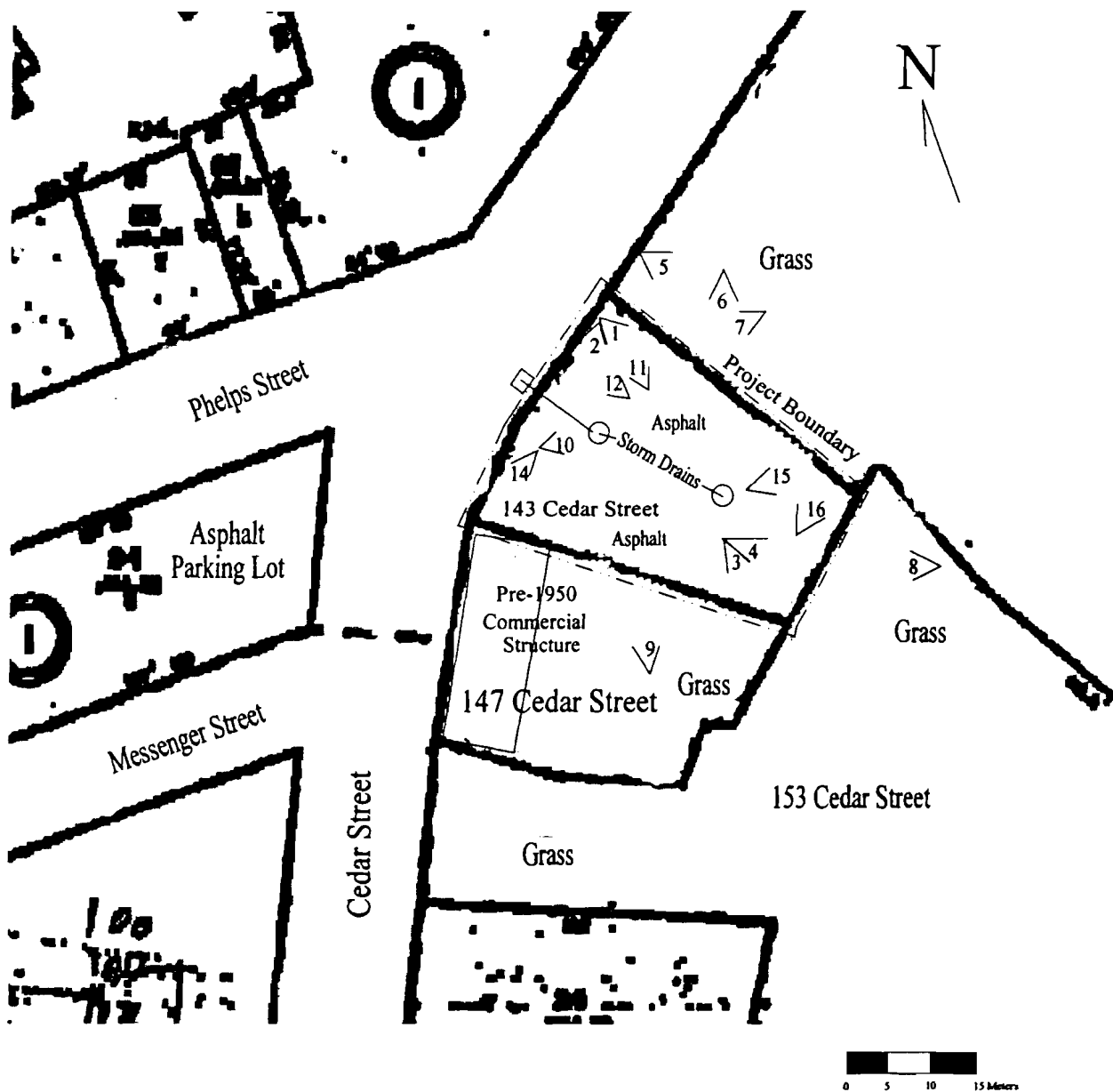
**D.12: 1930/1956 Sanborn map, with project area highlighted**



**ATTACHMENT E:**  
**Map of the project area showing testing locations and results**



City of Oneida Bond Act Applications



Stage 1A Cultural Resource Assessment  
Oneida MGP Project (143 Cedar St.)  
City of Oneida, Madison County, NY



**ATTACHMENT F:  
PROJECT AREA  
PHOTOGRAPHS**

Photo 1. Facing east, asphalt paving on Oneida MGP project area.



Photo 2. Facing south, asphalt paving on Oneida MGP project area with reconstructed 19<sup>th</sup> century structure (147 Cedar Street) to south.



Photo 3. Facing southeast, Oneida MGP project area above surrounding terrain to southeast.



Photo 4. Facing east, Oneida MGP project area above surrounding terrain to east.



Photo 5. Facing northeast, Oneida MGP project area above surrounding terrain to northeast.



Photo 6. Facing southeast, fieldstone and concrete retaining wall on north side of project area.



Photo 7. Facing southwest,  
fieldstone and concrete  
retaining wall on north side of  
project area.



Photo 8. Facing east, concrete  
retaining wall on east side of  
project area.



Photo 9. Facing northwest,  
concrete and fieldstone  
retaining wall on south side of  
project area.



Photo 10. Facing east, storm drain system located in middle of Oneida MGP lot.



Photo 11. Facing northwest, structures along Cedar Street.



Photo 12. Facing northwest, structures at intersection of Cedar Street and Messenger Street.





Photo 13. Facing west,  
structures along Messenger  
Street.



Photo 14. Facing west,  
parking lot at intersection of  
Messenger Street and Cedar  
Street.



Photo 15. Facing northeast,  
structures along Linden Street.



Photo 16. Facing north, view of structures along Linden Street from project area.

## **Attachment 3**

### **Subsurface Soil Information**

**Date Start/Finish:** 7/18/00  
**Drilling Company:** Nothnagle Drilling  
**Driller's Name:** Kevin Busch and Troy Bennett  
**Drilling Method:** Hollow Stem Auger  
**Bit Size:** 8.25-inch OD  
**Auger Size:** 4.25-inch ID  
**Rig Type:** BK-81  
**Sampling Method:** 2-inch split-spoons (SS)

**Northing:** 4931.7312  
**Easting:** 5005.7304  
**Casing Elevation:**  
**Borehole Depth:** 22 ft. bgs  
**Surface Elevation:** 436.3 ft. AMSL  
**Geologist:** Michael K. Cobb

**Well/Boring ID:** SB-1  
**Client:** Niagara Mohawk Power Corporation  
**Location:** 141 Cedar Street  
 Oneida, New York

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| Depth (feet) | Elevation (ft. AMSL) | Sample Run Number | Sample/In/Type | Recovery (feet) | Blows / 6 Inches    | N - Value | PID Headspace (ppm) | Analytical Sample | Geologic Column | Stratigraphic Description   | Hydrostratigraphy | Well Construction  |
|--------------|----------------------|-------------------|----------------|-----------------|---------------------|-----------|---------------------|-------------------|-----------------|---|-------------------|--|
| 0            |                      |                   |                |                 |                     |           |                     |                   |                 |   |                   |  |
| 435          |                      | 1                 | SS             | 0.5             | 4<br>6<br>4<br>5    | 10        | 0.4                 | X                 | X X X X         | Asphalt cover, 0-0.2' bgs. Brown SILT and fine SAND, trace Brick fragments, moist.  | Fill              |  |
|              |                      | 2                 | SS             | 1.2             | 2<br>9<br>19<br>9   | 28        | 109                 | X                 | X X X X         | Fine to medium SAND and brown SILT, fragments of Brick and Glass.<br>Timbers, black staining, slight odor, moist, Wood appears to be treated. |                   |  |
| -5           |                      | 3                 | SS             | 1.5             | 9<br>5<br>4<br>3    | 9         | 8.0                 | X                 | X X X X         | Medium red-brown fine SAND, moist.  | Silt and Sand     | Borehole tremie-grouted to grade with cement/bentonite slurry. |
| 430          |                      | 4                 | SS             | 1.7             | 3<br>3<br>4<br>5    | 7         | 20.7                |                   | X X X X         | Medium red-brown SILT and fine SAND.  |                   |  |
|              |                      | 5                 | SS             | 1.8             | 5<br>5<br>6<br>6    | 11        | 22.9                |                   | X X X X         | Medium red-brown fine SAND, little Silt.  |                   |  |
| -10          |                      | 6                 | SS             | 1.3             | 2<br>4<br>5<br>4    | 9         | 27.6                |                   | X X X X         | Medium red-brown fine to medium SAND, moist to wet.   |                   |  |
| 425          |                      | 7                 | SS             | 1.5             | 6<br>5<br>5<br>9    | 10        | 3.3                 |                   | X X X X         | Medium red-brown Silty fine SAND, moist to wet.   |                   |  |
|              |                      | 8                 | SS             | 2.0             | 6<br>16<br>20<br>21 | 36        | 6.0                 |                   | X X X X         | Medium red-brown Silty very fine SAND, medium stiff, moist to wet.  |                   |  |
| -15          |                      |                   |                |                 |                     |           |                     |                   |                 |   |                   |  |

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**Remarks:** ID/OD = inside/outside diameter ppm = parts per million  
 ags/bgs = above/below ground surface AMSL = above Mean Sea Level  
 PID = photoionization detector NA = not available

**Water Level Data**

**Date Depth Elev.**

Depth measured from top of casing.



**Well/Boring ID:** SB-2 **DRAFT**  
**Client:** Niagara Mohawk Power Corporation  
**Location:** 141 Cedar Street  
 Oneida, New York

**Location:** 141 Cedar Street  
Oneida, New York

Borehole tremie-grouted to grade with cement/bentonite slurry.

*Depth measured from top of casing.*



**Well/Boring ID:** SB-3  
**Client:** Niagara Mohawk Power Corporation  
**Location:** 141 Cedar Street  
Oneida, New York

**DRAFT**

| Depth (feet) | Elevation (ft. AMSL) | Sample Run Number | Sample/Int/Type | Recovery (feet) | Blows / 6 Inches      | N - Value | PID Headspace (ppm) | Analytical Sample | Geologic Column | Stratigraphic Description  | Hydrostratigraphy | Well Construction |
|--------------|----------------------|-------------------|-----------------|-----------------|-----------------------|-----------|---------------------|-------------------|-----------------|--|-------------------|-------------------|
| 0            |                      |                   |                 |                 |                       |           |                     |                   |                 |  |                   |                   |
| 435          |                      | 1                 | SS              | 1.2             | 4<br>6<br>6<br>4      | 12        | 2.8                 | X                 | x x x x x x x x | Fine to coarse SAND, little Silt, fine to medium Gravel, Brick fragments, loose, moist.  |                   |                   |
|              |                      | 2                 | SS              | 1.3             | 4<br>4<br>5<br>5      | 9         | 1.5                 | X                 | x x x x x x x x |  |                   |                   |
| -5           |                      | 3                 | SS              | 0.6             | 8<br>12<br>9<br>6     | 21        | 1.4                 | X                 | x x x x x x x x | Fine to coarse SAND and GRAVEL, little Silt , Brick fragments, Cobble lodged in spoon tip.   |                   |                   |
| 430          |                      | 4                 | SS              | 0.4             | 6<br>6<br>5<br>4      | 11        | 4.2                 | X                 | Brick pattern   | Shattered BRICK  |                   |                   |
|              |                      | 5                 | SS              | 0.8             | 3<br>4<br>8<br>4      | 12        | 3.4                 | X                 | Brick pattern   |  |                   |                   |
| -10          |                      | 6                 | SS              | 2.0             | 3<br>2<br>3<br>2      | 5         | 1.8                 | X                 | x x x x x x x x | Brown to black fine to coarse SAND, trace Cinders and Brick fragments, moist. In spoon tip, becomes Silty, little organics and gravel. |                   |                   |
| 425          |                      | 7                 | SS              | 2.0             | 2<br>2<br>1<br>1<br>2 | 3         | 2.1                 | X                 | Dotted pattern  | Medium red-brown fine SAND, trace Silt and Organic material, wet.  |                   |                   |
| -15          |                      | 8                 | SS              | 2.0             | 1<br>1<br>1           | 1         | 1.0                 |                   | Dotted pattern  | Medium red-brown fine SAND, little Silt, Clay, and Organic material, wet.  |                   |                   |
| 420          |                      |                   |                 |                 |                       |           |                     |                   |                 |  |                   |                   |

Fill

Silt & Sand

Borehole tremie-grouted to grade with cement/bentonite slurry.

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**Remarks:** ID/OD = inside/outside diameter    ppm = parts per million  
ags/bgs = above/below ground surface    AMSL = above Mean Sea Level  
PID = photoionization detector    NA = not available  
H = weight of hammer

### Water Level Data

| Date | Depth | Elev. |
|------|-------|-------|
|------|-------|-------|

*Depth measured from top of casing.*



|  |  |   |
|--|--|---|
| <b>Date Start/Finish:</b> 7/20/00<br><b>Drilling Company:</b> Nothnagle Drilling<br><b>Driller's Name:</b> Kevin Busch and Troy Bennett<br><b>Drilling Method:</b> Hollow Stem Auger<br><b>Bit Size:</b> 8.25-inch OD<br><b>Auger Size:</b> 4.25-inch ID<br><b>Rig Type:</b> BK-81<br><b>Sampling Method:</b> 2 and 3-inch split-spoons (SS) | <b>Northing:</b> 4938.0376<br><b>Easting:</b> 5071.469<br><b>Casing Elevation:</b><br><br><b>Borehole Depth:</b> 17 ft. bgs<br><b>Surface Elevation:</b> 436.2 ft. AMSL<br><br><b>Geologist:</b> Michael K. Cobb | <b>Well/Boring ID:</b> SB-4 / SB-4A<br><b>Client:</b> Niagara Mohawk Power Corporation<br><br><b>Location:</b> 141 Cedar Street<br>Oneida, New York |
|--|--|---|

# DRAFT

| Depth (feet) | Elevation (ft. AMSL) | Sample Run Number | Sample/Int/Type | Recovery (feet) | Blows / 6 Inches | N - Value | PID Headspace (ppm) | Analytical Sample | Geologic Column | Stratigraphic Description   | Hydrostratigraphy | Well Construction |
|--------------|----------------------|-------------------|-----------------|-----------------|------------------|-----------|---------------------|-------------------|-----------------|---|-------------------|-------------------|
| 0            |                      |                   |                 |                 |                  |           |                     |                   |                 |   |                   |                   |
| 435          |                      | 1                 | SS              | 0.9             | 4 ref            | NA        | 3.2                 |                   | x x             | SB-4: Brown fine to coarse SAND and fine GRAVEL, little medium to coarse Gravel, little Silt, loose, dry to moist.<br>Auger refusal. Offset 2' and auger to 2' bgs without sampling.          |                   |                   |
|              |                      |                   |                 |                 | 6                |           |                     |                   | x x             |   |                   |                   |
|              |                      | 2                 | SS              | 1.2             | 11               | 21        | 1.2                 |                   | x x x           | Medium brown fine to medium SAND, some coarse Sand and fine Gravel, trace Brick, moist.   |                   |                   |
|              |                      |                   |                 |                 | 10               |           |                     |                   | x x x           |   |                   |                   |
|              |                      |                   |                 |                 | 15               |           |                     |                   | x x x           |   |                   |                   |
| 5            |                      | 3                 | SS              | 1.3             | 7                |           |                     |                   | x x x           |   |                   |                   |
|              |                      |                   |                 |                 | 15               | 35        | 2.8                 |                   | x x x           |   |                   |                   |
|              |                      |                   |                 |                 | 20               |           |                     |                   | x x x           |   |                   |                   |
| 430          |                      |                   |                 |                 | 30               |           |                     |                   | x x x           |   |                   |                   |
|              |                      | NA                | SS              | 1.3             | 16               | NA        | 1.3                 |                   |                 | 6-7.6' bgs, fine to coarse SAND and GRAVEL, trace broken Cobbles, moist (FILL). Spoon and auger refusal at 7.6' bgs. Offset 15 feet south east to SB-4A; auger to 7' bgs and resume sampling. |                   |                   |
|              |                      |                   |                 |                 | 38               |           |                     |                   |                 |   |                   |                   |
|              |                      | 4                 | SS              | 1.5             | 1                |           |                     |                   |                 | SB-4A: Fine SAND, little Silt, trace fine Gravel and Brick, moist.  |                   |                   |
|              |                      |                   |                 |                 | 2                | 5         | 1.8                 |                   |                 |   |                   |                   |
|              |                      |                   |                 |                 | 3                |           |                     |                   |                 |   |                   |                   |
|              |                      |                   |                 |                 | 2                |           |                     |                   |                 |   |                   |                   |
| 10           |                      | 5                 | SS              | 2.0             | 3                | 5         | 4.6                 |                   |                 |   |                   |                   |
|              |                      |                   |                 |                 | 3                |           |                     |                   |                 |   |                   |                   |
| 425          |                      |                   |                 |                 | 3                |           |                     |                   |                 | Red-brown fine Silty SAND, trace Organic Material, moist.   |                   |                   |
|              |                      | 6                 | SS              | 2.0             | 3                | 5         | 3.4                 |                   |                 |   |                   |                   |
|              |                      |                   |                 |                 | 2                |           |                     |                   |                 | Black organic rich SILT, moist.   |                   |                   |
|              |                      |                   |                 |                 | 3                |           |                     |                   |                 | Medium red-brown Silty fine SAND, trace Organics, trace black staining, moist to wet.   |                   |                   |
|              |                      | 7                 | SS              | 2.0             | 4                | 8         | 1.3                 |                   |                 |   |                   |                   |
|              |                      |                   |                 |                 | 4                |           |                     |                   |                 |   |                   |                   |
|              |                      |                   |                 |                 | 4                |           |                     |                   |                 |   |                   |                   |
| 15           |                      |                   |                 |                 | 6                |           |                     |                   |                 | Medium brown SILT and very fine SAND, stiff, wet.   |                   |                   |
|              |                      | 8                 |                 | 2.0             | 6                | 15        | 3.1                 |                   |                 |   |                   |                   |
|              |                      |                   |                 |                 | 7                |           |                     |                   |                 |   |                   |                   |

Borehole tremie-grouted to grade with cement/bentonite slurry.



**Remarks:** ID/OD = inside/outside diameter    ppm = parts per million  
 ags/bgs = above/below ground surface    AMSL = above Mean Sea Level  
 PID = photoionization detector    NA = not available  
 ref = spoon refusal  
 Sampling details 0-7' bgs from SB-4; 7-17' bgs from SB-4A.

| Water Level Data |       |       |
|------------------|-------|-------|
| Date             | Depth | Elev. |
|                  |       |       |
|                  |       |       |
|                  |       |       |
|                  |       |       |
|                  |       |       |
|                  |       |       |
|                  |       |       |

Depth measured from top of casing.

## Client:

Niagara Mohawk Power Corporation

Well/Boring ID: SB-4 / SB-4A

**DRAFT**

## Site Location:

141 Cedar Street  
Oneida, New York

Borehole Depth: 17 ft. bgs

| Depth (feet) | Elevation (ft. AMSL) | Sample Run Number | Sample/Int/Type | Recovery (feet) | Blows / 6 Inches | N - Value | PID Headspace (ppm) | Analytical Sample | Geologic Column | Stratigraphic Description | Hydrostratigraphy | Well Construction  |
|--------------|----------------------|-------------------|-----------------|-----------------|------------------|-----------|---------------------|-------------------|-----------------|---------------------------|-------------------|--|
| 420          |                      | SS                | SS              | 9               | 12               |           | 3.1                 |                   |                 | As above.                 | Sand              | Borehole tremie-grouted to grade with cement/bentonite slurry. |
| 20           |                      |                   |                 |                 |                  |           |                     |                   |                 |                           |                   |  |
| 415          |                      |                   |                 |                 |                  |           |                     |                   |                 |                           |                   |  |
| 25           |                      |                   |                 |                 |                  |           |                     |                   |                 |                           |                   |  |
| 410          |                      |                   |                 |                 |                  |           |                     |                   |                 |                           |                   |  |
| 30           |                      |                   |                 |                 |                  |           |                     |                   |                 |                           |                   |  |
| 405          |                      |                   |                 |                 |                  |           |                     |                   |                 |                           |                   |  |
| 35           |                      |                   |                 |                 |                  |           |                     |                   |                 |                           |                   |  |

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**Remarks:** ID/OD = inside/outside diameter ppm = parts per million  
 ags/bgs = above/below ground surface AMSL = above Mean Sea Level  
 PID = photoionization detector NA = not available  
 ref = spoon refusal  
 Sampling details 0-7' bgs from SB-4; 7-17' bgs from SB-4A.

## Water Level Data

| Date | Depth | Elev. |
|------|-------|-------|
|      |       |       |
|      |       |       |
|      |       |       |
|      |       |       |

Depth measured from top of casing.

# DRAFT

*Depth measured from top of casing*

**Date Start/Finish:** 8/8/00  
**Drilling Company:** Nothnagle Drilling  
**Driller's Name:** Kevin Busch and Troy Bennett  
**Drilling Method:** Hollow Stem Auger  
**Bit Size:** 8 1/4-inch OD  
**Auger Size:** 4 1/4-inch ID  
**Rig Type:** BK-81  
**Sampling Method:** 2-inch split-spoons (SS)

**Northing:** 4905.297  
**Easting:** 5062.486  
**Casing Elevation:**  
  
**Borehole Depth:** 20 ft. bgs  
**Surface Elevation:** 435.8 ft. AMSL  
  
**Geologist:** Jason C. Sents

**Well/Boring ID:** SB-6  
**Client:** Niagara Mohawk Power Corporation  
  
**Location:** 141 Cedar Street  
 Oneida, New York

**DRAFT**

| Depth (feet) | Elevation (ft. AMSL) | Sample Run Number | Sample Int/Type | Recovery (feet) | Blows / 6 Inches  | N - Value | PID Headspace (ppm) | Analytical Sample | Geologic Column | Stratigraphic Description  | Hydrostratigraphy | Well Construction  |
|--------------|----------------------|-------------------|-----------------|-----------------|-------------------|-----------|---------------------|-------------------|-----------------|--|-------------------|--|
| 0            |                      |                   |                 |                 |                   |           |                     |                   |                 |  |                   |  |
| 435          |                      | 1                 | SS              | 2.0             | 8<br>10<br>8<br>9 | 18        | 2.4                 |                   |                 | x x Topsoil, grass and roots.<br>x x CLAY, CRUSHED STONE, BRICK, BLACK ASH, and CINDERS, damp.<br>x x CRUSHED CONCRETE, damp.<br>x x Black fine SAND, SILT, CINDERS, ASH, and CRUSHED BRICK, damp.<br>x x CRUSHED RED BRICK, damp.<br>x x Brown fine to medium SAND, SILT with crushed Sandstone Cobbles and red Brick, black staining, dry.<br>x x Dark brown to black fine to medium SAND and CLAY, with Cinders and Ash, wet. Broken Cobble at 4.3' bgs. Spoon refusal at 5' bgs. |                   |  |
| 5            |                      | 2                 | SS              | 1.5             | 7<br>7<br>5<br>5  | 12        | 2.7                 |                   |                 |  |                   |  |
| 430          |                      | 3                 | SS              | 0.4             | 8<br>ref          | NA        | 1.8                 |                   |                 |  |                   |  |
|              |                      | 4                 | SS              | 2.0             | 3<br>4<br>8<br>5  | 12        | 3.2                 |                   |                 | x x Dark red-brown fine to medium SAND, SILT, CRUSHED BRICK, SLAG, WOOD, CINDERS, ASH and COAL, trace roofing shingle, damp.<br>x x<br>x x<br>x x<br>x x   |                   |  |
|              |                      | 5                 | SS              | 1.0             | 4<br>1<br>2<br>7  | 3         | 2.9                 |                   |                 | x x Black CINDERS, ASH, COAL, and CINDER BLOCK, trace roofing shingle, wet.<br>x x<br>x x<br>x x<br>x x  |                   |  |
| 10           |                      | 6                 | SS              | 2.0             | 2<br>4<br>5<br>3  | 9         | 4.1                 |                   |                 | x x Dark brown fine SAND and SILT, some black Ash, Cinders, and Coal, trace white/tan Cinder Block, soft, wet.<br>x x<br>x x<br>x x  |                   |  |
|              |                      | 7                 | SS              | 2.0             | 2<br>2<br>3<br>4  | 5         | 9.5                 |                   |                 | Brown CLAY, trace fine Sand, soft, plastic, wet.<br>Brown fine SAND and CLAY, trace medium to coarse Sand, trace Gravel.<br>Dark red WOOD (possibly cedar).<br>Dark brown PEAT and WOOD, roots, cocoa-like odor, damp.<br>Brown CLAY, stiff, wet. Fine to medium Sand seams at 14.4, 14.8, 15.6 and 15.8' bgs.   | Fill<br>Peat/Clay | Borehole tremie-grouted to grade with cement/bentonite slurry. |
| 15           |                      | 8                 | SS              | 2.0             | 3<br>4<br>3<br>4  | 7         | 7.2                 |                   |                 |  |                   |  |
| 420          |                      |                   |                 |                 |                   |           |                     |                   |                 |  |                   |  |

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**Remarks:** ID/OD = inside/outside diameter ppm = parts per million  
 ags/bgs = above/below ground surface AMSL = above Mean Sea Level  
 PID = photoionization detector NA = not available  
 ref = spoon refusal

**Water Level Data**

**Date Depth Elev.**

Depth measured from top of casing.

**DRAFT**

**Borehole Depth:** 20 ft. bgs

[illegible]

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**Remarks:** ID/OD = inside/outside diameter      ppm = parts per million  
 ags/bgs = above/below ground surface      AMSL = above Mean Sea Level  
 PID = photoionization detector      NA = not available  
 ref = spoon refusal

| Water Level Data |       |       |
|------------------|-------|-------|
| Date             | Depth | Elev. |
|                  |       |       |
|                  |       |       |
|                  |       |       |

Depth measured from top of casing.

**NIAGARA MOHAWK POWER CORPORATION  
ONEIDA (141 CEDAR STREET) FORMER MGP SITE  
SITE INVESTIGATION ACTIVITIES**

**OFFSITE BORING SUMMARY**

| Boring ID | Sample Description  | Sample Interval | Headspace (ppm) |
|-----------|---|-----------------|-----------------|
| SB-7      | 0.0-0.7' - Tan brown fine SAND, little medium to coarse Sand, fine Gravel, loose, organics.   | (0-4')          | 0.0             |
|           | 0.7-1.2' - Medium to coarse SAND and fine GRAVEL, [FILL], brick fragments, loose, glass fragments.  |                 |                 |
|           | 1.2-4.0' - No recovery.   |                 |                 |
|           | 4.0-8.0' - No recovery.   | (4-8')          | NA              |
| SB-8      | 0.0-1.8' - Light brown fine to medium SAND and roots.   | (0-4')          | 0.0             |
|           | 1.8-4.0' - No recovery.   |                 |                 |
|           | 4.0-5.2' - Light brown fine to coarse SAND, some Gravel and organics; black tar throughout sample.  | (4-6')          | 206             |
|           | 5.2-8.0' - Olive-brown to dark brown PEAT, little black staining.   | (6-8')          | 25.2            |
| SB-9      | 0.0-1.0' - Medium reddish-brown fine to medium SAND, little coarse Sand, fine Gravel.   | (0-2')          | 0.9             |
|           | 1.0-2.1' - Loose fine to coarse SAND, fine Gravel, black, gray, yellow, and red Fill, trace Slag and Brick fragments.   | (2-4')          | 2.1             |
|           | 2.1-3.5' - Medium brown fine SAND, some Silt, little medium to coarse Sand, fine Gravel, Brick, and Cinders [FILL], moist to wet.                                   |                 |                 |
|           | 3.5-4.0' - No recovery.   |                 |                 |
|           | 4.0-5.0' - Medium to light brown fine SAND and SILT, little Clay, trace coarse Sand, stiff, wet.  | (4-6')          | 0.0             |
|           | 5.0'-7.0' - Black Peaty SILT, grading to dark brown-black PEAT.   |                 |                 |
| SB-10     | 7.0-8.0' - No recovery.   | (6-8)           | 1.0             |
|           | 0.0-4.0' - Red-brown fine to coarse SAND, organics, and broken concrete [FILL], black staining.   | (0-4')          | 2.5             |
|           | 4.0-5.4' - Red-brown to tan fine to medium SAND, trace to little Silt and Gravel.   | (4-7')          | 0.0             |
|           | 5.4-8.0' - Olive-gray to dark brown PEAT, black staining.   | (7-8')          | 0.0             |
| SB-11     | 0.0-4.0' - Red-brown fine to coarse SAND, organics, and slag [FILL].  | (0-4')          | 0.0             |
|           | 4.0-5.0' - No recovery in sleeve; black tar on gravel found in shoe. Refusal at 5' bgs.   | (4-5')          | NA              |
| SB-12     | 0.0-3.5' - Red-brown fine to medium SAND, organics, and little Gravel; trace oily sheen; black staining, and odor. Possible tar at 3-3.5' bgs. Refusal at 3.5' bgs. | (0-3')          | 0.0             |
|           |   | (3-3.5')        | 38.7            |
| SB-13     | 0.0-4.0' - Brown fine to medium SAND, organics, slag, and brick [FILL]; yellow staining.  | (0-3')          | 0.4             |
|           |   | (3-4')          | 0.0             |
|           | 4.0-7.2' - Olive/yellow-brown to red FILL, fine to coarse Sand; black staining.   | (4-6')          | 0.0             |
|           | 7.2-8.0' - Dark brown to black PEAT.  | (6-8')          | 0.0             |
| SB-14     | 0.0-1.8' - Medium brown fine to medium SAND, trace coarse Sand and fine Gravel, loose [FILL].   | (0-2')          | 0.8             |
|           | 1.8-2.5' - Coarse SAND and fine GRAVEL, [FILL], brick fragments, slag, wet, yellowish brown, red brick.   | (2-4')          | 0.0             |
|           | 2.5-4.0' - No recovery.   |                 |                 |
|           | 4.0-5.8' - Medium brown fine SAND, some Silt, wet, soft, trace medium to coarse Sand, possible brick fragments.   | (4-6')          | 4.3             |
|           | 5.8-6.2' - Dark brown-black Peaty SILT, little black wood fragments, possible faint odor.   | (6-8')          | 28.2            |
|           | 6.2-7.5' - Dark brown-black PEAT, low density, moist.   |                 |                 |
| SB-15     | 7.5-8.0' - No recovery.   |                 |                 |
|           | 0.0-1.3' - Medium brown fine to medium SAND, trace coarse Sand, loose, moist.   | (0-2')          | 0.0             |
|           | 1.3-2.2' - Fine to coarse SAND, trace Gravel, red Brick, black-gray Fill, possible Slag, broken rock, trace organics, wet.  | (2-4')          | 0.0             |
|           | 2.2-4.0' - No recovery.   |                 |                 |
|           | 4.0-5.3' - Medium brown fine SAND, little Silt and medium Sand, trace coarse Sand and fine Gravel, possible brick fragments, wet.                                   | (4-6')          | 5.5             |
|           | 5.3-7.6' - Medium brown-black PEAT, low density, moist.   | (6-8')          | 0.0             |
|           | 7.6-8.0' - No recovery.   |                 |                 |

**Notes:**

1. Headspace screening measurements of the relative concentration of volatile organic compounds in soil in ppm (parts per million) obtained with a MiniRae(TM) photoionization detector.
2. Soil borings completed with a direct-push rig.
3. Sample Interval - Indicates the depth interval (in feet below ground surface) penetrated by the direct push sampling device. The intervals indicated in the soil description represent actual soil recovery.

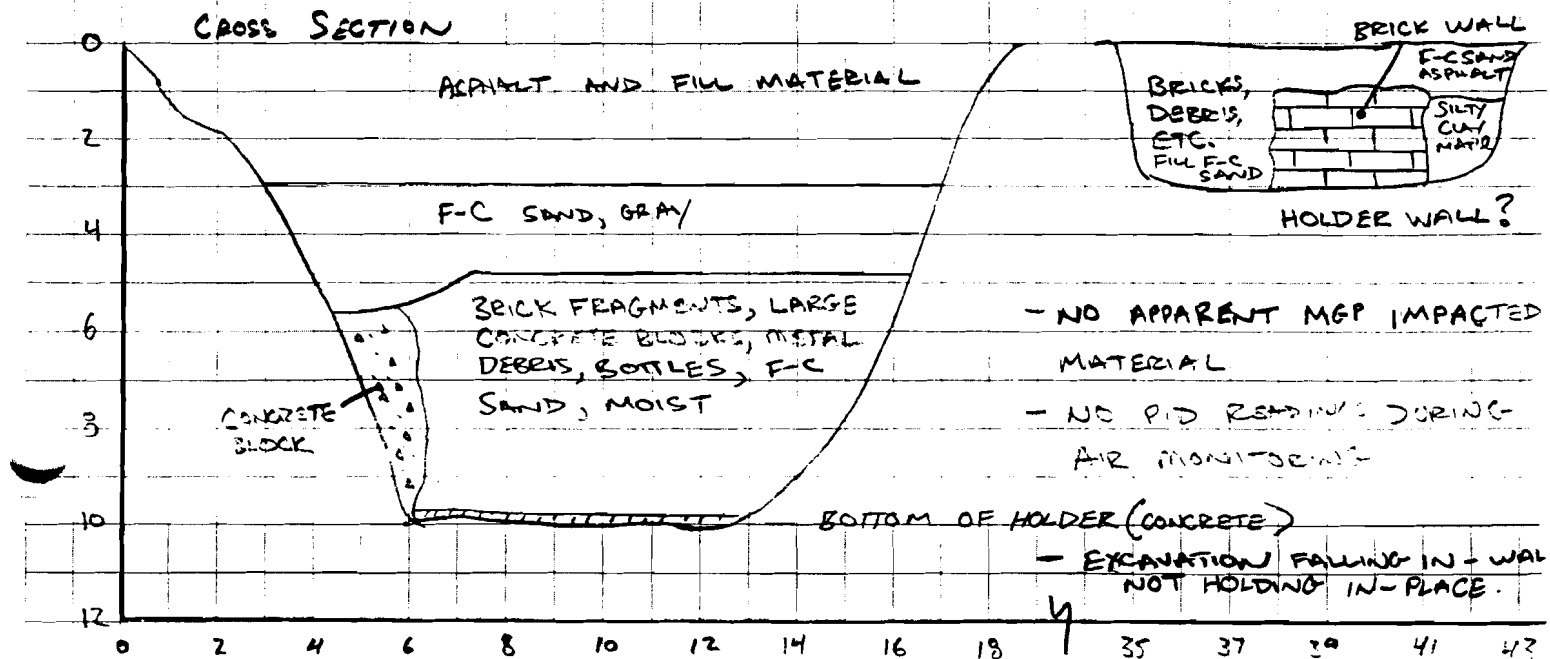
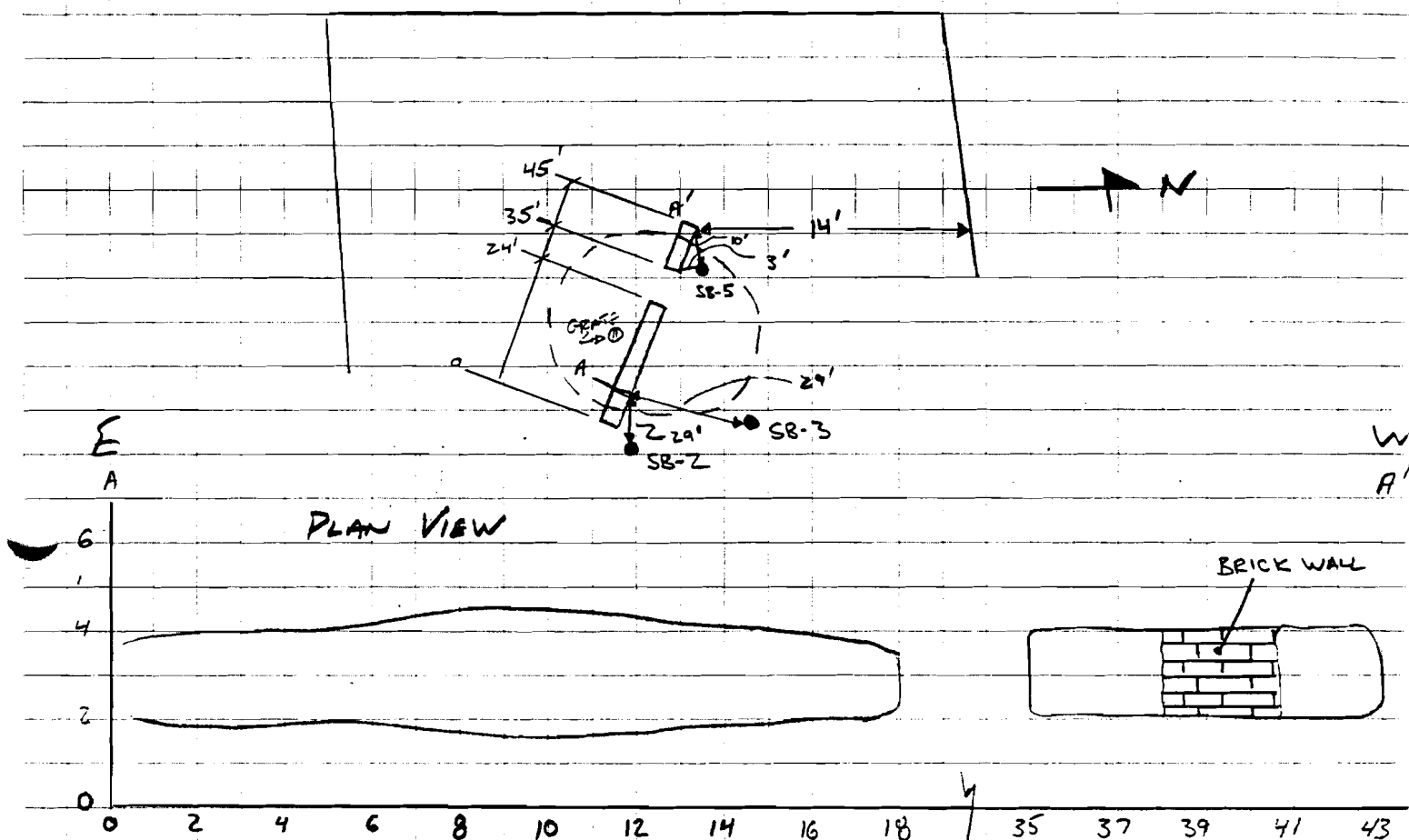
|                                    |               |           |     |        |       |
|------------------------------------|---------------|-----------|-----|--------|-------|
| SUBJECT                            | ONEIDA        | PROJ. NO. | BY  | DATE   | SHEET |
| NIAGARA MOHAWK POWER CORPORATION - | 141 CEDAR ST. | 364.56    | TMM | 8/2/00 |       |

CS. BY \_\_\_\_\_; DATE \_\_\_\_\_

CHECKED BY \_\_\_\_\_; DATE \_\_\_\_\_

TP-1

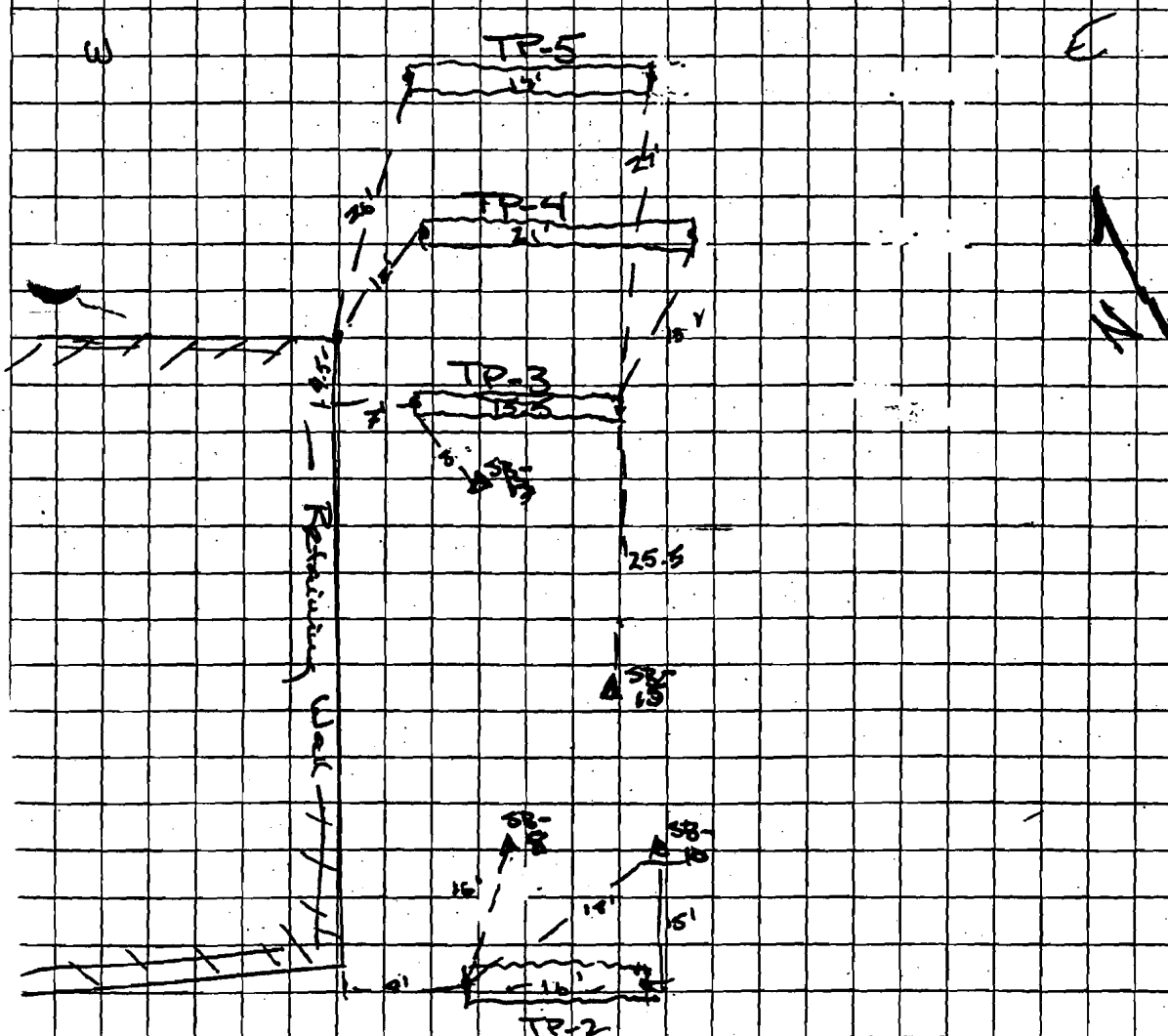
CEDAR ST.



|                           |           |     |         |       |
|---------------------------|-----------|-----|---------|-------|
| SUBJECT                   | PROJ. NO. | BY  | DATE    | SHEET |
| Onida Cedar St Test Plots | 36456     | MHC | 12/7/00 |       |

CALCS. BY \_\_\_\_\_; DATE \_\_\_\_\_ CHECKED BY \_\_\_\_\_; DATE \_\_\_\_\_

TP-2 16'  
TP-3 15.5'  
TP-4 20'  
TP-5 17'



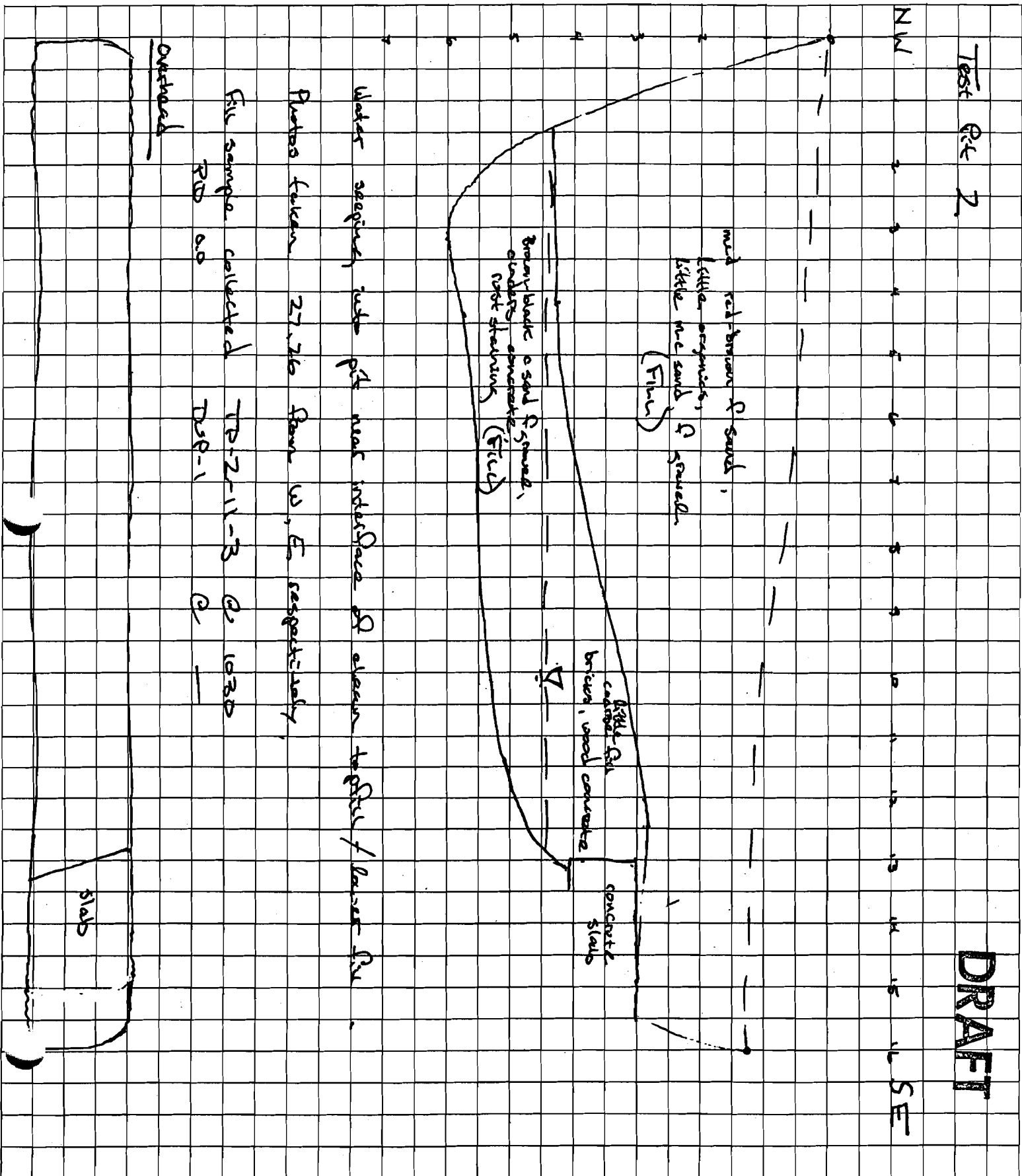
Not to scale.

DRAFT



|                                       |                           |                  |                        |       |
|---------------------------------------|---------------------------|------------------|------------------------|-------|
| SUBJECT<br><b>Oneida Cedar Street</b> | PROJ. NO.<br><b>36456</b> | BY<br><b>MKC</b> | DATE<br><b>12/2/80</b> | SHEET |
|                                       |                           |                  |                        |       |

CS. BY \_\_\_\_\_; DATE \_\_\_\_\_ CHECKED BY \_\_\_\_\_; DATE \_\_\_\_\_



**DRAFT**

**SUBJECT**

Oneida Cedar Street

PROJ. NO.

36456

BY

MVC

DATE \_\_\_\_\_

12/7/00

**SHEET**

CS. BY \_\_\_\_\_; DATE \_\_\_\_\_

CHECKED BY \_\_\_\_\_; DATE \_\_\_\_\_

|      |    |   |
|------|----|---|
| last | of | w |
|------|----|---|

NE

17

med. red-brown & sand  
loose little inc. sand,  
& gravel (fill)

mic sand, gravel  
cinders, concrete, brick,  
wire (PVC)

Test. bad. soft

Brown-belted charge still  
up toward broken floor.  
Said no flies 7

|          |        |
|----------|--------|
| Concrete | 5 beds |
|----------|--------|

slight  
broom  
or water

Water seeping into pit near Topail/Least Oil contacts, shows slow  
Slight carbonate adit upon aquifers, railroad ties.  
Topail Q sand is very loose, collapses into pit.

Sample of pest collected  
P.D. 11

TP-3-8-6 @ 120

Overhead

२५७

**DRAFT**

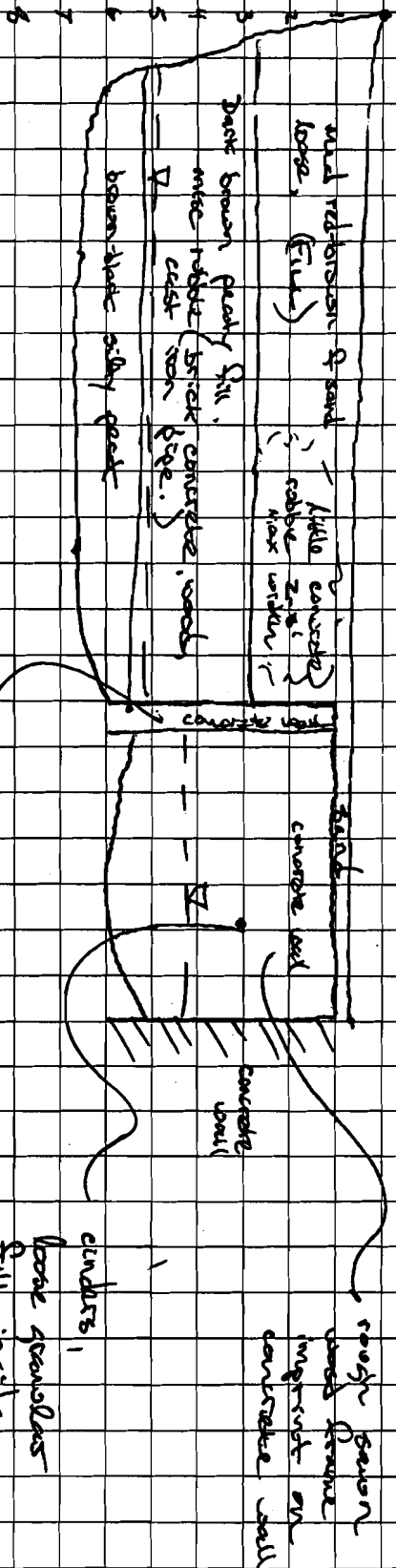
|                     |           |     |         |       |
|---------------------|-----------|-----|---------|-------|
| SUBJECT             | PROJ. NO. | BY  | DATE    | SHEET |
| Oneida Cedar Street | 36456     | MKC | 12/7/00 |       |

CS. BY \_\_\_\_\_; DATE \_\_\_\_\_ CHECKED BY \_\_\_\_\_; DATE \_\_\_\_\_

**DRAFT**

Test Pit 4

N 110° 2' 4" 6" 8" 10" 12" 14" 16" 18" 20" 22" SE



cinders,  
loose gravel  
fill inside  
chamber  
(P.D. 0-5)

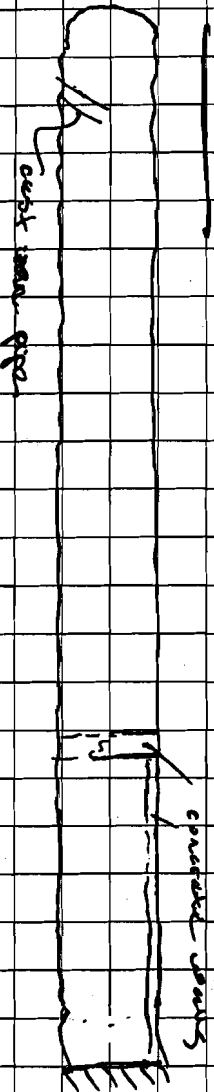
wall broken to approx 4' by.

Samples:  
TR-4-3-3 @ 1150  
(peaty fill, P.D. 0-5)  
TR-4-10-6 @ 1200 (w/ins/mass)  
(silty peat, P.D. 0-1)

Under slab to deep into western part of pit.

Water apparently perched in concrete chambers &

Overhead



cast iron pipe

concrete walls

SUBJECT

Oneida Cedar St.

PROJ. NO.

30450

BY

MLC

DATE

12/7/00

SHEET

CALCS. BY \_\_\_\_\_; DATE \_\_\_\_\_

CHECKED BY \_\_\_\_\_; DATE \_\_\_\_\_

Test Pit 5

DRAFT

N

S E

red brown fine sand (fill)

large broken bricks,  
light wood,  
sand in voids  
(fill)

gray brick c sand gravel  
small chunks, fragments  
(fill)

gray-white  
brick ash,  
light cinders, large  
Red, light  
w/iron silt and clay

large bottle cinders,  
some fine sand and  
gravel (fill)

Samples: TP-5-2-5 brown/black m.c. sand / c gravel. All cinders, brick, concrete  
P/B 0.0

TP-5-11-4 light brown/black peat, silt and clay, peaty odor.  
P/B 4.1

TP-5-5-5 gray black cinders / ash loose, no odors.  
P/B 0.0  
sample standard 1250

Water seeping into pit at approx 5.5' bgs.  
trace shown on water.

Photos 21, 20 taken from W, E, respectively.

## **Attachment 4**

# **Laboratory Analytical Results**



# Galson Laboratories

## VOLATILE ANALYTICAL REPORT

Client : Blasland, Bouck & Lee  
Account # : 10624  
Site : NMPC ONEIDA CEDAR ST

Date Received : 15-AUG-00  
Date Sampled : 15-AUG-00  
Date Extracted: NA

Matrix : Water  
Method : NYSDEC ASP 95-1  
Units : ug/l

| Galson ID: | L62754-1 | L62754-2       | WG24857-1    |
|------------|----------|----------------|--------------|
| Client ID: | PZ-1     | PZ-TRIPBLANK-2 | Method Blank |

|                            |     |     |     |
|----------------------------|-----|-----|-----|
| Chloromethane              | <10 | <10 | <10 |
| Bromomethane               | <10 | <10 | <10 |
| Vinyl Chloride             | <10 | <10 | <10 |
| Chloroethane               | <10 | <10 | <10 |
| Methylene Chloride         | <10 | <10 | <10 |
| Acetone                    | <10 | <10 | <10 |
| Carbon Disulfide           | <10 | <10 | <10 |
| 1,1-Dichloroethene         | <10 | <10 | <10 |
| 1,1-Dichloroethane         | <10 | <10 | <10 |
| 1,2-Dichloroethene (Total) | <10 | <10 | <10 |
| Chloroform                 | <10 | <10 | <10 |
| 1,2-Dichloroethane         | <10 | <10 | <10 |
| 2-Butanone                 | <10 | <10 | <10 |
| 1,1,1-Trichloroethane      | <10 | <10 | <10 |
| Carbon Tetrachloride       | <10 | <10 | <10 |
| Bromodichloromethane       | <10 | <10 | <10 |
| 1,2-Dichloropropane        | <10 | <10 | <10 |
| cis-1,3-Dichloropropene    | <10 | <10 | <10 |
| Trichloroethene            | <10 | <10 | <10 |
| Dibromochloromethane       | <10 | <10 | <10 |
| 1,1,2-Trichloroethane      | <10 | <10 | <10 |
| Benzene                    | <10 | <10 | <10 |
| trans-1,3-Dichloropropene  | <10 | <10 | <10 |
| Bromoform                  | <10 | <10 | <10 |
| 4-Methyl-2-Pentanone       | <10 | <10 | <10 |
| 2-Hexanone                 | <10 | <10 | <10 |
| Tetrachloroethene          | <10 | <10 | <10 |

Approved by : PJT  
Date : 30-AUG-00  
QC by : *[Signature]*  
Date : 9/6/00  
NYS DOH # : 11626  
Footnotes:



## WATER VOLATILE SURROGATE RECOVERY

Client : Blasland, Bouck & Lee

Login # : L62754

[illegible]

SMC1 (TOL) = Toluene-d8  
SMC2 (BFB) = Bromofluorobenzene  
SMC3 (DCE) = 1,2-Dichloroethane-d4

QC LIMITS  
(62-127)  
(73-120)  
(68-121)

```
# Column to be used to flag recovery values
* Values outside of QC limits
D Surrogate diluted out
```

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

PZ-1

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62754

Matrix: (soil/water) WATER

Lab Sample ID: L62754-1

Sample wt/vol: 1050 (g/mL) mL

Lab File ID: ED82104

Level: (low/med) LOW

Date Received: 08/15/00

% Moisture: decanted: (Y/N) N

Date Extracted: 08/18/00

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: 7.0

| CAS NO.        | COMPOUND                   | CONCENTRATION UNITS: |   |
|----------------|----------------------------|----------------------|---|
|                |                            | (ug/L or ug/Kg) ug/l | Q |
| 51-28-5-----   | 2,4-Dinitrophenol          | 24                   | U |
| 100-02-7-----  | 4-Nitrophenol              | 24                   | 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             | 24                   | U |
| 534-52-1-----  | 4,6-Dinitro-2-methylphenol | 24                   | U |
| 86-30-6-----   | N-Nitrosodiphenylamine     | 10                   | U |
| 101-55-3-----  | 4-Bromophenyl-phenylether  | 10                   | U |
| 118-74-1-----  | Hexachlorobenzene          | 10                   | U |
| 87-86-5-----   | Pentachlorophenol          | 24                   | 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 |
| 56-55-3-----   | Benzo(a)anthracene         | 10                   | U |
| 91-94-1-----   | 3,3'-Dichlorobenzidine     | 10                   | U |
| 218-01-9-----  | Chrysene                   | 10                   | U |
| 117-81-7-----  | bis(2-Ethylhexyl)phthalate | 10                   | U |
| 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-----   | Dibenzo(a,h)anthracene     | 10                   | U |
| 191-24-2-----  | Benzo(g,h,i)perylene       | 10                   | U |



1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SBLK WG24913

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62754

Matrix: (soil/water) WATER

Lab Sample ID: WG24913-1

Sample wt/vol: 1000 (g/mL) mL

Lab File ID: ED82103

Level: (low/med) LOW

Date Received: 00/00/00

% Moisture: decanted: (Y/N) N

Date Extracted: 08/18/00

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: 7.0

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     | 10 | U |
| 101-55-3-----  | 4-Bromophenyl-phenylether  | 10 | U |
| 118-74-1-----  | Hexachlorobenzene          | 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 |
| 56-55-3-----   | Benzo(a)anthracene         | 10 | U |
| 91-94-1-----   | 3,3'-Dichlorobenzidine     | 10 | U |
| 218-01-9-----  | Chrysene                   | 10 | U |
| 117-81-7-----  | bis(2-Ethylhexyl)phthalate | 10 | U |
| 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-----   | Dibenzo(a,h)anthracene     | 10 | U |
| 191-24-2-----  | Benzo(g,h,i)perylene       | 10 | U |

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

PZ-1

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62754

Matrix: (soil/water) WATER

Lab Sample ID: L62754-1

Sample wt/vol: 1000 (g/mL) mL

Lab File ID: HP13A\0817C020

% Moisture:           decanted: (Y/N) N

Date Received: 08/15/00

Extraction:   (SepF/Cont/Sonc) SEPF

Date Extracted: 08/17/00

Concentrated Extract Volume: 10000 (uL)

Date Analyzed: 08/18/00

Injection Volume: 1.0 (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

CAS NO.

COMPOUND

Q

|                 |                     |      |   |
|-----------------|---------------------|------|---|
| 319-84-6-----   | alpha-BHC           | 0.05 | U |
| 319-85-7-----   | beta-BHC            | 0.05 | U |
| 319-86-8-----   | delta-BHC           | 0.05 | U |
| 58-89-9-----    | gamma-BHC (Lindane) | 0.05 | U |
| 76-44-8-----    | Heptachlor          | 0.05 | U |
| 309-00-2-----   | Aldrin              | 0.05 | U |
| 1024-57-3-----  | Heptachlor epoxide  | 0.05 | U |
| 959-98-8-----   | Endosulfan I        | 0.05 | U |
| 60-57-1-----    | Dieldrin            | 0.1  | U |
| 72-55-9-----    | 4,4'-DDE            | 0.1  | U |
| 72-20-8-----    | Endrin              | 0.1  | U |
| 332-13-659----- | Endosulfan II       | 0.1  | U |
| 72-54-8-----    | 4,4'-DDD            | 0.1  | U |
| 103-10-78-----  | Endosulfan sulfate  | 0.1  | U |
| 50-29-3-----    | 4,4'-DDT            | 0.1  | U |
| 72-43-5-----    | Methoxychlor        | 0.5  | U |
| 53494-70-5----- | Endrin ketone       | 0.1  | U |
| 7421-36-3-----  | Endrin aldehyde     | 0.1  | U |
| 5103-71-9-----  | alpha-Chlordane     | 0.05 | U |
| 5103-7-42-----  | gamma-Chlordane     | 0.05 | 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 |

FORM I PEST

**Contract:**

Lab Code: Case No.: 1 SAS No.: SDG No.: L62754

GC Column(1): DB-608 ID: .53 (mm) GC Column(2): DB-1701 ID: .53 (mm)

[illegible]

ADVISORY  
QC LIMITS  
(30-150)  
(30-150)

```
# Column to be used to flag recovery values
* Values outside of QC limits
D Surrogate diluted out
```



## METALS ANALYTICAL REPORT

Client : Blasland, Bouck & Lee  
Account # : 10624  
Site : NMPC ONEIDA CEDAR ST

Date Received : 15-AUG-00  
Date Sampled : 15-AUG-00

Matrix : Water  
Method : SW846 6010B/7470A; ICP/CVAA

| Galson ID:           | L62754-1 | L62754-3 | WG24971-1    |
|----------------------|----------|----------|--------------|
| Client ID:           | PZ-1     | PZ-1     | Method Blank |
|                      | Units    |          |              |
| Beryllium, dissolved | mg/l NR  | <0.0050  | NR           |
| Zinc                 | mg/l 1.3 | NR       | <0.01        |
| Cadmium, dissolved   | mg/l NR  | <0.0050  | NR           |
| Calcium, dissolved   | mg/l NR  | 440      | NR           |
| Chromium, dissolved  | mg/l NR  | <0.010   | NR           |
| Cobalt, dissolved    | mg/l NR  | <0.010   | NR           |
| Copper, dissolved    | mg/l NR  | <0.010   | NR           |
| Iron, dissolved      | mg/l NR  | 0.79     | NR           |
| Lead, dissolved      | mg/l NR  | <0.0030  | NR           |
| Magnesium, dissolved | mg/l NR  | 20       | NR           |
| Manganese, dissolved | mg/l NR  | 1.8      | NR           |
| Nickel, dissolved    | mg/l NR  | <0.020   | NR           |
| Potassium, dissolved | mg/l NR  | 12       | NR           |
| Selenium, dissolved  | mg/l NR  | 0.019    | NR           |
| Silver, dissolved    | mg/l NR  | <0.010   | NR           |
| Sodium, dissolved    | mg/l NR  | 5.3      | NR           |
| Thallium, dissolved  | mg/l NR  | <0.010   | NR           |
| Vanadium, dissolved  | mg/l NR  | <0.010   | NR           |
| Zinc, dissolved      | mg/l NR  | 1.1      | NR           |

Approved by : JK  
Date : 06-SEP-00  
QC by : *[Signature]*  
Date : 9/6/00  
NYS DOH # : 11626  
Footnotes:





## INORGANIC ANALYTICAL REPORT



Client : Blasland, Bouck & Lee  
Account # : 10624  
Site : NMPC ONEIDA CEDAR ST

Date Received : 15-AUG-00  
Date Sampled : 15-AUG-00

Matrix : Water

Galson ID: L62754-1  
Client ID: PZ-1

|                | Method | Units |       |
|----------------|--------|-------|-------|
| Cyanide, Total | CLP-M  | mg/l  | <0.01 |

Approved by : DM  
Date : 24-AUG-00  
QC by :   
Date :   
NYS DOH # : 11626  
Footnotes:







Environmental Laboratories, Inc.

587 East Middle Turnpike, P.O. Box 418, Manchester, CT 06040  
Tel. (860) 645-1102 Fax (860) 645-0823

# Analysis Report

August 14, 2000

FOR: Mr. Mike Fifield  
Blasland & Bouck Engineers, PC  
6723 Towpath Road Box 66  
Syracuse NY 13214-0066

## Sample Information

Matrix: SOLID  
Location Code: BLASLAND  
Project Code:  
P.O.#: 36456

## Custody Information

Collected by: MC  
Received by: SW  
Analyzed by: see below

Date Time  
07/18/00 15:45  
07/20/00 11:00

## Laboratory Data

Client ID: NM PC ONEIDA CEDAR ST SB-1 0-4

Phoenix I.D. AC80361

| Parameter                        | Result    | RL    | Units | Date     | by  | Reference    |
|----------------------------------|-----------|-------|-------|----------|-----|--------------|
| TCLP Mercury                     | BDL       | 0.001 | mg/L  | 07/24/00 | RS  | E1311/E245.1 |
| Benzene                          | 4400      | 250   | ug/Kg | 07/24/00 | R/B | SW8021       |
| PP Silver                        | BDL       | 0.01  | mg/L  | 07/25/00 | EK  | E1311/SW6010 |
| TCLP Arsenic                     | 0.03      | 0.01  | mg/L  | 07/25/00 | EK  | E1311/SW6010 |
| TCLP Barium                      | 0.49      | 0.01  | mg/L  | 07/25/00 | EK  | E1311/SW6010 |
| TCLP Cadmium                     | BDL       | 0.005 | mg/L  | 07/25/00 | EK  | E1311/SW6010 |
| TCLP Chromium                    | BDL       | 0.01  | mg/L  | 07/25/00 | EK  | E1311/SW6010 |
| TCLP Lead                        | 0.101     | 0.015 | mg/L  | 07/25/00 | EK  | E1311/SW6010 |
| TCLP Selenium                    | BDL       | 0.05  | mg/L  | 07/25/00 | EK  | E1311/SW6010 |
| Sulfur                           | 0.20      | 0.01  | %     | 08/11/00 | OL  | ASTMD129     |
| Total Organic Halogens           | ND        | 25    | mg/kg | 07/31/00 | OL  | SW9020       |
| Sonication Ext. For PCB          | Completed |       |       | 07/20/00 | PL  | SW846-3550   |
| Soil Ext. for Semi- Vol          | Completed |       |       | 07/20/00 | B/D | SW3550/3545  |
| Percent Solid                    | 84        |       | %     | 07/21/00 | CF  | E160.3       |
| TCLP Digestion Mercury           | Completed |       |       | 07/24/00 | JA  | E1311/7470   |
| TCLP Extraction Metals           | Completed |       |       | 07/20/00 | T/A | EPA 1311     |
| Tot.Petroleum HC                 | 240       | 40    | mg/Kg | 07/26/00 | OL  | E418.1       |
| Total Cyanide                    | 2.87      | 0.23  | mg/Kg | 07/25/00 | PJ  | SW9010       |
| <b>Polychlorinated Biphenyls</b> |           |       |       |          |     |              |
| PCB-1016                         | ND        | 400   | ug/Kg | 07/21/00 | KCA | SW 8082      |
| PCB-1221                         | ND        | 400   | ug/Kg | 07/21/00 | KCA | SW 8082      |
| PCB-1232                         | ND        | 400   | ug/Kg | 07/21/00 | KCA | SW 8082      |
| 3-1242                           | ND        | 400   | ug/Kg | 07/21/00 | KCA | SW 8082      |
| PCB-1248                         | ND        | 400   | ug/Kg | 07/21/00 | KCA | SW 8082      |

| Parameter                   | Result | RL     | Units | Date     | by  | Reference |
|-----------------------------|--------|--------|-------|----------|-----|-----------|
| PCB-1254                    | ND     | 400    | ug/Kg | 07/21/00 | KCA | SW 8082   |
| PCB-1260                    | ND     | 400    | ug/Kg | 07/21/00 | KCA | SW 8082   |
| PCB-1262                    | ND     | 400    | ug/Kg | 07/21/00 | KCA | SW 8082   |
| PCB-1268                    | ND     | 400    | ug/Kg | 07/21/00 | KCA | SW 8082   |
| <b><u>Semivolatiles</u></b> |        |        |       |          |     |           |
| 1,2,4-Trichlorobenzene      | ND     | 33000  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 1,2-Dichlorobenzene         | ND     | 33000  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 1,2-Diphenylhydrazine       | ND     | 33000  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 1,3-Dichlorobenzene         | ND     | 33000  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 1,4-Dichlorobenzene         | ND     | 33000  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2,4,5-Trichlorophenol       | ND     | 33000  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2,4,6-Trichlorophenol       | ND     | 33000  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2,4-Dichlorophenol          | ND     | 33000  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2,4-Dimethylphenol          | ND     | 33000  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2,4-Dinitrophenol           | ND     | 160000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2,4-Dinitrotoluene          | ND     | 33000  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2,6-Dichlorophenol          | ND     | 33000  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2,6-Dinitrotoluene          | ND     | 33000  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2-Chloronaphthalene         | ND     | 33000  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2-Chlorophenol              | ND     | 33000  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2-Methylnaphthalene         | ND     | 33000  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2-Methylphenol (o-cresol)   | ND     | 33000  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2-Nitroaniline              | ND     | 160000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2-Nitrophenol               | ND     | 33000  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 3,3'-Dichlorobenzidine      | ND     | 66000  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 3-Nitroaniline              | ND     | 160000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 4,6-Dinitro-2-methylphenol  | ND     | 160000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 4-Bromophenyl phenyl ether  | ND     | 33000  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 4-Chloro-3-methylphenol     | ND     | 66000  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 4-Chloroaniline             | ND     | 66000  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 4-Chlorophenyl phenyl ether | ND     | 33000  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 4-Methylphenol (p-cresol)   | ND     | 33000  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 4-Nitroaniline              | ND     | 160000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 4-Nitrophenol               | ND     | 160000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Acenaphthene                | ND     | 33000  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Acenaphthylene              | 53000  | 33000  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Anthracene                  | 87000  | 33000  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Benzidine                   | ND     | 33000  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Benzo(a)anthracene          | 160000 | 33000  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Benzo(a)pyrene              | 110000 | 33000  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Benzo(b)fluoranthene        | 160000 | 33000  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Benzo(g,h,i)perylene        | 34000  | 33000  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Benzo(k)fluoranthene        | 76000  | 33000  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Benzoic acid                | ND     | 160000 | ug/Kg | 07/22/00 | DRC | SW 8270   |




| Parameter                           | Result      | RL    | Units | Date     | by  | Reference |
|-------------------------------------|-------------|-------|-------|----------|-----|-----------|
| Benzyl alcohol                      | ND          | 66000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Butyl butyl phthalate               | ND          | 33000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Bis(2-chloroethoxy)methane          | ND          | 33000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Bis(2-chloroethyl)ether             | ND          | 33000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Bis(2-chloroisopropyl)ether         | ND          | 33000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Bis(2-ethylhexyl)phthalate          | ND          | 33000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Chrysene                            | 130000      | 33000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Di-n-butylphthalate                 | ND          | 33000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Di-n-octyl phthalate                | ND          | 33000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Dibenz(a,h)anthracene               | ND          | 33000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Dibenzofuran                        | 37000       | 33000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Diethyl phthalate                   | ND          | 33000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Dimethyl phthalate                  | ND          | 33000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Fluoranthene                        | 330000      | 33000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Fluorene                            | 47000       | 33000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Hexachlorobenzene                   | ND          | 33000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Hexachlorobutadiene                 | ND          | 33000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Hexachlorocyclopentadiene           | ND          | 33000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Hexachloroethane                    | ND          | 33000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Indeno(1,2,3-c,d)pyrene             | 38000       | 33000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Isophorone                          | ND          | 33000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| N-Nitrosodi-n-propylamine           | ND          | 33000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| N-Nitrosodimethylamine              | ND          | 33000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| N-Nitrosodiphenylamine              | ND          | 33000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Naphthalene                         | 180000      | 33000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Nitrobenzene                        | ND          | 33000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Pentachlorophenol                   | ND          | 33000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Phenanthrene                        | 290000      | 33000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Phenol                              | ND          | 33000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Pyrene                              | 260000      | 33000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| % 2,4,6-Tribromophenol (Surrog Rec) | Diluted Out |       | %     | 07/22/00 | DRC | SW 8270   |
| % 2-Fluorobiphenyl (Surrogate Rec)  | Diluted Out |       | %     | 07/22/00 | DRC | SW 8270   |
| % 2-Fluorophenol (Surrogate Rec)    | Diluted Out |       | %     | 07/22/00 | DRC | SW 8270   |
| % Nitrobenzene-d5 (Surrogate Rec)   | Diluted Out |       | %     | 07/22/00 | DRC | SW 8270   |
| % Phenol-d5 (Surrogate Rec)         | Diluted Out |       | %     | 07/22/00 | DRC | SW 8270   |
| % Terphenyl-d14 (Surrogate Rec)     | Diluted Out |       | %     | 07/22/00 | DRC | SW 8270   |

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**Comments:**

ND=Not detected BDL = Below Detection Limit RL=Reporting Limit

If there are any questions regarding this data, please call Phoenix Client Services at extension 200.

  
Phyllis Shiller, Laboratory Director  
August 14, 2000



Environmental Laboratories, Inc.

587 East Middle Turnpike, P.O. Box 418 Manchester, CT 06040  
Tel. (860) 645-1102 Fax (860) 645-0823

# Analysis Report

August 14, 2000

FOR: Mr. Mike Fifield  
Blasland & Bouck Engineers, PC  
6723 Towpath Road Box 66  
Syracuse NY 13214-0066

## Sample Information

Matrix: SOLID  
Location Code: BLASLAND  
Project Code:  
P.O.#: 36456

## Custody Information

Collected by: MC  
Received by: SW  
Analyzed by: see below

Date Time  
07/19/00 10:40  
07/20/00 11:00

## Laboratory Data

Client ID: NM PC ONEIDA CEDAR ST SB-2 0-6

Phoenix I.D. AC80362

| Parameter                        | Result    | RL    | Units | Date     | by  | Reference    |
|----------------------------------|-----------|-------|-------|----------|-----|--------------|
| TCLP Mercury                     | BDL       | 0.001 | mg/L  | 07/24/00 | RS  | E1311/E245.1 |
| Benzene                          | ND        | 5.0   | ug/Kg | 07/24/00 | R/B | SW8021       |
| LP Silver                        | BDL       | 0.01  | mg/L  | 07/25/00 | EK  | E1311/SW6010 |
| TCLP Arsenic                     | BDL       | 0.01  | mg/L  | 07/25/00 | EK  | E1311/SW6010 |
| TCLP Barium                      | 0.77      | 0.01  | mg/L  | 07/25/00 | EK  | E1311/SW6010 |
| TCLP Cadmium                     | BDL       | 0.005 | mg/L  | 07/25/00 | EK  | E1311/SW6010 |
| TCLP Chromium                    | BDL       | 0.01  | mg/L  | 07/25/00 | EK  | E1311/SW6010 |
| TCLP Lead                        | 0.101     | 0.015 | mg/L  | 07/25/00 | EK  | E1311/SW6010 |
| TCLP Selenium                    | BDL       | 0.05  | mg/L  | 07/25/00 | EK  | E1311/SW6010 |
| Sulfur                           | 0.065     | 0.01  | %     | 08/11/00 | OL  | ASTMD129     |
| Total Organic Halogens           | ND        | 25    | mg/kg | 07/31/00 | OL  | SW9020       |
| Sonication Ext. For PCB          | Completed |       |       | 07/20/00 | PL  | SW846-3550   |
| Soil Ext. for Semi- Vol          | Completed |       |       | 07/20/00 | B/D | SW3550/3545  |
| Percent Solid                    | 89        |       | %     | 07/21/00 | CF  | E160.3       |
| TCLP Digestion Mercury           | Completed |       |       | 07/24/00 | JA  | E1311/7470   |
| TCLP Extraction Metals           | Completed |       |       | 07/20/00 | T/A | EPA 1311     |
| Tot.Petroleum HC                 | BDL       | 40    | mg/Kg | 07/26/00 | OL  | E418.1       |
| Total Cyanide                    | BDL       | 0.22  | mg/Kg | 07/25/00 | PJ  | SW9010       |
| <b>Polychlorinated Biphenyls</b> |           |       |       |          |     |              |
| PCB-1016                         | ND        | 400   | ug/Kg | 07/21/00 | KCA | SW 8082      |
| PCB-1221                         | ND        | 400   | ug/Kg | 07/21/00 | KCA | SW 8082      |
| PCB-1232                         | ND        | 400   | ug/Kg | 07/21/00 | KCA | SW 8082      |
| B-1242                           | ND        | 400   | ug/Kg | 07/21/00 | KCA | SW 8082      |
| PCB-1248                         | ND        | 400   | ug/Kg | 07/21/00 | KCA | SW 8082      |

| Parameter                   | Result | RL   | Units | Date by  | Reference   |
|-----------------------------|--------|------|-------|----------|-------------|
| PCB-1254                    | ND     | 400  | ug/Kg | 07/21/00 | KCA SW 8082 |
| PCB-1260                    | ND     | 400  | ug/Kg | 07/21/00 | KCA SW 8082 |
| PCB-1262                    | ND     | 400  | ug/Kg | 07/21/00 | KCA SW 8082 |
| PCB-1268                    | ND     | 400  | ug/Kg | 07/21/00 | KCA SW 8082 |
| <b><u>Semivolatiles</u></b> |        |      |       |          |             |
| 1,2,4-Trichlorobenzene      | ND     | 330  | ug/Kg | 07/21/00 | DRC SW 8270 |
| 1,2-Dichlorobenzene         | ND     | 330  | ug/Kg | 07/21/00 | DRC SW 8270 |
| 1,2-Diphenylhydrazine       | ND     | 330  | ug/Kg | 07/21/00 | DRC SW 8270 |
| 1,3-Dichlorobenzene         | ND     | 330  | ug/Kg | 07/21/00 | DRC SW 8270 |
| 1,4-Dichlorobenzene         | ND     | 330  | ug/Kg | 07/21/00 | DRC SW 8270 |
| 2,4,5-Trichlorophenol       | ND     | 330  | ug/Kg | 07/21/00 | DRC SW 8270 |
| 2,4,6-Trichlorophenol       | ND     | 330  | ug/Kg | 07/21/00 | DRC SW 8270 |
| 2,4-Dichlorophenol          | ND     | 330  | ug/Kg | 07/21/00 | DRC SW 8270 |
| 2,4-Dimethylphenol          | ND     | 330  | ug/Kg | 07/21/00 | DRC SW 8270 |
| 2,4-Dinitrophenol           | ND     | 1600 | ug/Kg | 07/21/00 | DRC SW 8270 |
| 2,4-Dinitrotoluene          | ND     | 330  | ug/Kg | 07/21/00 | DRC SW 8270 |
| 2,6-Dichlorophenol          | ND     | 330  | ug/Kg | 07/21/00 | DRC SW 8270 |
| 2,6-Dinitrotoluene          | ND     | 330  | ug/Kg | 07/21/00 | DRC SW 8270 |
| 2-Chloronaphthalene         | ND     | 330  | ug/Kg | 07/21/00 | DRC SW 8270 |
| 2-Chlorophenol              | ND     | 330  | ug/Kg | 07/21/00 | DRC SW 8270 |
| 2-Methylnaphthalene         | ND     | 330  | ug/Kg | 07/21/00 | DRC SW 8270 |
| 2-Methylphenol (o-cresol)   | ND     | 330  | ug/Kg | 07/21/00 | DRC SW 8270 |
| 2-Nitroaniline              | ND     | 1600 | ug/Kg | 07/21/00 | DRC SW 8270 |
| 2-Nitrophenol               | ND     | 330  | ug/Kg | 07/21/00 | DRC SW 8270 |
| 3,3'-Dichlorobenzidine      | ND     | 660  | ug/Kg | 07/21/00 | DRC SW 8270 |
| 3-Nitroaniline              | ND     | 1600 | ug/Kg | 07/21/00 | DRC SW 8270 |
| 4,6-Dinitro-2-methylphenol  | ND     | 1600 | ug/Kg | 07/21/00 | DRC SW 8270 |
| 4-Bromophenyl phenyl ether  | ND     | 330  | ug/Kg | 07/21/00 | DRC SW 8270 |
| 4-Chloro-3-methylphenol     | ND     | 660  | ug/Kg | 07/21/00 | DRC SW 8270 |
| 4-Chloroaniline             | ND     | 660  | ug/Kg | 07/21/00 | DRC SW 8270 |
| 4-Chlorophenyl phenyl ether | ND     | 330  | ug/Kg | 07/21/00 | DRC SW 8270 |
| 4-Methylphenol (p-cresol)   | ND     | 330  | ug/Kg | 07/21/00 | DRC SW 8270 |
| 4-Nitroaniline              | ND     | 1600 | ug/Kg | 07/21/00 | DRC SW 8270 |
| 4-Nitrophenol               | ND     | 1600 | ug/Kg | 07/21/00 | DRC SW 8270 |
| Acenaphthylene              | ND     | 330  | ug/Kg | 07/21/00 | DRC SW 8270 |
| Acenaphthene                | ND     | 330  | ug/Kg | 07/21/00 | DRC SW 8270 |
| Anthracene                  | ND     | 330  | ug/Kg | 07/21/00 | DRC SW 8270 |
| Benidine                    | ND     | 330  | ug/Kg | 07/21/00 | DRC SW 8270 |
| Benzo(a)anthracene          | ND     | 330  | ug/Kg | 07/21/00 | DRC SW 8270 |
| Benzo(a)pyrene              | ND     | 330  | ug/Kg | 07/21/00 | DRC SW 8270 |
| Benzo(b)fluoranthene        | ND     | 330  | ug/Kg | 07/21/00 | DRC SW 8270 |
| Benzo(g,h,i)perylene        | ND     | 330  | ug/Kg | 07/21/00 | DRC SW 8270 |
| Benzo(k)fluoranthene        | ND     | 330  | ug/Kg | 07/21/00 | DRC SW 8270 |
| Benzoic acid                | ND     | 1600 | ug/Kg | 07/21/00 | DRC SW 8270 |

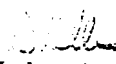
| Parameter                           | Result | RL  | Units | Date     | by  | Reference |
|-------------------------------------|--------|-----|-------|----------|-----|-----------|
| Benzyl alcohol                      | ND     | 660 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| yl butyl phthalate                  | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Bis(2-chloroethoxy)methane          | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Bis(2-chloroethyl)ether             | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Bis(2-chloroisopropyl)ether         | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Bis(2-ethylhexyl)phthalate          | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Chrysene                            | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Di-n-butylphthalate                 | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Di-n-octyl phthalate                | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Dibenz(a,h)anthracene               | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Dibenzofuran                        | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Diethyl phthalate                   | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Dimethyl phthalate                  | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Fluoranthene                        | 360    | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Fluorene                            | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Hexachlorobenzene                   | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Hexachlorobutadiene                 | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Hexachlorocyclopentadiene           | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Hexachloroethane                    | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Indeno(1,2,3-c,d)pyrene             | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Isophorone                          | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| N-Nitrosodi-n-propylamine           | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| N-Nitrosodimethylamine              | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| N-Nitrosodiphenylamine              | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Naphthalene                         | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Nitrobenzene                        | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Pentachlorophenol                   | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Phenanthrene                        | 400    | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Phenol                              | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Pyrene                              | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| % 2,4,6-Tribromophenol (Surrog Rec) | 49     |     | %     | 07/21/00 | DRC | SW 8270   |
| % 2-Fluorobiphenyl (Surrogate Rec)  | 58     |     | %     | 07/21/00 | DRC | SW 8270   |
| % 2-Fluorophenol (Surrogate Rec)    | 77     |     | %     | 07/21/00 | DRC | SW 8270   |
| % Nitrobenzene-d5 (Surrogate Rec)   | 71     |     | %     | 07/21/00 | DRC | SW 8270   |
| % Phenol-d5 (Surrogate Rec)         | 73     |     | %     | 07/21/00 | DRC | SW 8270   |
| % Terphenyl-d14 (Surrogate Rec)     | 86     |     | %     | 07/21/00 | DRC | SW 8270   |

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**Comments:**

ND=Not detected BDL = Below Detection Limit RL=Reporting Limit

If there are any questions regarding this data, please call Phoenix Client Services at extension 200.

  
Phyllis Shiller, Laboratory Director  
August 14, 2000



Environmental Laboratories, Inc.

567 East Middle Turnpike, P.O. Box 418, Manchester, CT 06040  
Tel. (860) 645-1102 Fax (860) 645-0823

# Analysis Report

August 14, 2000

FOR: Mr. Mike Fifield  
Blasland & Bouck Engineers, PC  
6723 Towpath Road Box 66  
Syracuse NY 13214-0066

## Sample Information

Matrix: SOLID  
Location Code: BLASLAND  
Project Code:  
P.O.#: 36456

## Custody Information

Collected by: MC  
Received by: SW  
Analyzed by: see below

Date Time  
07/19/00 10:50  
07/20/00 11:00

## Laboratory Data

Client ID: NM PC ONEIDA CEDAR SB-2 6-10

Phoenix I.D. AC80363

| Parameter               | Result    | RL    | Units | Date     | by  | Reference    |
|-------------------------|-----------|-------|-------|----------|-----|--------------|
| TCLP Mercury            | BDL       | 0.001 | mg/L  | 07/24/00 | RS  | E1311/E245.1 |
| Benzene                 | ND        | 500   | ug/Kg | 07/24/00 | R/B | SW8021       |
| TCLP Silver             | BDL       | 0.01  | mg/L  | 07/25/00 | EK  | E1311/SW6010 |
| TCLP Arsenic            | BDL       | 0.01  | mg/L  | 07/25/00 | EK  | E1311/SW6010 |
| TCLP Barium             | 1.61      | 0.01  | mg/L  | 07/25/00 | EK  | E1311/SW6010 |
| TCLP Cadmium            | 0.019     | 0.005 | mg/L  | 07/25/00 | EK  | E1311/SW6010 |
| TCLP Chromium           | BDL       | 0.01  | mg/L  | 07/25/00 | EK  | E1311/SW6010 |
| TCLP Lead               | 0.406     | 0.015 | mg/L  | 07/25/00 | EK  | E1311/SW6010 |
| TCLP Selenium           | BDL       | 0.05  | mg/L  | 07/25/00 | EK  | E1311/SW6010 |
| Sulfur                  | 0.056     | 0.01  | %     | 08/11/00 | OL  | ASTMD129     |
| Total Organic Halogens  | ND        | 25    | mg/kg | 07/31/00 | OL  | SW9020       |
| Sonication Ext. For PCB | Completed |       |       | 07/20/00 | PL  | SW846-3550   |
| Soil Ext. for Semi- Vol | Completed |       |       | 07/20/00 | B/D | SW3550/3545  |
| Percent Solid           | 83        |       | %     | 07/21/00 | CF  | E160.3       |
| TCLP Digestion Mercury  | Completed |       |       | 07/24/00 | JA  | E1311/7470   |
| TCLP Extraction Metals  | Completed |       |       | 07/20/00 | T/A | EPA 1311     |
| Tot. Petroleum HC       | 160       | 40    | mg/Kg | 07/26/00 | OL  | E418.1       |
| Total Cyanide           | 51.8      | 2.40  | mg/Kg | 07/25/00 | PJ  | SW9010       |

## Polychlorinated Biphenyls

|          |    |     |       |          |     |         |
|----------|----|-----|-------|----------|-----|---------|
| PCB-1016 | ND | 400 | ug/Kg | 07/21/00 | KCA | SW 8082 |
| PCB-1221 | ND | 400 | ug/Kg | 07/21/00 | KCA | SW 8082 |
| PCB-1232 | ND | 400 | ug/Kg | 07/21/00 | KCA | SW 8082 |
| PCB-1242 | ND | 400 | ug/Kg | 07/21/00 | KCA | SW 8082 |
| PCB-1248 | ND | 400 | ug/Kg | 07/21/00 | KCA | SW 8082 |

| Parameter                   | Result | RL    | Units | Date     | by  | Reference |
|-----------------------------|--------|-------|-------|----------|-----|-----------|
| PCB-1254                    | ND     | 400   | ug/Kg | 07/21/00 | KCA | SW 8082   |
| PCB-1260                    | ND     | 400   | ug/Kg | 07/21/00 | KCA | SW 8082   |
| PCB-1262                    | ND     | 400   | ug/Kg | 07/21/00 | KCA | SW 8082   |
| PCB-1268                    | ND     | 400   | ug/Kg | 07/21/00 | KCA | SW 8082   |
| <b><u>Semivolatiles</u></b> |        |       |       |          |     |           |
| 1,2,4-Trichlorobenzene      | ND     | 3300  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 1,2-Dichlorobenzene         | ND     | 3300  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 1,2-Diphenylhydrazine       | ND     | 3300  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 1,3-Dichlorobenzene         | ND     | 3300  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 1,4-Dichlorobenzene         | ND     | 3300  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2,4,5-Trichlorophenol       | ND     | 3300  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2,4,6-Trichlorophenol       | ND     | 3300  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2,4-Dichlorophenol          | ND     | 3300  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2,4-Dimethylphenol          | ND     | 3300  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2,4-Dinitrophenol           | ND     | 16000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2,4-Dinitrotoluene          | ND     | 3300  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2,6-Dichlorophenol          | ND     | 3300  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2,6-Dinitrotoluene          | ND     | 3300  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2-Chloronaphthalene         | ND     | 3300  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2-Chlorophenol              | ND     | 3300  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2-Methylnaphthalene         | 3400   | 3300  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2-Methylphenol (o-cresol)   | ND     | 3300  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2-Nitroaniline              | ND     | 16000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2-Nitrophenol               | ND     | 3300  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 3,3'-Dichlorobenzidine      | ND     | 6600  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 3-Nitroaniline              | ND     | 16000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 4,6-Dinitro-2-methylphenol  | ND     | 16000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 4-Bromophenyl phenyl ether  | ND     | 3300  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 4-Chloro-3-methylphenol     | ND     | 6600  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 4-Chloroaniline             | ND     | 6600  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 4-Chlorophenyl phenyl ether | ND     | 3300  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 4-Methylphenol (p-cresol)   | ND     | 3300  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 4-Nitroaniline              | ND     | 16000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 4-Nitrophenol               | ND     | 16000 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Acenaphthene                | ND     | 3300  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Acenaphthylene              | ND     | 3300  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Anthracene                  | 8500   | 3300  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Benzidine                   | ND     | 3300  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Benzo(a)anthracene          | 12000  | 3300  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Benzo(a)pyrene              | 7600   | 3300  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Benzo(b)fluoranthene        | 9000   | 3300  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Benzo(g,h,i)perylene        | ND     | 3300  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Benzo(k)fluoranthene        | 8400   | 3300  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Benzoic acid                | ND     | 16000 | ug/Kg | 07/22/00 | DRC | SW 8270   |



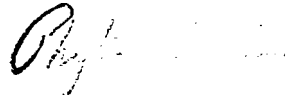
| Parameter                           | Result | RL   | Units | Date     | by  | Reference |
|-------------------------------------|--------|------|-------|----------|-----|-----------|
| Benzyl alcohol                      | ND     | 6600 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Benzyl butyl phthalate              | ND     | 3300 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Bis(2-chloroethoxy)methane          | ND     | 3300 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Bis(2-chloroethyl)ether             | ND     | 3300 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Bis(2-chloroisopropyl)ether         | ND     | 3300 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Bis(2-ethylhexyl)phthalate          | ND     | 3300 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Chrysene                            | 9900   | 3300 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Di-n-butylphthalate                 | ND     | 3300 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Di-n-octyl phthalate                | ND     | 3300 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Dibenz(a,h)anthracene               | ND     | 3300 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Dibenzofuran                        | 4500   | 3300 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Diethyl phthalate                   | ND     | 3300 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Dimethyl phthalate                  | ND     | 3300 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Fluoranthene                        | 23000  | 3300 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Fluorene                            | ND     | 3300 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Hexachlorobenzene                   | ND     | 3300 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Hexachlorobutadiene                 | ND     | 3300 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Hexachlorocyclopentadiene           | ND     | 3300 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Hexachloroethane                    | ND     | 3300 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Indeno(1,2,3-c,d)pyrene             | ND     | 3300 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Isophorone                          | ND     | 3300 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| N-Nitrosodi-n-propylamine           | ND     | 3300 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Nitrosodimethylamine                | ND     | 3300 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| N-Nitrosodiphenylamine              | ND     | 3300 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Naphthalene                         | 6900   | 3300 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Nitrobenzene                        | ND     | 3300 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Pentachlorophenol                   | ND     | 3300 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Phenanthrene                        | 25000  | 3300 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Phenol                              | ND     | 3300 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Pyrene                              | 17000  | 3300 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| % 2,4,6-Tribromophenol (Surrog Rec) | 45     |      | %     | 07/22/00 | DRC | SW 8270   |
| % 2-Fluorobiphenyl (Surrogate Rec)  | 44     |      | %     | 07/22/00 | DRC | SW 8270   |
| % 2-Fluorophenol (Surrogate Rec)    | 47     |      | %     | 07/22/00 | DRC | SW 8270   |
| % Nitrobenzene-d5 (Surrogate Rec)   | 45     |      | %     | 07/22/00 | DRC | SW 8270   |
| % Phenol-d5 (Surrogate Rec)         | 46     |      | %     | 07/22/00 | DRC | SW 8270   |
| % Terphenyl-d14 (Surrogate Rec)     | 62     |      | %     | 07/22/00 | DRC | SW 8270   |

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**Comments:**

ND=Not detected BDL = Below Detection Limit RL=Reporting Limit

If there are any questions regarding this data, please call Phoenix Client Services at extension 200.



Phyllis Shiller, Laboratory Director  
August 14, 2000



Environmental Laboratories, Inc.  
587 East Middle Turnpike, P.O. Box 418 Manchester, CT 06040  
Tel. (860) 645-1102 Fax (860) 645-0823

## Analysis Report

August 14, 2000

FOR: Mr. Mike Fifield  
Blasland & Bouck Engineers, PC  
6723 Towpath Road Box 66  
Syracuse NY 13214-0066

### Sample Information

Matrix: SOLID  
Location Code: BLASLAND  
Project Code:  
P.O.#: 36456

### Custody Information

Collected by: MC  
Received by: SW  
Analyzed by: see below

Date Time  
07/19/00 14:30  
07/20/00 11:00

## Laboratory Data

Client ID: NM PC ONEIDA CEDAR ST SB-3 0-6

Phoenix I.D. AC80364

| Parameter                               | Result    | RL    | Units | Date     | by  | Reference    |
|---|-----------|-------|-------|----------|-----|--------------|
| TCLP Mercury                            | BDL       | 0.001 | mg/L  | 07/24/00 | RS  | E1311/E245.1 |
| Benzene                                 | BDL       | 5.0   | ug/Kg | 07/25/00 | RM  | SW8021       |
| LP Silver                               | BDL       | 0.01  | mg/L  | 07/25/00 | EK  | E1311/SW6010 |
| TCLP Arsenic                            | 0.02      | 0.01  | mg/L  | 07/25/00 | EK  | E1311/SW6010 |
| TCLP Barium                             | 0.47      | 0.01  | mg/L  | 07/25/00 | EK  | E1311/SW6010 |
| TCLP Cadmium                            | BDL       | 0.005 | mg/L  | 07/25/00 | EK  | E1311/SW6010 |
| TCLP Chromium                           | BDL       | 0.01  | mg/L  | 07/25/00 | EK  | E1311/SW6010 |
| TCLP Lead                               | 0.129     | 0.015 | mg/L  | 07/25/00 | EK  | E1311/SW6010 |
| TCLP Selenium                           | BDL       | 0.05  | mg/L  | 07/25/00 | EK  | E1311/SW6010 |
| Sulfur                                  | 0.087     | 0.01  | %     | 08/11/00 | OL  | ASTMD129     |
| Total Organic Halogens                  | ND        | 25    | mg/kg | 07/31/00 | OL  | SW9020       |
| Sonication Ext. For PCB                 | Completed |       |       | 07/20/00 | PL  | SW846-3550   |
| Soil Ext. for Semi- Vol                 | Completed |       |       | 07/20/00 | B/D | SW3550/3545  |
| Percent Solid                           | 91        |       | %     | 07/21/00 | CF  | E160.3       |
| TCLP Digestion Mercury                  | Completed |       |       | 07/24/00 | JA  | E1311/7470   |
| TCLP Extraction Metals                  | Completed |       |       | 07/20/00 | T/A | EPA 1311     |
| Tot. Petroleum HC                       | 720       | 40    | mg/Kg | 07/26/00 | OL  | E418.1       |
| Total Cyanide                           | 0.296     | 0.22  | mg/Kg | 07/25/00 | PJ  | SW9010       |
| <b><u>Polychlorinated Biphenyls</u></b> |           |       |       |          |     |              |
| PCB-1016                                | ND        | 400   | ug/Kg | 07/21/00 | KCA | SW 8082      |
| PCB-1221                                | ND        | 400   | ug/Kg | 07/21/00 | KCA | SW 8082      |
| PCB-1232                                | ND        | 400   | ug/Kg | 07/21/00 | KCA | SW 8082      |
| PCB-1242                                | ND        | 400   | ug/Kg | 07/21/00 | KCA | SW 8082      |
| PCB-1248                                | ND        | 400   | ug/Kg | 07/21/00 | KCA | SW 8082      |

| Parameter                   | Result | RL   | Units | Date     | by  | Reference |
|-----------------------------|--------|------|-------|----------|-----|-----------|
| PCB-1254                    | ND     | 400  | ug/Kg | 07/21/00 | KCA | SW 8082   |
| PCB-1260                    | ND     | 400  | ug/Kg | 07/21/00 | KCA | SW 8082   |
| PCB-1262                    | ND     | 400  | ug/Kg | 07/21/00 | KCA | SW 8082   |
| PCB-1268                    | ND     | 400  | ug/Kg | 07/21/00 | KCA | SW 8082   |
| <b>Semivolatiles</b>        |        |      |       |          |     |           |
| 1,2,4-Trichlorobenzene      | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 1,2-Dichlorobenzene         | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 1,2-Diphenylhydrazine       | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 1,3-Dichlorobenzene         | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 1,4-Dichlorobenzene         | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 2,4,5-Trichlorophenol       | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 2,4,6-Trichlorophenol       | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 2,4-Dichlorophenol          | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 2,4-Dimethylphenol          | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 2,4-Dinitrophenol           | ND     | 1600 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 2,4-Dinitrotoluene          | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 2,6-Dichlorophenol          | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 2,6-Dinitrotoluene          | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 2-Chloronaphthalene         | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 2-Chlorophenol              | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 2-Methylnaphthalene         | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 2-Methylphenol (o-cresol)   | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 2-Nitroaniline              | ND     | 1600 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 2-Nitrophenol               | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 3,3'-Dichlorobenzidine      | ND     | 660  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 3-Nitroaniline              | ND     | 1600 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 4,6-Dinitro-2-methylphenol  | ND     | 1600 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 4-Bromophenyl phenyl ether  | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 4-Chloro-3-methylphenol     | ND     | 660  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 4-Chloroaniline             | ND     | 660  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 4-Chlorophenyl phenyl ether | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 4-Methylphenol (p-cresol)   | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 4-Nitroaniline              | ND     | 1600 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 4-Nitrophenol               | ND     | 1600 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Acenaphthene                | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Acenaphthylene              | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Anthracene                  | 430    | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Benzidine                   | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Benzo(a)anthracene          | 1500   | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Benzo(a)pyrene              | 2000   | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Benzo(b)fluoranthene        | 2400   | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Benzo(g,h,i)perylene        | 380    | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Benzo(k)fluoranthene        | 2400   | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Benzoic acid                | ND     | 1600 | ug/Kg | 07/21/00 | DRC | SW 8270   |


| Parameter                           | Result | RL  | Units | Date     | by  | Reference |
|-------------------------------------|--------|-----|-------|----------|-----|-----------|
| Benzyl alcohol                      | ND     | 660 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Benzyl butyl phthalate              | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Bis(2-chloroethoxy)methane          | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Bis(2-chloroethyl)ether             | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Bis(2-chloroisopropyl)ether         | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Bis(2-ethylhexyl)phthalate          | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Chrysene                            | 1200   | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Di-n-butylphthalate                 | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Di-n-octyl phthalate                | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Dibenz(a,h)anthracene               | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Dibenzofuran                        | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Diethyl phthalate                   | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Dimethyl phthalate                  | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Fluoranthene                        | 2200   | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Fluorene                            | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Hexachlorobenzene                   | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Hexachlorobutadiene                 | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Hexachlorocyclopentadiene           | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Hexachloroethane                    | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Indeno(1,2,3-c,d)pyrene             | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Isophorone                          | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| N-Nitrosodi-n-propylamine           | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| N-Nitrosodimethylamine              | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| N-Nitrosodiphenylamine              | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Naphthalene                         | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Nitrobenzene                        | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Pentachlorophenol                   | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Phenanthrene                        | 1600   | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Phenol                              | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Pyrene                              | 1800   | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| % 2,4,6-Tribromophenol (Surrog Rec) | 64     |     | %     | 07/21/00 | DRC | SW 8270   |
| % 2-Fluorobiphenyl (Surrogate Rec)  | 59     |     | %     | 07/21/00 | DRC | SW 8270   |
| % 2-Fluorophenol (Surrogate Rec)    | 84     |     | %     | 07/21/00 | DRC | SW 8270   |
| % Nitrobenzene-d5 (Surrogate Rec)   | 78     |     | %     | 07/21/00 | DRC | SW 8270   |
| % Phenol-d5 (Surrogate Rec)         | 81     |     | %     | 07/21/00 | DRC | SW 8270   |
| % Terphenyl-d14 (Surrogate Rec)     | 99     |     | %     | 07/21/00 | DRC | SW 8270   |

---

**Comments:**

ND=Not detected BDL = Below Detection Limit RL=Reporting Limit

If there are any questions regarding this data, please call Phoenix Client Services at extension 200.



Phyllis Shiller, Laboratory Director  
August 14, 2000



Environmental Laboratories, Inc.

587 East Middle Turnpike, P.O. Box 418, Manchester, CT 06040  
Tel. (860) 645-1102 Fax (860) 645-0823

## Analysis Report

August 14, 2000

FOR: Mr. Mike Fifield  
Blasland & Bouck Engineers, PC  
6723 Towpath Road Box 66  
Syracuse NY 13214-0066

### Sample Information

Matrix: SOLID  
Location Code: BLASLAND  
Project Code:  
P.O.#: 36456

### Custody Information

Collected by: MC  
Received by: SW  
Analyzed by: see below

Date Time  
07/19/00 15:00  
07/20/00 11:00

## Laboratory Data

Client ID: NM PC ONEIDA CEDAR SB-3 6-12

Phoenix I.D. AC80365

| Parameter               | Result    | RL    | Units | Date     | by  | Reference    |
|-------------------------|-----------|-------|-------|----------|-----|--------------|
| TCLP Mercury            | BDL       | 0.001 | mg/L  | 07/24/00 | RS  | E1311/E245.1 |
| Benzene                 | BDL       | 5.0   | ug/Kg | 07/25/00 | RM  | SW8021       |
| TCLP Silver             | BDL       | 0.01  | mg/L  | 07/25/00 | EK  | E1311/SW6010 |
| TCLP Arsenic            | BDL       | 0.01  | mg/L  | 07/25/00 | EK  | E1311/SW6010 |
| TCLP Barium             | 0.49      | 0.01  | mg/L  | 07/25/00 | EK  | E1311/SW6010 |
| TCLP Cadmium            | 0.008     | 0.005 | mg/L  | 07/25/00 | EK  | E1311/SW6010 |
| TCLP Chromium           | BDL       | 0.01  | mg/L  | 07/25/00 | EK  | E1311/SW6010 |
| TCLP Lead               | 5.46      | 0.015 | mg/L  | 07/25/00 | EK  | E1311/SW6010 |
| TCLP Selenium           | BDL       | 0.05  | mg/L  | 07/25/00 | EK  | E1311/SW6010 |
| Sulfur                  | 0.082     | 0.01  | %     | 08/11/00 | OL  | ASTMD129     |
| Total Organic Halogens  | ND        | 25    | mg/kg | 07/31/00 | OL  | SW9020       |
| Sonication Ext. For PCB | Completed |       |       | 07/20/00 | PL  | SW846-3550   |
| Soil Ext. for Semi- Vol | Completed |       |       | 07/20/00 | B/D | SW3550/3545  |
| Percent Solid           | 73        |       | %     | 07/21/00 | CF  | E160.3       |
| TCLP Digestion Mercury  | Completed |       |       | 07/24/00 | JA  | E1311/7470   |
| TCLP Extraction Metals  | Completed |       |       | 07/20/00 | T/A | EPA 1311     |
| Tot.Petroleum HC        | BDL       | 40    | mg/Kg | 07/26/00 | OL  | E418.1       |
| Total Cyanide           | 2.24      | 0.27  | mg/Kg | 07/25/00 | PJ  | SW9010       |

### Polychlorinated Biphenyls

|          |    |     |       |          |     |         |
|----------|----|-----|-------|----------|-----|---------|
| PCB-1016 | ND | 400 | ug/Kg | 07/21/00 | KCA | SW 8082 |
| PCB-1221 | ND | 400 | ug/Kg | 07/21/00 | KCA | SW 8082 |
| PCB-1232 | ND | 400 | ug/Kg | 07/21/00 | KCA | SW 8082 |
| PCB-1242 | ND | 400 | ug/Kg | 07/21/00 | KCA | SW 8082 |
| PCB-1248 | ND | 400 | ug/Kg | 07/21/00 | KCA | SW 8082 |

| Parameter                   | Result | RL   | Units | Date     | by  | Reference |
|-----------------------------|--------|------|-------|----------|-----|-----------|
| PCB-1254                    | ND     | 400  | ug/Kg | 07/21/00 | KCA | SW 8082   |
| PCB-1260                    | ND     | 400  | ug/Kg | 07/21/00 | KCA | SW 8082   |
| PCB-1262                    | ND     | 400  | ug/Kg | 07/21/00 | KCA | SW 8082   |
| PCB-1268                    | ND     | 400  | ug/Kg | 07/21/00 | KCA | SW 8082   |
| <b>Semivolatiles</b>        |        |      |       |          |     |           |
| 1,2,4-Trichlorobenzene      | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 1,2-Dichlorobenzene         | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 1,2-Diphenylhydrazine       | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 1,3-Dichlorobenzene         | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 1,4-Dichlorobenzene         | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 2,4,5-Trichlorophenol       | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 2,4,6-Trichlorophenol       | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 2,4-Dichlorophenol          | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 2,4-Dimethylphenol          | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 2,4-Dinitrophenol           | ND     | 1600 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 2,4-Dinitrotoluene          | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 2,6-Dichlorophenol          | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 2,6-Dinitrotoluene          | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 2-Chloronaphthalene         | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 2-Chlorophenol              | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 2-Methylnaphthalene         | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 2-Methylphenol (o-cresol)   | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 2-Nitroaniline              | ND     | 1600 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 2-Nitrophenol               | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 3,3'-Dichlorobenzidine      | ND     | 660  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 3-Nitroaniline              | ND     | 1600 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 4,6-Dinitro-2-methylphenol  | ND     | 1600 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 4-Bromophenyl phenyl ether  | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 4-Chloro-3-methylphenol     | ND     | 660  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 4-Chloroaniline             | ND     | 660  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 4-Chlorophenyl phenyl ether | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 4-Methylphenol (p-cresol)   | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 4-Nitroaniline              | ND     | 1600 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| 4-Nitrophenol               | ND     | 1600 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Acenaphthene                | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Acenaphthylene              | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Anthracene                  | 370    | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Benidine                    | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Benzo(a)anthracene          | 1700   | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Benzo(a)pyrene              | 1700   | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Benzo(b)fluoranthene        | 2800   | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Benzo(g,h,i)perylene        | ND     | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Benzo(k)fluoranthene        | 1800   | 330  | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Benzoic acid                | ND     | 1600 | ug/Kg | 07/21/00 | DRC | SW 8270   |




| Parameter                           | Result | RL  | Units | Date     | by  | Reference |
|-------------------------------------|--------|-----|-------|----------|-----|-----------|
| Benzyl alcohol                      | ND     | 660 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| zyl butyl phthalate                 | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Bis(2-chloroethoxy)methane          | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Bis(2-chloroethyl)ether             | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Bis(2-chloroisopropyl)ether         | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Bis(2-ethylhexyl)phthalate          | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Chrysene                            | 1300   | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Di-n-butylphthalate                 | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Di-n-octyl phthalate                | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Dibenz(a,h)anthracene               | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Dibenzofuran                        | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Diethyl phthalate                   | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Dimethyl phthalate                  | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Fluoranthene                        | 3100   | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Fluorene                            | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Hexachlorobenzene                   | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Hexachlorobutadiene                 | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Hexachlorocyclopentadiene           | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Hexachloroethane                    | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Indeno(1,2,3-c,d)pyrene             | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Isophorone                          | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Nitrosodi-n-propylamine             | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Nitrosodimethylamine                | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| N-Nitrosodiphenylamine              | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Naphthalene                         | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Nitrobenzene                        | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Pentachlorophenol                   | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Phenanthrene                        | 1400   | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Phenol                              | ND     | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| Pyrene                              | 2600   | 330 | ug/Kg | 07/21/00 | DRC | SW 8270   |
| % 2,4,6-Tribromophenol (Surrog Rec) | 51     |     | %     | 07/21/00 | DRC | SW 8270   |
| % 2-Fluorobiphenyl (Surrogate Rec)  | 49     |     | %     | 07/21/00 | DRC | SW 8270   |
| % 2-Fluorophenol (Surrogate Rec)    | 59     |     | %     | 07/21/00 | DRC | SW 8270   |
| % Nitrobenzene-d5 (Surrogate Rec)   | 62     |     | %     | 07/21/00 | DRC | SW 8270   |
| % Phenol-d5 (Surrogate Rec)         | 61     |     | %     | 07/21/00 | DRC | SW 8270   |
| % Terphenyl-d14 (Surrogate Rec)     | 81     |     | %     | 07/21/00 | DRC | SW 8270   |

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**Comments:**

ND=Not detected BDL = Below Detection Limit RL=Reporting Limit

If there are any questions regarding this data, please call Phoenix Client Services at extension 200.



Phyllis Shiller, Laboratory Director  
August 14, 2000



Environmental Laboratories, Inc.

567 East Middle Turnpike, P.O. Box 418 Manchester, CT 06040  
Tel. (860) 645-1102 Fax (860) 645-0823

## QC Report

AC80366

August 14, 2000

Sample ID AC80366

Analysis: AA Metals Analysis QC

AC80366

QC Source: ERA 9986  
NY 1706

Blank  
PPM

QC Check  
( % Rec.)

QC Spike  
( % Rec.)

QC Sample  
Replicate  
(% change)

Analyte

|             |         |     |     |    |
|-------------|---------|-----|-----|----|
| AS Arsenic  | <0.005  | 108 | 86  | NC |
| Hg Mercury  | <0.0005 | 108 | 103 | NC |
| Pb Lead     | <0.001  | 107 | 110 | NC |
| Sb Antimony | <0.005  | 104 | 85  | NC |
| Se Selenium | <0.005  | 99  | 110 | NC |
| Tl Thallium | <0.005  | 99  | 85  | NC |

Analysis: ICP Metals Analysis QC

AC80366

QC Source: ERA 9989  
 MIN QCI 702 Blank  
 ICP 0700  
 Analyte

QC Check  
 Sample  
 (% Rec.)

QC Spike  
 Sample  
 ( % Rec.)

QC Sample  
 Replicate  
 (% change)

| Analyte       | Blank | QC Check Sample (% Rec.) | QC Spike Sample ( % Rec.) | QC Sample Replicate (% change) |
|---------------|-------|--------------------------|---------------------------|--------------------------------|
| Ag Silver     | <0.01 | 98                       | 94                        | NC                             |
| Al Aluminum   | <0.05 | 97                       | 103                       | NC                             |
| As Arsenic    | <0.05 | 100                      | 116                       | NC                             |
| Ba Barium     | <0.01 | 100                      | 103                       | NC                             |
| Be Beryllium  | <0.01 | 99                       | 107                       | NC                             |
| Ca Calcium    | <0.10 | 100                      | 81                        | 0.2                            |
| Cd Cadmium    | <0.01 | 100                      | 111                       | NC                             |
| Co Cobalt     | <0.01 | 100                      | 104                       | NC                             |
| Cr Chromium   | <0.01 | 101                      | 109                       | NC                             |
| Cu Copper     | <0.01 | 99                       | 100                       | 11.5                           |
| Fe Iron       | <0.05 | 100                      | 104                       | 14.0                           |
| K Potassium   | <0.10 | 98                       | 110                       | 0.0                            |
| Mg Magnesium  | <0.01 | 102                      | 98                        | 0.0                            |
| Mn Manganese  | <0.01 | 101                      | 104                       | 0.0                            |
| Mo Molybdenum | <0.05 | 99                       | 91                        | NC                             |
| Na Sodium     | <0.10 | 100                      | 82                        | 0.5                            |
| Ni Nickel     | <0.01 | 100                      | 107                       | NC                             |
| P Phosphorus  | <0.05 | 92                       | -                         | NC                             |
| Pb Lead       | <0.01 | 100                      | 96                        | NC                             |
| Se Selenium   | <0.05 | 104                      | 116                       | NC                             |
| Sn Tin        | <0.25 | 103                      | 97                        | NC                             |
| V Vanadium    | <0.01 | 99                       | 112                       | NC                             |
| Zn Zinc       | <0.01 | 101                      | 121                       | NC                             |

Analysis: PCB QC

AC80366

| Analyte  | Method<br>Blank<br>(ppb) | Matrix<br>Spike<br>(% Rec ) | Matrix<br>Spike<br>Dup.<br>(% Rec) | RPD |
|----------|--------------------------|-----------------------------|------------------------------------|-----|
| PCB-1260 | ND                       | 97%                         | 94%                                | 3%  |

| Semivolatile<br>Analyte | Matrix<br>Spike<br>(%Rec) | Spike<br>Dup.<br>(%Rec) | % Diff.<br>(% D) |
|-------------------------|---------------------------|-------------------------|------------------|
| Phenol                  | 69                        | 68                      | 1                |
| 2-Chlorophenol          | 68                        | 72                      | 5                |
| 1,4-Dichlorobenzene     | 34                        | 36                      | 6                |
| N-Nitroso-di-n-prop.    | 52                        | 55                      | 7                |
| 1,2,4- Trichlorobenzene | 42                        | 45                      | 8                |
| 4-Chloro-3-methylphenol | 79                        | 70                      | 12               |
| Acenaphthene            | 60                        | 64                      | 7                |
| 2,4-Dinitrotoluene      | 65                        | 68                      | 3                |
| Pentachlorophenol       | 44                        | 36                      | 19               |
| Pyrene                  | 38                        | 43                      | 13               |

No target analytes were detected to the stated detection limits in the applicable method blanks with the following exceptions:

NONE

## Analysis: Total Cyanide Analysis QC

AC80366

QC BLANK:&lt;0.010

UNITS:MG/L

QC CHECK SAMPLE % RECOVERY:100

QC SOURCE:ULTRA#79754

QC SAMPLE SPIKE % RECOVERY:90.0

SPIKED SAMPLE:AC80384

QC SAMPLE REPLICATE % CHANGE:N/C

REPLICATED SAMPLE:AC80384

## Analysis: Volatile (GC) Analysis QC

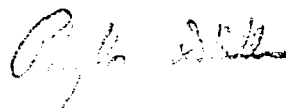
AC80366

| Analyte              | Matrix<br>Spike<br>(%Rec) | Matrix<br>Spike<br>Dup<br>(%Rec) | Relative<br>%Diff<br>(%D) |
|----------------------|---------------------------|----------------------------------|---------------------------|
| Benzene              | 108%                      | 111%                             | 3%                        |
| Chlorobenzene        | 94%                       | 96%                              | 2%                        |
| 1,1-Dichloroethylene | 108%                      | 111%                             | 17%                       |
| Toluene              | 84%                       | 84%                              | 0%                        |
| Trichloroethylene    | 94%                       | 90%                              | 4%                        |

No analytes were detected in the applicable method blanks above the stated detection limits with the following exceptions:

(NONE)

If there are any questions regarding this data, please call Phoenix Client Services at extension 200.

A handwritten signature in dark ink, appearing to read "Phyllis Shiller". The signature is written in a cursive, flowing style.

**Phyllis Shiller**  
**Laboratory Director**



6723 Towpath Road, P.O. Box 66  
Syracuse, New York 13214-0066  
TEL: (315) 446-9120

## CHAIN OF CUSTODY RECORD

10/23/95  
5951188L.CDR

**Distribution: Original Accompanies:**

**ent; Copy to Coordinator Field Files**

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Environmental Laboratories, Inc.  
587 East Middle Turnpike, P.O. Box 418, Manchester, CT 06040  
Tel. (860) 645-1102 Fax (860) 645-0823

## Analysis Report

August 02, 2000

FOR: Ms. Cathy Garaci  
Blasland & Bouck Engineers, PC  
6723 Towpath Road Box 66  
Syracuse NY 13214-0066

### Sample Information

Matrix: SOLID  
Location Code: BLASLAND  
Project Code:  
P.O.#: 36456

### Custody Information

Collected by: MC  
Received by: SW  
Analyzed by: see below

### Date

07/20/00  
07/21/00

### Time

10:40  
11:00

## Laboratory Data

Client ID: NMPC CEDAR ST SB-4A (7-13)

Phoenix I.D. AC80620

| Parameter                               | Result    | RL    | Units | Date     | by  | Reference    |
|---|-----------|-------|-------|----------|-----|--------------|
| TCLP Mercury                            | BDL       | 0.001 | mg/L  | 07/24/00 | RS  | E1311/E245.1 |
| zene                                    | ND        | 5.0   | ug/Kg | 07/21/00 | R/B | SW8021       |
| TCLP Silver                             | BDL       | 0.01  | mg/L  | 07/25/00 | TH  | E1311/SW6010 |
| TCLP Arsenic                            | 0.04      | 0.01  | mg/L  | 07/25/00 | TH  | E1311/SW6010 |
| TCLP Barium                             | 0.49      | 0.01  | mg/L  | 07/25/00 | TH  | E1311/SW6010 |
| TCLP Cadmium                            | BDL       | 0.005 | mg/L  | 07/25/00 | TH  | E1311/SW6010 |
| TCLP Chromium                           | BDL       | 0.01  | mg/L  | 07/25/00 | TH  | E1311/SW6010 |
| TCLP Lead                               | 0.096     | 0.015 | mg/L  | 07/25/00 | TH  | E1311/SW6010 |
| TCLP Selenium                           | BDL       | 0.05  | mg/L  | 07/25/00 | TH  | E1311/SW6010 |
| Sulfur                                  | 0.02      | 0.01  | %     | 08/01/00 | OL  | ASTMD129     |
| Total Organic Halogens                  | ND        | 25.0  | mg/kg | 08/01/00 | OL  | SW9020       |
| Sonication Ext. For PCB                 | Completed |       |       | 07/21/00 | PL  | SW846-3550   |
| Soil Ext. for Semi- Vol                 | Completed |       |       | 07/21/00 | PL  | SW3550/3545  |
| Percent Solid                           | 86        |       | %     | 07/21/00 | A/D | E160.3       |
| TCLP Digestion Mercury                  | Completed |       |       | 07/24/00 | JA  | E1311/7470   |
| TCLP Extraction Metals                  | Completed |       |       | 07/21/00 | PL  | EPA 1311     |
| Tot.Petroleum HC                        | BDL       | 23    | mg/Kg | 07/28/00 | EB  | E418.1       |
| Total Cyanide                           | BDL       | 0.21  | mg/Kg | 07/25/00 | PJ  | SW9010       |
| <b><u>Polychlorinated Biphenyls</u></b> |           |       |       |          |     |              |
| PCB-1016                                | ND        | 400   | ug/Kg | 07/24/00 | KCA | SW 8082      |
| PCB-1221                                | ND        | 400   | ug/Kg | 07/24/00 | KCA | SW 8082      |
| 3-1232                                  | ND        | 400   | ug/Kg | 07/24/00 | KCA | SW 8082      |
| PCB-1242                                | ND        | 400   | ug/Kg | 07/24/00 | KCA | SW 8082      |
| PCB-1248                                | ND        | 400   | ug/Kg | 07/24/00 | KCA | SW 8082      |

| Parameter                   | Result | RL   | Units | Date     | by  | Reference |
|-----------------------------|--------|------|-------|----------|-----|-----------|
| PCB-1254                    | ND     | 400  | ug/Kg | 07/24/00 | KCA | SW 8082   |
| PCB-1260                    | ND     | 400  | ug/Kg | 07/24/00 | KCA | SW 8082   |
| PCB-1262                    | ND     | 400  | ug/Kg | 07/24/00 | KCA | SW 8082   |
| PCB-1268                    | ND     | 400  | ug/Kg | 07/24/00 | KCA | SW 8082   |
| <b>Semivolatiles</b>        |        |      |       |          |     |           |
| 1,2,4-Trichlorobenzene      | ND     | 330  | ug/Kg | 07/24/00 | DRC | SW 8270   |
| 1,2-Dichlorobenzene         | ND     | 330  | ug/Kg | 07/24/00 | DRC | SW 8270   |
| 1,2-Diphenylhydrazine       | ND     | 330  | ug/Kg | 07/24/00 | DRC | SW 8270   |
| 1,3-Dichlorobenzene         | ND     | 330  | ug/Kg | 07/24/00 | DRC | SW 8270   |
| 1,4-Dichlorobenzene         | ND     | 330  | ug/Kg | 07/24/00 | DRC | SW 8270   |
| 2,4,5-Trichlorophenol       | ND     | 330  | ug/Kg | 07/24/00 | DRC | SW 8270   |
| 2,4,6-Trichlorophenol       | ND     | 330  | ug/Kg | 07/24/00 | DRC | SW 8270   |
| 2,4-Dichlorophenol          | ND     | 330  | ug/Kg | 07/24/00 | DRC | SW 8270   |
| 2,4-Dimethylphenol          | ND     | 330  | ug/Kg | 07/24/00 | DRC | SW 8270   |
| 2,4-Dinitrophenol           | ND     | 1600 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| 2,4-Dinitrotoluene          | ND     | 330  | ug/Kg | 07/24/00 | DRC | SW 8270   |
| 2,6-Dichlorophenol          | ND     | 330  | ug/Kg | 07/24/00 | DRC | SW 8270   |
| 2,6-Dinitrotoluene          | ND     | 330  | ug/Kg | 07/24/00 | DRC | SW 8270   |
| 2-Chloronaphthalene         | ND     | 330  | ug/Kg | 07/24/00 | DRC | SW 8270   |
| 2-Chlorophenol              | ND     | 330  | ug/Kg | 07/24/00 | DRC | SW 8270   |
| 2-Methylnaphthalene         | ND     | 330  | ug/Kg | 07/24/00 | DRC | SW 8270   |
| 2-Methylphenol (o-cresol)   | ND     | 330  | ug/Kg | 07/24/00 | DRC | SW 8270   |
| 2-Nitroaniline              | ND     | 1600 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| 2-Nitrophenol               | ND     | 330  | ug/Kg | 07/24/00 | DRC | SW 8270   |
| 3,3'-Dichlorobenzidine      | ND     | 660  | ug/Kg | 07/24/00 | DRC | SW 8270   |
| 3-Nitroaniline              | ND     | 1600 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| 4,6-Dinitro-2-methylphenol  | ND     | 1600 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| 4-Bromophenyl phenyl ether  | ND     | 330  | ug/Kg | 07/24/00 | DRC | SW 8270   |
| 4-Chloro-3-methylphenol     | ND     | 660  | ug/Kg | 07/24/00 | DRC | SW 8270   |
| 4-Chloroaniline             | ND     | 660  | ug/Kg | 07/24/00 | DRC | SW 8270   |
| 4-Chlorophenyl phenyl ether | ND     | 330  | ug/Kg | 07/24/00 | DRC | SW 8270   |
| 4-Methylphenol (p-cresol)   | ND     | 330  | ug/Kg | 07/24/00 | DRC | SW 8270   |
| 4-Nitroaniline              | ND     | 1600 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| 4-Nitrophenol               | ND     | 1600 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| Acenaphthene                | ND     | 330  | ug/Kg | 07/24/00 | DRC | SW 8270   |
| Acenaphthylene              | ND     | 330  | ug/Kg | 07/24/00 | DRC | SW 8270   |
| Anthracene                  | ND     | 330  | ug/Kg | 07/24/00 | DRC | SW 8270   |
| Benidine                    | ND     | 330  | ug/Kg | 07/24/00 | DRC | SW 8270   |
| Benzo(a)anthracene          | ND     | 330  | ug/Kg | 07/24/00 | DRC | SW 8270   |
| Benzo(a)pyrene              | ND     | 330  | ug/Kg | 07/24/00 | DRC | SW 8270   |
| Benzo(b)fluoranthene        | ND     | 330  | ug/Kg | 07/24/00 | DRC | SW 8270   |
| Benzo(g,h,i)perylene        | ND     | 330  | ug/Kg | 07/24/00 | DRC | SW 8270   |
| Benzo(k)fluoranthene        | ND     | 330  | ug/Kg | 07/24/00 | DRC | SW 8270   |
| Benzoic acid                | ND     | 1600 | ug/Kg | 07/24/00 | DRC | SW 8270   |

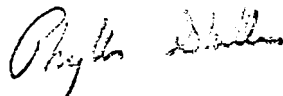
| Parameter                           | Result | RL  | Units | Date     | by  | Reference |
|-------------------------------------|--------|-----|-------|----------|-----|-----------|
| 1-Propyl alcohol                    | ND     | 660 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| 1-Propyl butyl phthalate            | ND     | 330 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| Bis(2-chloroethoxy)methane          | ND     | 330 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| Bis(2-chloroethyl)ether             | ND     | 330 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| Bis(2-chloroisopropyl)ether         | ND     | 330 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| Bis(2-ethylhexyl)phthalate          | ND     | 330 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| Chrysene                            | ND     | 330 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| Di-n-butylphthalate                 | ND     | 330 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| Di-n-octyl phthalate                | ND     | 330 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| Dibenz(a,h)anthracene               | ND     | 330 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| Dibenzofuran                        | ND     | 330 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| Diethyl phthalate                   | ND     | 330 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| Dimethyl phthalate                  | ND     | 330 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| Fluoranthene                        | 330    | 330 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| Fluorene                            | ND     | 330 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| Hexachlorobenzene                   | ND     | 330 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| Hexachlorobutadiene                 | ND     | 330 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| Hexachlorocyclopentadiene           | ND     | 330 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| Hexachloroethane                    | ND     | 330 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| Indeno(1,2,3-c,d)pyrene             | ND     | 330 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| Isophorone                          | ND     | 330 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| N-Nitrosodi-n-propylamine           | ND     | 330 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| N-Nitrosodimethylamine              | ND     | 330 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| N-Nitrosodiphenylamine              | ND     | 330 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| Naphthalene                         | ND     | 330 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| Nitrobenzene                        | ND     | 330 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| Pentachlorophenol                   | ND     | 330 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| Phenanthrene                        | ND     | 330 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| Phenol                              | ND     | 330 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| Pyrene                              | ND     | 330 | ug/Kg | 07/24/00 | DRC | SW 8270   |
| % 2,4,6-Tribromophenol (Surrog Rec) | 52     |     | %     | 07/24/00 | DRC | SW 8270   |
| % 2-Fluorobiphenyl (Surrogate Rec)  | 57     |     | %     | 07/24/00 | DRC | SW 8270   |
| % 2-Fluorophenol (Surrogate Rec)    | 73     |     | %     | 07/24/00 | DRC | SW 8270   |
| % Nitrobenzene-d5 (Surrogate Rec)   | 70     |     | %     | 07/24/00 | DRC | SW 8270   |
| % Phenol-d5 (Surrogate Rec)         | 71     |     | %     | 07/24/00 | DRC | SW 8270   |
| % Terphenyl-d14 (Surrogate Rec)     | 85     |     | %     | 07/24/00 | DRC | SW 8270   |

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**Comments:**

ND=Not detected BDL = Below Detection Limit RL=Reporting Limit

If there are any questions regarding this data, please call Phoenix Client Services at extension 200.



Phyllis Shiller, Laboratory Director  
August 02, 2000



Environmental Laboratories, Inc.  
587 East Middle Turnpike, P.O. Box 418, Manchester, CT 06040  
Tel. (860) 645-1102 Fax (860) 645-0823

## Analysis Report

August 02, 2000

FOR: Ms. Cathy Garaci  
Blasland & Bouck Engineers, PC  
6723 Towpath Road Box 66  
Syracuse NY 13214-0066

### Sample Information

Matrix: SOLID  
Location Code: BLASLAND  
Project Code:  
P.O.#: 36456

### Custody Information

Collected by: MC  
Received by: SW  
Analyzed by: see below

### Date

07/20/00  
07/21/00

### Time

11:30  
11:00

## Laboratory Data

Client ID: NMPC CEDAR ST SB-5 (0-8)

Phoenix I.D. AC80621

| Parameter                               | Result    | RL    | Units | Date     | by  | Reference    |
|---|-----------|-------|-------|----------|-----|--------------|
| TCLP Mercury                            | BDL       | 0.001 | mg/L  | 07/24/00 | RS  | E1311/E245.1 |
| zene                                    | ND        | 5.0   | ug/Kg | 07/21/00 | R/B | SW8021       |
| TCLP Silver                             | BDL       | 0.01  | mg/L  | 07/25/00 | TH  | E1311/SW6010 |
| TCLP Arsenic                            | BDL       | 0.01  | mg/L  | 07/25/00 | TH  | E1311/SW6010 |
| TCLP Barium                             | 1.64      | 0.01  | mg/L  | 07/25/00 | TH  | E1311/SW6010 |
| TCLP Cadmium                            | 0.006     | 0.005 | mg/L  | 07/25/00 | TH  | E1311/SW6010 |
| TCLP Chromium                           | BDL       | 0.01  | mg/L  | 07/25/00 | TH  | E1311/SW6010 |
| TCLP Lead                               | 0.147     | 0.015 | mg/L  | 07/25/00 | TH  | E1311/SW6010 |
| TCLP Selenium                           | BDL       | 0.05  | mg/L  | 07/25/00 | TH  | E1311/SW6010 |
| Sulfur                                  | 0.08      | 0.01  | %     | 08/01/00 | OL  | ASTMD129     |
| Total Organic Halogens                  | ND        | 25.0  | mg/kg | 08/01/00 | OL  | SW9020       |
| Sonication Ext. For PCB                 | Completed |       |       | 07/21/00 | PL  | SW846-3550   |
| Soil Ext. for Semi- Vol                 | Completed |       |       | 07/21/00 | PL  | SW3550/3545  |
| Percent Solid                           | 87        |       | %     | 07/21/00 | A/D | E160.3       |
| TCLP Digestion Mercury                  | Completed |       |       | 07/24/00 | JA  | E1311/7470   |
| TCLP Extraction Metals                  | Completed |       |       | 07/21/00 | PL  | EPA 1311     |
| Tot.Petroleum HC                        | 540       | 46    | mg/Kg | 07/28/00 | EB  | E418.1       |
| Total Cyanide                           | 1.12      | 0.22  | mg/Kg | 07/25/00 | PJ  | SW9010       |
| <b><u>Polychlorinated Biphenyls</u></b> |           |       |       |          |     |              |
| PCB-1016                                | ND        | 400   | ug/Kg | 07/24/00 | KCA | SW 8082      |
| PCB-1221                                | ND        | 400   | ug/Kg | 07/24/00 | KCA | SW 8082      |
| -1232                                   | ND        | 400   | ug/Kg | 07/24/00 | KCA | SW 8082      |
| PCB-1242                                | ND        | 400   | ug/Kg | 07/24/00 | KCA | SW 8082      |
| PCB-1248                                | ND        | 400   | ug/Kg | 07/24/00 | KCA | SW 8082      |

| Parameter                   | Result | RL   | Units | Date     | by  | Reference |
|-----------------------------|--------|------|-------|----------|-----|-----------|
| PCB-1254                    | ND     | 400  | ug/Kg | 07/24/00 | KCA | SW 8082   |
| PCB-1260                    | ND     | 400  | ug/Kg | 07/24/00 | KCA | SW 8082   |
| PCB-1262                    | ND     | 400  | ug/Kg | 07/24/00 | KCA | SW 8082   |
| PCB-1268                    | ND     | 400  | ug/Kg | 07/24/00 | KCA | SW 8082   |
| <b><u>Semivolatiles</u></b> |        |      |       |          |     |           |
| 1,2,4-Trichlorobenzene      | ND     | 330  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 1,2-Dichlorobenzene         | ND     | 330  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 1,2-Diphenylhydrazine       | ND     | 330  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 1,3-Dichlorobenzene         | ND     | 330  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 1,4-Dichlorobenzene         | ND     | 330  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2,4,5-Trichlorophenol       | ND     | 330  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2,4,6-Trichlorophenol       | ND     | 330  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2,4-Dichlorophenol          | ND     | 330  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2,4-Dimethylphenol          | ND     | 330  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2,4-Dinitrophenol           | ND     | 1600 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2,4-Dinitrotoluene          | ND     | 330  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2,6-Dichlorophenol          | ND     | 330  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2,6-Dinitrotoluene          | ND     | 330  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2-Chloronaphthalene         | ND     | 330  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2-Chlorophenol              | ND     | 330  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2-Methylnaphthalene         | ND     | 330  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2-Methylphenol (o-cresol)   | ND     | 330  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2-Nitroaniline              | ND     | 1600 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 2-Nitrophenol               | ND     | 330  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 3,3'-Dichlorobenzidine      | ND     | 660  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 3-Nitroaniline              | ND     | 1600 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 4,6-Dinitro-2-methylphenol  | ND     | 1600 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 4-Bromophenyl phenyl ether  | ND     | 330  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 4-Chloro-3-methylphenol     | ND     | 660  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 4-Chloroaniline             | ND     | 660  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 4-Chlorophenyl phenyl ether | ND     | 330  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 4-Methylphenol (p-cresol)   | ND     | 330  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 4-Nitroaniline              | ND     | 1600 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| 4-Nitrophenol               | ND     | 1600 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Acenaphthene                | ND     | 330  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Acenaphthylene              | ND     | 330  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Anthracene                  | ND     | 330  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Benzidine                   | ND     | 330  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Benzo(a)anthracene          | 1100   | 330  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Benzo(a)pyrene              | 1200   | 330  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Benzo(b)fluoranthene        | 1500   | 330  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Benzo(g,h,i)perylene        | ND     | 330  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Benzo(k)fluoranthene        | 2400   | 330  | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Benzoic acid                | ND     | 1600 | ug/Kg | 07/22/00 | DRC | SW 8270   |



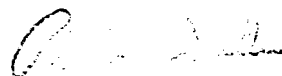
| Parameter                           | Result | RL  | Units | Date     | by  | Reference |
|-------------------------------------|--------|-----|-------|----------|-----|-----------|
| Benzyl alcohol                      | ND     | 660 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Benzyl butyl phthalate              | ND     | 330 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Bis(2-chloroethoxy)methane          | ND     | 330 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Bis(2-chloroethyl)ether             | ND     | 330 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Bis(2-chloroisopropyl)ether         | ND     | 330 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Bis(2-ethylhexyl)phthalate          | ND     | 330 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Chrysene                            | 650    | 330 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Di-n-butylphthalate                 | ND     | 330 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Di-n-octyl phthalate                | ND     | 330 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Dibenz(a,h)anthracene               | ND     | 330 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Dibenzofuran                        | ND     | 330 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Diethyl phthalate                   | ND     | 330 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Dimethyl phthalate                  | ND     | 330 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Fluoranthene                        | 1200   | 330 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Fluorene                            | ND     | 330 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Hexachlorobenzene                   | ND     | 330 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Hexachlorobutadiene                 | ND     | 330 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Hexachlorocyclopentadiene           | ND     | 330 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Hexachloroethane                    | ND     | 330 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Indeno(1,2,3-c,d)pyrene             | ND     | 330 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Isophorone                          | ND     | 330 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| N-Nitrosodi-n-propylamine           | ND     | 330 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| N-Nitrosodimethylamine              | ND     | 330 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| N-Nitrosodiphenylamine              | ND     | 330 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Naphthalene                         | ND     | 330 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Nitrobenzene                        | ND     | 330 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Pentachlorophenol                   | ND     | 330 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Phenanthrene                        | 630    | 330 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Phenol                              | ND     | 330 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| Pyrene                              | 990    | 330 | ug/Kg | 07/22/00 | DRC | SW 8270   |
| % 2,4,6-Tribromophenol (Surrog Rec) | 33     |     | %     | 07/22/00 | DRC | SW 8270   |
| % 2-Fluorobiphenyl (Surrogate Rec)  | 29     |     | %     | 07/22/00 | DRC | SW 8270   |
| % 2-Fluorophenol (Surrogate Rec)    | 38     |     | %     | 07/22/00 | DRC | SW 8270   |
| % Nitrobenzene-d5 (Surrogate Rec)   | 36     |     | %     | 07/22/00 | DRC | SW 8270   |
| % Phenol-d5 (Surrogate Rec)         | 37     |     | %     | 07/22/00 | DRC | SW 8270   |
| % Terphenyl-d14 (Surrogate Rec)     | 41     |     | %     | 07/22/00 | DRC | SW 8270   |

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**Comments:**

ND=Not detected BDL = Below Detection Limit RL=Reporting Limit

If there are any questions regarding this data, please call Phoenix Client Services at extension 200.



Phyllis Shiller, Laboratory Director  
August 02, 2000



Environmental Laboratories, Inc.

587 East Middle Turnpike, P.O. Box 418, Manchester, CT 06040  
Tel. (860) 645-1102 Fax (860) 645-0823

## QC Report

AC80622

August 02, 2000

Sample ID AC80622

Analysis: AA Metals Analysis QC

AC80622

|                                |              |                       |                       |                                      |
|--------------------------------|--------------|-----------------------|-----------------------|--------------------------------------|
| QC Source: ERA 9986<br>NY 1706 | Blank<br>PPM | QC Check<br>( % Rec.) | QC Spike<br>( % Rec.) | QC Sample<br>Replicate<br>(% change) |
| Analyte                        |              |                       |                       |                                      |
| AS Arsenic                     | <0.005       | 108                   | 86                    | NC                                   |
| Hg Mercury                     | <0.0005      | 108                   | 103                   | NC                                   |
| Pb Lead                        | <0.001       | 107                   | 110                   | NC                                   |
| Sb Antimony                    | <0.005       | 104                   | 85                    | NC                                   |
| Se Selenium                    | <0.005       | 99                    | 110                   | NC                                   |
| Tl Thallium                    | <0.005       | 99                    | 85                    | NC                                   |

Analysis: ICP Metals Analysis QC

AC80622

|  |       |                                |                                 |                                      |
|--|-------|--------------------------------|---------------------------------|--------------------------------------|
| QC Source: ERA 9989<br>MIN QCI 702<br>ICP 0700 | Blank | QC Check<br>Sample<br>(% Rec.) | QC Spike<br>Sample<br>( % Rec.) | QC Sample<br>Replicate<br>(% change) |
| Analyte  |       |                                |                                 |                                      |
| Ag Silver                                      | <0.01 | 98                             | 94                              | NC                                   |
| Al Aluminum                                    | <0.05 | 97                             | 103                             | NC                                   |
| As Arsenic                                     | <0.05 | 100                            | 116                             | NC                                   |
| Ba Barium                                      | <0.01 | 100                            | 103                             | NC                                   |
| Be Beryllium                                   | <0.01 | 99                             | 107                             | NC                                   |
| Ca Calcium                                     | <0.10 | 100                            | 81                              | 0.2                                  |
| Cd Cadmium                                     | <0.01 | 100                            | 111                             | NC                                   |
| Co Cobalt                                      | <0.01 | 100                            | 104                             | NC                                   |
| Cr Chromium                                    | <0.01 | 101                            | 109                             | NC                                   |
| Cu Copper                                      | <0.01 | 99                             | 100                             | 11.5                                 |
| Fe Iron  | <0.05 | 100                            | 104                             | 14.0                                 |
| K Potassium                                    | <0.10 | 98                             | 110                             | 0.0                                  |
| Mg Magnesium                                   | <0.01 | 102                            | 98                              | 0.0                                  |
| Mn Manganese                                   | <0.01 | 101                            | 104                             | 0.0                                  |
| Mo Molybdenum                                  | <0.05 | 99                             | 91                              | NC                                   |
| Na Sodium                                      | <0.10 | 100                            | 82                              | 0.5                                  |
| Ni Nickel                                      | <0.01 | 100                            | 107                             | NC                                   |
| P Phosphorus                                   | <0.05 | 92                             | -                               | NC                                   |
| Pb Lead  | <0.01 | 100                            | 96                              | NC                                   |
| Se Selenium                                    | <0.05 | 104                            | 116                             | NC                                   |
| Sn Tin   | <0.25 | 103                            | 97                              | NC                                   |
| V Vanadium                                     | <0.01 | 99                             | 112                             | NC                                   |
| Zn Zinc  | <0.01 | 101                            | 121                             | NC                                   |

Analysis: PCB QC

AC80622

| Analyte  | Method<br>Blank<br>(ppb) | Matrix<br>Spike<br>(% Rec ) | Matrix<br>Spike<br>Dup.<br>(% Rec) | RPD |
|----------|--------------------------|-----------------------------|------------------------------------|-----|
| PCB-1260 | ND                       | 97%                         | 94%                                | 3%  |

Analysis: Semivolatile (MS) Analysis QC

AC80622

| Semivolatile<br>Analyte | Matrix<br>Spike<br>(%Rec) | Spike<br>Dup.<br>(%Rec) | % Diff.<br>(% D) |
|-------------------------|---------------------------|-------------------------|------------------|
| Phenol                  | 69                        | 68                      | 1                |
| 2-Chlorophenol          | 68                        | 72                      | 5                |
| 1,4-Dichlorobenzene     | 34                        | 36                      | 6                |
| N-Nitroso-di-n-prop.    | 52                        | 55                      | 7                |
| 1,2,4- Trichlorobenzene | 42                        | 45                      | 8                |
| 4-Chloro-3-methylphenol | 79                        | 70                      | 12               |
| Acenaphthene            | 60                        | 64                      | 7                |
| 2,4-Dinitrotoluene      | 65                        | 68                      | 3                |
| Pentachlorophenol       | 44                        | 36                      | 19               |
| Pyrene                  | 38                        | 43                      | 13               |

No target analytes were detected to the stated detection limits in the applicable method blanks with the following exceptions:

NONE

Analysis: Total Cyanide Analysis QC

AC80622

|                                  |                           |
|----------------------------------|---------------------------|
| QC BLANK:<0.010                  | UNITS:MG/L                |
| QC CHECK SAMPLE % RECOVERY:100   | QC SOURCE:ULTRA#79754     |
| QC SAMPLE SPIKE % RECOVERY:76.0  | SPIKED SAMPLE:AC80379     |
| QC SAMPLE REPLICATE % CHANGE:N/C | REPLICATED SAMPLE:AC80379 |

Analysis: Total Petroleum HC/IR Q

AC80622

|                                   |                            |
|-----------------------------------|----------------------------|
| QC BLANK: <20.0                   | UNITS: mg/KG               |
| QC CHECK SAMPLE % RECOVERY:104    | QC SOURCE: IN HOUSE        |
| SP SAMPLE SPIKE % RECOVERY: 94.7  | SPIKE SAMPLE: AC80620      |
| QC SAMPLE REPLICATE % CHANGE: N/C | REPLICATED SAMPLE: AC80620 |

| Analyte              | Matrix<br>Spike<br>(%Rec) | Matrix<br>Spike<br>Dup<br>(%Rec) | Relative<br>%Diff<br>(%D) |
|----------------------|---------------------------|----------------------------------|---------------------------|
| Benzene              | 108%                      | 111%                             | 3%                        |
| Chlorobenzene        | 94%                       | 96%                              | 2%                        |
| 1,1-Dichloroethylene | 108%                      | 111%                             | 17%                       |
| Toluene              | 84%                       | 84%                              | 0%                        |
| Trichloroethylene    | 94%                       | 90%                              | 4%                        |

No analytes were detected in the applicable method blanks above the stated detection limits with the following exceptions:

(NONE)

If there are any questions regarding this data, please call Phoenix Client Services at extension 200.

Phyllis Shiller  
Laboratory Director



6723 Towpath Road, P.O. Box 66  
Syracuse, New York 13214-0066  
TEL: (315) 446-9120

Fedex A3 # 8166 0668 4328

## CHAIN OF CUSTODY RECORD

[illegible]









## INORGANIC ANALYTICAL REPORT

Client : Blasland, Bouck & Lee  
Account # : 10624  
Site : NMPC ONEIDA CEDAR ST

Date Received : 20-JUL-00  
Date Sampled : 20-JUL-00

Matrix : Soil

Galson ID: L62040-1  
Client ID: SB-5 (8-10)

|                  | Method     | Units |      |
|------------------|------------|-------|------|
| Corrosivity/pH   | SW846 9045 | SU    | 7.4  |
| Reactive Cyanide | SW846 Ch.  | mg/kg | <100 |
| Ignitability     | SW846 1030 |       | NEG  |
| Reactive Sulfide | SW846 Ch.  | mg/kg | <100 |

Approved by : DM

Date : 09-AUG-00

QC by : *[Signature]*

Date : 8/10/00

NYS DOH # : 11626

Footnotes:

- \* The bulk pH was performed using SW846 method 9045. A sample is corrosive if pH is less than or equal to 2, or greater than or equal to 12.5 Standard Units (SU). Under these conditions, the sample is not corrosive.
- \* The sample does not exceed the USEPA action levels of 250mg HCN/kg waste and/or 500 mg H<sub>2</sub>S/kg waste as stated in SW846; therefore it is not reactive.
- \* The sample does not ignite or support combustion. Under these conditions the sample is non-ignitable.





## METALS ANALYTICAL REPORT

Client : Blasland, Bouck & Lee  
Account # : 10624  
Site : NMPC ONEIDA CEDAR ST

Date Received : 20-JUL-00  
Date Sampled : 20-JUL-00

Matrix : Leachate  
Method : SW846 6010B/ 7470A; ICP/ CVAA

| Galson ID:    | L62040-1     | WG24230-1    | WG24362-1    |
|---------------|--------------|--------------|--------------|
| Client ID:    | SB-5 (8-10)  | Method Blank | Method Blank |
|               | Units        |              |              |
| Mercury TCLP  | mg/l <0.0003 | NR           | <0.0003      |
| Arsenic TCLP  | mg/l <0.01   | <0.01        | NR           |
| Barium TCLP   | mg/l 1.3     | <1           | NR           |
| Cadmium TCLP  | mg/l 0.037   | <0.005       | NR           |
| Chromium TCLP | mg/l <0.01   | <0.01        | NR           |
| Lead TCLP     | mg/l 0.78    | <0.02        | NR           |
| Selenium TCLP | mg/l <0.02   | <0.02        | NR           |
| Silver TCLP   | mg/l <0.01   | <0.01        | NR           |

Approved by : LK  
Date : 01-AUG-00  
QC by : *[Signature]*  
Date : 8/10/00  
NYS DOH # : 11626  
Footnotes:



Contract:

SDG No.: L62040

S1 (2FP) = 2-Fluorophenol  
S2 (PHL) = Phenol-d6  
S3 (NBZ) = Nitrobenzene-d5  
S4 (FBP) = 2-Fluorobiphenyl  
S5 (TBP) = 2,4,6-Tribromophenol  
S6 (TPH) = Terphenyl-d14  
S7 (2CP) = 2-Chlorophenol-d4  
S8 (DCB) = 1,2-Dichlorobenzene-d4

(23-120)  
(30-146)  
(42-132)  
(32-127)  
(31-134)  
(43-132)  
(38-135) (advisory)  
(39-132) (advisory)

page 1 of 1



## SEMIVOLATILE ANALYTICAL REPORT

Client : Blasland, Bouck & Lee  
Account # : 10624  
Site : NMPC ONEIDA CEDAR ST

Date Received : 20-JUL-00  
Date Sampled : 20-JUL-00  
Date Extracted : 31-JUL-00

Matrix : Leachate  
Method : SW846 1311 8270-TCLP  
Units : ug/l

| Galson ID: | L62040-1    | WG24409-1    | WG24409-2  |
|------------|-------------|--------------|------------|
| Client ID: | SB-5 (8-10) | Method Blank | TCLP Blank |

|                       |      |      |      |
|-----------------------|------|------|------|
| Pyridine              | <100 | <100 | <100 |
| 1,4-Dichlorobenzene   | <100 | <100 | <100 |
| 2-Methylphenol        | <100 | <100 | <100 |
| 3 & 4-Methylphenol    | <200 | <200 | <200 |
| Hexachloroethane      | <100 | <100 | <100 |
| Nitrobenzene          | <100 | <100 | <100 |
| Hexachlorobutadiene   | <100 | <100 | <100 |
| 2,4,6-Trichlorophenol | <100 | <100 | <100 |
| 2,4,5-Trichlorophenol | <100 | <100 | <100 |
| 2,4-Dinitrotoluene    | <100 | <100 | <100 |
| Hexachlorobenzene     | <100 | <100 | <100 |
| Pentachlorophenol     | <250 | <250 | <250 |

|                 |          |          |          |
|-----------------|----------|----------|----------|
| Analysis Date   | 08/03/00 | 08/03/00 | 08/03/00 |
| Dilution Factor | 1        | 1        | 1        |

Approved by : Oommen Kappil  
Date : 10-AUG-00  
QC by : *[Signature]*  
Date : 8/14/00  
NYS DOH # : 11626  
Footnotes:

TCLP extraction performed 7/24/00.



## LEACHATE VOLATILE SURROGATE RECOVERY

Client : Blasland, Bouck & Lee

Login # : L62040

[illegible]

SMC1 (TOL) = Toluene-d8  
SMC2 (BFB) = Bromofluorobenzene  
SMC3 (DCE) = 1,2-Dichloroethane-d4

QC LIMITS  
(74-123)  
(72-118)  
(58-134)

```
# Column to be used to flag recovery values
* Values outside of QC limits
D Surrogate diluted out
```



## VOLATILE ANALYTICAL REPORT

Client : Blasland, Bouck & Lee  
Account # : 10624  
Site : NMPC ONEIDA CEDAR ST

Date Received : 20-JUL-00  
Date Sampled : 20-JUL-00  
Date Extracted: NA

Matrix : Leachate  
Method : SW846 1311 8260  
Units : ug/l

| Galson ID:           | *L62040-1   | WG24361-1    | WG24361-2  |
|----------------------|-------------|--------------|------------|
| Client ID:           | SB-5 (8-10) | Method Blank | TCLP Blank |
| Benzene              | <50         | <50          | <50        |
| 2-Butanone           | 110         | <100         | 100        |
| Carbon Tetrachloride | <50         | <50          | <50        |
| Chlorobenzene        | <50         | <50          | <50        |
| Chloroform           | <50         | <50          | <50        |
| 1,2-Dichloroethane   | <50         | <50          | <50        |
| 1,1-Dichloroethene   | <50         | <50          | <50        |
| Tetrachloroethene    | <50         | <50          | <50        |
| Trichloroethene      | <50         | <50          | <50        |
| Vinyl Chloride       | <50         | <50          | <50        |
| Analysis Date        | 07/28/00    | 07/28/00     | 07/28/00   |
| Dilution Factor      | 10          | 10           | 10         |

Approved by : PJT  
Date : 10-AUG-00  
QC by : *[Signature]*  
Date : 8/10/00  
NYS DOH # : 11626  
Footnotes:

TCLP extraction performed 07/24/00.

\* : 2-Butanone was also detected in the TCLP blank, and should not be considered as originally present in the sample.





**6601 Kirkville Road East  
E. Syracuse, New York 13057  
315 437-7252 • 888-577-5227**

Company Name

BBL, Inc.

Project Name / Number

36456

NMPC Over the Cedar St

### Turn-Around Time

☒ - Standard Service

☐ - \* Rush Service

Date requested by:

Ph # 3151-446 4120

Fax # ( ) -

Send Report to: Cathy Geraci  
BBL, Inc.

Send Invoice to: Bills Payable  
RBI

P.O. # 36456

Page 3 of 3

### PARAMETERS FOR ANALYSIS

[illegible]

REMARKS: Samples stored on ice. Hand delivered

Total Containers - 3

SAMPLER'S NAME: Michael Cobb

SIGNATURE: Yh [Signature]

| VOC Pres | U | P | AU | NA |
|----------|---|---|----|----|
|----------|---|---|----|----|

**SAMPLES RELINQUISHED BY:**

**SAMPLES RECEIVED BY:**

Custody Seal Intact? ☐ Yes ☐ No ☐ N.A.

Shipment Complete? ☐ Yes ☐ No

NAME: Michael Cohn DATE: 7/20/00  
SIGNATURE: [Signature] TIME: 1925

NAME: Shawn S Noor DATE: 7/20/00  
SIGNATURE: Shawn S Noor TIME: 4:20 pm

Temp 4.7 °C TS TB TM

NAME: \_\_\_\_\_ DATE: \_\_\_\_\_  
SIGNATURE: \_\_\_\_\_ TIME: \_\_\_\_\_

Received For Laboratory By: \_\_\_\_\_ DATE: \_\_\_\_\_  
(Signature) TIME: \_\_\_\_\_

Airbill # 5ano delivered

NAME: \_\_\_\_\_ DATE: \_\_\_\_\_  
SIGNATURE: \_\_\_\_\_ TIME: \_\_\_\_\_

Received For Laboratory By: DATE:  
(Signature) TIME:



1 Mustard ST.  
Suite 250  
Rochester, NY 14609

**THIS IS AN ANALYTICAL TEST REPORT FOR:**

Client : Galson Laboratories  
Project Reference: L62040  
Lab Submission # : R2003092  
Reported : 08/09/00

Report Contains a total of 8 pages

The results reported herein relate only to the samples received by the laboratory. This report may not be reproduced except in full, without the approval of Columbia Analytical Services.

This package has been reviewed by Columbia Analytical Services' QA Department/Laboratory Director to comply with NELAC standards prior to report submittal. *Michael E. Perry*

00001





# CASE NARRATIVE

This report contains analytical results for the following samples:

Submission #: R2003092

Lab ID

396317

Client ID

SB-5 (8-10)

All samples were received in good condition.

All samples have been analyzed by the approved methods cited on the analytical results pages.

All holding times and associated QC were within limits.

No analytical or QC problems were encountered.



Effective 04/01/96

### CAS LIST OF QUALIFIERS

(The basis of this proposal are the EPA-CLP Qualifiers)

- U - Indicates compound was analyzed for but was not detected. The sample quantitation limit must be corrected for dilution and for percent moisture.
- J - Indicates an estimated value. For further explanation see case narrative / cover letter.
- 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.
- A - This flag indicates that a TIC is a suspected aldol-condensation product.
- N - Spiked sample recovery not within control limits.  
(Flag the entire batch - Inorganic analysis only)
- \* - Duplicate analysis not within control limits.  
(Flag the entire batch - Inorganic analysis only)
- Also used to qualify Organics QC data outside limits.
- D - Spike diluted out.
- S - Reported value determined by Method of Standard Additions. (MSA)
- X - As specified in the case narrative.

### **CAS Lab ID # for State Certifications**

NY ID # in Rochester: 10145  
CT ID # in Rochester: PH0556  
MA ID # in Rochester: M-NY032  
OH EPA # in Rochester: VAP

NJ ID # in Rochester: 73004  
RI ID # in Rochester: 158  
NH ID # in Rochester: 294198-A  
AIHA # in Rochester: 7889

COLUMBIA ANALYTICAL SERVICES**EXTRACTABLE ORGANICS**

METHOD 8081A TCLP

Reported: 08/09/00

Galson Laboratories

Project Reference: L62040

Client Sample ID : SB-5 (8-10)

Date Sampled : 07/20/00 11:40 Order #: 396317      Sample Matrix: SOIL/SEDIMENT  
Date Received: 07/26/00 Submission #: R2003092      Analytical Run 53569

| ANALYTE              | PQL        | RESULT | UNITS |
|----------------------|------------|--------|-------|
| DATE EXTRACTED       | : 07/27/00 |        |       |
| DATE ANALYZED        | : 08/07/00 |        |       |
| ANALYTICAL DILUTION: | 10.00      |        |       |
| GAMMA-BHC (LINDANE)  | 0.50       | 5.0 U  | UG/L  |
| CHLORDANE            | 2.0        | 20 U   | UG/L  |
| ENDRIN               | 0.50       | 5.0 U  | UG/L  |
| HEPTACHLOR           | 0.50       | 5.0 U  | UG/L  |
| HEPTACHLOR EPOXIDE   | 0.50       | 5.0 U  | UG/L  |
| METHOXYCHLOR         | 2.0        | 20 U   | UG/L  |
| TOXAPHENE            | 10         | 100 U  | UG/L  |

SURROGATE RECOVERIESQC LIMITS

|                                |              |    |   |
|--------------------------------|--------------|----|---|
| DECACHLOROBIPHENYL (DCB)       | (24 - 154 %) | 79 | % |
| TETRACHLORO-META-XYLENE (TCMX) | (30 - 150 %) | 96 | % |

Data Reported following TCLP Toxicity Characteristics Leaching Procedure.  
Federal Register, Part 261, Vol. 55, NO 126, June 29, 1990.

00004

COLUMBIA ANALYTICAL SERVICES**EXTRACTABLE ORGANICS**

METHOD 8151A

Reported: 08/09/00

Galson Laboratories

Project Reference: L62040

Client Sample ID : SB-5 (8-10)

Date Sampled : 07/20/00 11:40 Order #: 396317      Sample Matrix: SOIL/SEDIMENT  
Date Received: 07/26/00 Submission #: R2003092      Analytical Run 53627

| ANALYTE | PQL | RESULT | UNITS |
|---------|-----|--------|-------|
|---------|-----|--------|-------|

DATE EXTRACTED : 07/28/00  
DATE ANALYZED : 08/05/00  
ANALYTICAL DILUTION: 100.00

|                   |      |      |      |
|-------------------|------|------|------|
| 2,4-D             | 0.50 | 50 U | UG/L |
| 2,4,5-TP (SILVEX) | 0.50 | 50 U | UG/L |

SURROGATE RECOVERIESQC LIMITS

|      |              |     |   |
|------|--------------|-----|---|
| DCAA | (18 - 152 %) | 111 | % |
|------|--------------|-----|---|

Data Reported following TCLP Toxicity Characteristics Leaching Procedure.  
Federal Register, Part 261, Vol. 55, NO 126, June 29, 1990.

00005

COLUMBIA ANALYTICAL SERVICES**EXTRACTABLE ORGANICS**

METHOD 8081A TCLP

Reported: 08/09/00

**Project Reference:**

Client Sample ID : METHOD BLANK

|                |                 |                              |
|----------------|-----------------|------------------------------|
| Date Sampled : | Order #: 398260 | Sample Matrix: SOIL/SEDIMENT |
| Date Received: | Submission #:   | Analytical Run 53569         |

| ANALYTE | PQL | RESULT | UNITS |
|---------|-----|--------|-------|
|---------|-----|--------|-------|

|                      |            |
|----------------------|------------|
| DATE EXTRACTED       | : 07/27/00 |
| DATE ANALYZED        | : 08/05/00 |
| ANALYTICAL DILUTION: | 10.00      |

|                     |      |       |      |
|---------------------|------|-------|------|
| GAMMA-BHC (LINDANE) | 0.50 | 5.0 U | UG/L |
| CHLORDANE           | 2.0  | 20 U  | UG/L |
| ENDRIN              | 0.50 | 5.0 U | UG/L |
| HEPTACHLOR          | 0.50 | 5.0 U | UG/L |
| HEPTACHLOR EPOXIDE  | 0.50 | 5.0 U | UG/L |
| METHOXYCHLOR        | 2.0  | 20 U  | UG/L |
| TOXAPHENE           | 10   | 100 U | UG/L |

SURROGATE RECOVERIESQC LIMITS

|                                |              |     |   |
|--------------------------------|--------------|-----|---|
| DECACHLOROBIPHENYL (DCB)       | (24 - 154 %) | 125 | % |
| TETRACHLORO-META-XYLENE (TCMX) | (30 - 150 %) | 98  | % |

Data Reported following TCLP Toxicity Characteristics Leaching Procedure.  
Federal Register, Part 261, Vol. 55, NO 126, June 29, 1990.

COLUMBIA ANALYTICAL SERVICES

EXTRACTABLE ORGANICS

METHOD 8151A

Reported: 08/09/00

Project Reference:

Client Sample ID : METHOD BLANK

|                |                 |                              |
|----------------|-----------------|------------------------------|
| Date Sampled : | Order #: 398563 | Sample Matrix: SOIL/SEDIMENT |
| Date Received: | Submission #:   | Analytical Run 53627         |

| ANALYTE | PQL | RESULT | UNITS |
|---------|-----|--------|-------|
|---------|-----|--------|-------|

DATE EXTRACTED : 07/28/00  
DATE ANALYZED : 08/04/00  
ANALYTICAL DILUTION: 100.00

|                   |      |      |      |
|-------------------|------|------|------|
| 2,4-D             | 0.50 | 50 U | UG/L |
| 2,4,5-TP (SILVEX) | 0.50 | 50 U | UG/L |

SURROGATE RECOVERIES

QC LIMITS

|      |              |     |   |
|------|--------------|-----|---|
| DCAA | (18 - 152 %) | 110 | % |
|------|--------------|-----|---|

Data Reported following TCLP Toxicity Characteristics Leaching Procedure.  
Federal Register, Part 261, Vol. 55, NO 126, June 29, 1990.

00007







1

2

3





## VOLATILE ANALYTICAL REPORT

Client : Blasland, Bouck & Lee  
Account # : 10624  
Site : NMPC ONEIDA CEDAR ST

Date Received : 19-JUL-00  
Date Sampled : 19-JUL-00  
Date Extracted: NA

Matrix : Leachate  
Method : SW846 1311 8260  
Units : ug/l

| Galson ID: | L62002-1   | L62002-2     | WG24361-1    |
|------------|------------|--------------|--------------|
| Client ID: | SB-2 (0-2) | SB-3 (10-12) | Method Blank |

|                      |          |          |          |
|----------------------|----------|----------|----------|
| Benzene              | <50      | <50      | <50      |
| 2-Butanone           | <100     | <100     | <100     |
| Carbon Tetrachloride | <50      | <50      | <50      |
| Chlorobenzene        | <50      | <50      | <50      |
| Chloroform           | <50      | <50      | <50      |
| 1,2-Dichloroethane   | <50      | <50      | <50      |
| 1,1-Dichloroethene   | <50      | <50      | <50      |
| Tetrachloroethene    | <50      | <50      | <50      |
| Trichloroethene      | <50      | <50      | <50      |
| Vinyl Chloride       | <50      | <50      | <50      |
| Analysis Date        | 07/28/00 | 07/28/00 | 07/28/00 |
| Dilution Factor      | 10       | 10       | 10       |

Approved by : PJT

Date : 09-AUG-00

QC by : *[Signature]*

Date : 8/9/00

NYS DOH # : 11626

Footnotes:

TCLP extraction performed 07/24/00.



## LEACHATE VOLATILE SURROGATE RECOVERY

Client : Blasland, Bouck & Lee

Login # : L62002

[illegible]

SMC1 (TOL) = Toluene-d8  
SMC2 (BFB) = Bromofluorobenzene  
SMC3 (DCE) = 1,2-Dichloroethane-d4

QC LIMITS  
(74-123)  
(72-118)  
(58-134)

```
# Column to be used to flag recovery values
* Values outside of QC limits
D Surrogate diluted out
```



## SEMIVOLATILE ANALYTICAL REPORT

Client : Blasland, Bouck & Lee  
Account # : 10624  
Site : NMPC ONEIDA CEDAR ST

Date Received : 19-JUL-00  
Date Sampled : 19-JUL-00  
Date Extracted: 31-JUL-00

Matrix : Leachate  
Method : SW846 1311 8270-TCLP  
Units : ug/l

| Galson ID: | L62002-1   | L62002-2     | WG24409-1    |
|------------|------------|--------------|--------------|
| Client ID: | SB-2 (0-2) | SB-3 (10-12) | Method Blank |

|                       |      |      |      |
|-----------------------|------|------|------|
| Pyridine              | <100 | <100 | <100 |
| 1,4-Dichlorobenzene   | <100 | <100 | <100 |
| 2-Methylphenol        | <100 | <100 | <100 |
| 3 & 4-Methylphenol    | <200 | <200 | <200 |
| Hexachloroethane      | <100 | <100 | <100 |
| Nitrobenzene          | <100 | <100 | <100 |
| Hexachlorobutadiene   | <100 | <100 | <100 |
| 2,4,6-Trichlorophenol | <100 | <100 | <100 |
| 2,4,5-Trichlorophenol | <100 | <100 | <100 |
| 2,4-Dinitrotoluene    | <100 | <100 | <100 |
| Hexachlorobenzene     | <100 | <100 | <100 |
| ntachlorophenol       | <250 | <250 | <250 |

|                 |          |          |          |
|-----------------|----------|----------|----------|
| Analysis Date   | 08/03/00 | 08/03/00 | 08/03/00 |
| Dilution Factor | 1        | 1        | 1        |

Approved by : Oommen Kappil  
Date : 09-AUG-00  
QC by : *[Signature]*  
Date : 8/10/00  
NYS DOH # : 11626  
Footnotes:

TCLP extraction performed 7/24/00.





## SEMIVOLATILE ANALYTICAL REPORT

Client : Blasland, Bouck & Lee  
Account # : 10624  
Site : NMPC ONEIDA CEDAR ST

Date Received : 19-JUL-00  
Date Sampled : 19-JUL-00  
Date Extracted: 31-JUL-00

Matrix : Leachate  
Method : SW846 1311 8270-TCLP  
Units : ug/l

Galson ID: WG24409-2  
Client ID: TCLP Blank

---

|                       |      |
|-----------------------|------|
| Pyridine              | <100 |
| 1,4-Dichlorobenzene   | <100 |
| 2-Methylphenol        | <100 |
| 3 & 4-Methylphenol    | <200 |
| Hexachloroethane      | <100 |
| Nitrobenzene          | <100 |
| Hexachlorobutadiene   | <100 |
| 2,4,6-Trichlorophenol | <100 |
| 2,4,5-Trichlorophenol | <100 |
| 2,4-Dinitrotoluene    | <100 |
| Hexachlorobenzene     | <100 |
| Pentachlorophenol     | <250 |

Analysis Date 08/03/00  
Dilution Factor 1

---

Approved by : Oommen Kappil  
Date : 09 AUG-00  
QC by : *[Signature]*  
Date : 8/10/00  
NYS DOH # : 11626  
Footnotes:

TCLP extraction performed 7/24/00.



Contract:

SDG No.: L62002

S1 (2FP) = 2-Fluorophenol  
S2 (PHL) = Phenol-d6  
S3 (NBZ) = Nitrobenzene-d5  
S4 (FBP) = 2-Fluorobiphenyl  
S5 (TBP) = 2,4,6-Tribromophenol  
S6 (TPH) = Terphenyl-d14  
S7 (2CP) = 2-Chlorophenol-d4  
S8 (DCB) = 1,2-Dichlorobenzene-d4

(23-120)  
(30-146)  
(42-132)  
(32-127)  
(31-134)  
(43-132)  
(38-135) (advisory)  
(39-132) (advisory)

```
# Column to be used to flag recovery values
* Values outside of QC limits
D Surrogate diluted out
```



## METALS ANALYTICAL REPORT

Client : Blasland, Bouck & Lee  
Account # : 10624  
Site : NMPC ONEIDA CEDAR ST

Date Received : 19-JUL-00  
Date Sampled : 19-JUL-00

Matrix : Leachate  
Method : SW846 6010B/ 7470A; ICP/ CVAA

| Galson ID:    |       | L62002-1   | L62002-2     | WG24230-1    |
|---------------|-------|------------|--------------|--------------|
| Client ID:    |       | SB-2 (0-2) | SB-3 (10-12) | Method Blank |
|               | Units |            |              |              |
| Mercury TCLP  | mg/l  | <0.0003    | <0.0003      | NR           |
| Arsenic TCLP  | mg/l  | <0.01      | 0.016        | <0.01        |
| Barium TCLP   | mg/l  | 1.8        | <1           | <1           |
| Cadmium TCLP  | mg/l  | 0.0087     | 0.0061       | <0.005       |
| Chromium TCLP | mg/l  | <0.01      | <0.01        | <0.01        |
| Lead TCLP     | mg/l  | 0.14       | 0.074        | <0.02        |
| Selenium TCLP | mg/l  | <0.02      | <0.02        | <0.02        |
| Silver TCLP   | mg/l  | <0.01      | <0.01        | <0.01        |

Approved by : LK  
Date : 01-AUG-00  
QC by : *[Signature]*  
Date : 8/9/00  
NYS DOH # : 11626  
Footnotes:







## METALS ANALYTICAL REPORT

Client : Blasland, Bouck & Lee  
Account # : 10624  
Site : NMPC ONEIDA CEDAR ST

Date Received : 19-JUL-00  
Date Sampled : 19-JUL-00

Matrix : Leachate  
Method : SW846 6010B/ 7470A; ICP/ CVAA

Galson ID: WG24362-1  
Client ID: Method Blank

## Units

---

|               |      |         |
|---------------|------|---------|
| Mercury TCLP  | mg/l | <0.0003 |
| Arsenic TCLP  | mg/l | NR      |
| Barium TCLP   | mg/l | NR      |
| Cadmium TCLP  | mg/l | NR      |
| Chromium TCLP | mg/l | NR      |
| Lead TCLP     | mg/l | NR      |
| Selenium TCLP | mg/l | NR      |
| Silver TCLP   | mg/l | NR      |

---

Approved by : LK  
Date : 01-AUG-00  
QC by : *[Signature]*  
Date : 8/1/00  
NYS DOH # : 11626  
Footnotes:





# Galson Laboratories

## INORGANIC ANALYTICAL REPORT

Client : Blasland, Bouck & Lee  
Account # : 10624  
Site : NMPC ONEIDA CEDAR ST

Date Received : 19-JUL-00  
Date Sampled : 19-JUL-00

Matrix : Soil

| Galson ID:       |            |       | L62002-1   | L62002-2     |
|------------------|------------|-------|------------|--------------|
| Client ID:       |            |       | SB-2 (0-2) | SB-3 (10-12) |
|                  | Method     | Units |            |              |
| Corrosivity/pH   | SW846 9045 | SU    | 10.3       | 8.0          |
| Reactive Cyanide | SW846 Ch.  | mg/kg | <100       | <100         |
| Ignitability     | SW846 1030 | NEG   | NEG        | NEG          |
| Reactive Sulfide | SW846 Ch.  | mg/kg | <100       | <100         |

Approved by : DM  
Date : 09-AUG-00  
QC by : *[Signature]*  
Date : 8/9/00  
NYS DOH # : 11626  
Footnotes:

- \* The bulk pH was performed using SW846 method 9045. A sample is corrosive if p is less than or equal to 2, or greater than or equal to 12.5 Standard Units (SU). Under these conditions, the samples are not corrosive.
- \* The samples do not exceed the USEPA action levels of 250mg HCN/kg waste and/or 500 mg H<sub>2</sub>S/kg waste as stated in SW846; therefore they are not reactive.
- \* The samples do not ignite or support combustion.  
Under these conditions the samples are non-ignitable.







1 Mustard ST.  
Suite 250  
Rochester, NY 14609

**THIS IS AN ANALYTICAL TEST REPORT FOR:**

Client : Galson Laboratories  
Project Reference: L62002  
Lab Submission # : R2003091  
Reported : 08/09/00

Report Contains a total of 13 pages

The results reported herein relate only to the samples received by the laboratory. This report may not be reproduced except in full, without the approval of Columbia Analytical Services.

This package has been reviewed by Columbia Analytical Services' QA Department/Laboratory Director to comply with NELAC standards prior to report submittal. *Michael K. Pe...*

00001



#### CASE NARRATIVE

This report contains analytical results for the following samples:

Submission #: R2003091

| <u>Lab ID</u> | <u>Client ID</u> |
|---------------|------------------|
| 396314        | SB-2 (0-2)       |
| 396315        | SB-3 (10-12)     |
| 396316        | TCLP BLANK       |

All samples were received in good condition.

All samples have been analyzed by the approved methods cited on the analytical results pages.

All holding times and associated QC were within limits.

No analytical or QC problems were encountered.



Effective 04/01/96

### CAS LIST OF QUALIFIERS

(The basis of this proposal are the EPA-CLP Qualifiers)

- U - Indicates compound was analyzed for but was not detected. The sample quantitation limit must be corrected for dilution and for percent moisture.
- J - Indicates an estimated value. For further explanation see case narrative / cover letter.
- 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.
- A - This flag indicates that a TIC is a suspected aldol-condensation product.
- N - Spiked sample recovery not within control limits.  
(Flag the entire batch - Inorganic analysis only)
- \* - Duplicate analysis not within control limits.  
(Flag the entire batch - Inorganic analysis only)
- Also used to qualify Organics QC data outside limits.
- D - Spike diluted out.
- S - Reported value determined by Method of Standard Additions. (MSA)
- X - As specified in the case narrative.

### **CAS Lab ID # for State Certifications**

NY ID # in Rochester: 10145  
CT ID # in Rochester: PH0556  
MA ID # in Rochester: M-NY032  
OH EPA # in Rochester: VAP

NJ ID # in Rochester: 73004  
RI ID # in Rochester: 158  
NH ID # in Rochester: 294198-A  
AIHA # in Rochester: 7889

COLUMBIA ANALYTICAL SERVICES**EXTRACTABLE ORGANICS**

METHOD 8081A TCLP

Reported: 08/09/00

Galson Laboratories

Project Reference: L62002

Client Sample ID : SB-2 (0-2)

Date Sampled : 07/19/00 13:00 Order #: 396314      Sample Matrix: SOIL/SEDIMENT  
Date Received: 07/26/00 Submission #: R2003091      Analytical Run 53569

| ANALYTE              | PQL        | RESULT | UNITS |
|----------------------|------------|--------|-------|
| DATE EXTRACTED       | : 07/27/00 |        |       |
| DATE ANALYZED        | : 08/07/00 |        |       |
| ANALYTICAL DILUTION: | 10.00      |        |       |
| GAMMA-BHC (LINDANE)  | 0.50       | 5.0 U  | UG/L  |
| CHLORDANE            | 2.0        | 20 U   | UG/L  |
| ENDRIN               | 0.50       | 5.0 U  | UG/L  |
| HEPTACHLOR           | 0.50       | 5.0 U  | UG/L  |
| HEPTACHLOR EPOXIDE   | 0.50       | 5.0 U  | UG/L  |
| METHOXYCHLOR         | 2.0        | 20 U   | UG/L  |
| TOXAPHENE            | 10         | 100 U  | UG/L  |

SURROGATE RECOVERIESQC LIMITS

|                                |              |     |   |
|--------------------------------|--------------|-----|---|
| DECACHLOROBIPHENYL (DCB)       | (24 - 154 %) | 122 | % |
| TETRACHLORO-META-XYLENE (TCMX) | (30 - 150 %) | 100 | % |

Data Reported following TCLP Toxicity Characteristics Leaching Procedure.  
Federal Register, Part 261, Vol. 55, NO 126, June 29, 1990.

COLUMBIA ANALYTICAL SERVICES

EXTRACTABLE ORGANICS

METHOD 8151A

Reported: 08/09/00

Galson Laboratories

Project Reference: L62002

Client Sample ID : SB-2 (0-2)

Date Sampled : 07/19/00 13:00 Order #: 396314 Sample Matrix: SOIL/SEDIMENT  
Date Received: 07/26/00 Submission #: R2003091 Analytical Run 53627

| ANALYTE | PQL | RESULT | UNITS |
|---------|-----|--------|-------|
|---------|-----|--------|-------|

|                      |   |          |  |
|----------------------|---|----------|--|
| DATE EXTRACTED       | : | 07/28/00 |  |
| DATE ANALYZED        | : | 08/05/00 |  |
| ANALYTICAL DILUTION: |   | 100.00   |  |

|                   |      |      |      |
|-------------------|------|------|------|
| 2,4-D             | 0.50 | 50 U | UG/L |
| 2,4,5-TP (SILVEX) | 0.50 | 50 U | UG/L |

SURROGATE RECOVERIES

QC LIMITS

|      |              |     |   |
|------|--------------|-----|---|
| DCAA | (18 - 152 %) | 109 | % |
|------|--------------|-----|---|

Data Reported following TCLP Toxicity Characteristics Leaching Procedure.  
Federal Register, Part 261, Vol. 55, NO 126, June 29, 1990.

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COLUMBIA ANALYTICAL SERVICES**EXTRACTABLE ORGANICS**

METHOD 8081A TCLP

Reported: 08/09/00

Galson Laboratories

Project Reference: L62002

Client Sample ID : SB-3 (10-12)

Date Sampled : 07/19/00 14:50 Order #: 396315 Sample Matrix: SOIL/SEDIMENT  
Date Received: 07/26/00 Submission #: R2003091 Analytical Run 53569

| ANALYTE                    | PQL  | RESULT | UNITS |
|----------------------------|------|--------|-------|
| DATE EXTRACTED : 07/27/00  |      |        |       |
| DATE ANALYZED : 08/07/00   |      |        |       |
| ANALYTICAL DILUTION: 10.00 |      |        |       |
| GAMMA-BHC (LINDANE)        | 0.50 | 5.0 U  | UG/L  |
| CHLORDANE                  | 2.0  | 20 U   | UG/L  |
| ENDRIN                     | 0.50 | 5.0 U  | UG/L  |
| HEPTACHLOR                 | 0.50 | 5.0 U  | UG/L  |
| HEPTACHLOR EPOXIDE         | 0.50 | 5.0 U  | UG/L  |
| METHOXYCHLOR               | 2.0  | 20 U   | UG/L  |
| TOXAPHENE                  | 10   | 100 U  | UG/L  |

SURROGATE RECOVERIESQC LIMITS

|                                |              |     |   |
|--------------------------------|--------------|-----|---|
| DECACHLOROBIPHENYL (DCB)       | (24 - 154 %) | 129 | % |
| TETRACHLORO-META-XYLENE (TCMX) | (30 - 150 %) | 102 | % |

Data Reported following TCLP Toxicity Characteristics Leaching Procedure.  
Federal Register, Part 261, Vol. 55, NO 126, June 29, 1990.

COLUMBIA ANALYTICAL SERVICES

EXTRACTABLE ORGANICS

METHOD 8151A

Reported: 08/09/00

Galson Laboratories

Project Reference: L62002

Client Sample ID : SB-3 (10-12)

Date Sampled : 07/19/00 14:50 Order #: 396315 Sample Matrix: SOIL/SEDIMENT  
Date Received: 07/26/00 Submission #: R2003091 Analytical Run 53627

| ANALYTE                     | PQL  | RESULT | UNITS |
|-----------------------------|------|--------|-------|
| DATE EXTRACTED : 07/28/00   |      |        |       |
| DATE ANALYZED : 08/05/00    |      |        |       |
| ANALYTICAL DILUTION: 100.00 |      |        |       |
| 2,4-D                       | 0.50 | 50 U   | UG/L  |
| 2,4,5-TP (SILVEX)           | 0.50 | 50 U   | UG/L  |

SURROGATE RECOVERIES

QC LIMITS

|      |              |     |   |
|------|--------------|-----|---|
| DCAA | (18 - 152 %) | 111 | % |
|------|--------------|-----|---|

Data Reported following TCLP Toxicity Characteristics Leaching Procedure.  
Federal Register, Part 261, Vol. 55, NO 126, June 29, 1990.

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COLUMBIA ANALYTICAL SERVICES**EXTRACTABLE ORGANICS**

METHOD 8081A TCLP

Reported: 08/09/00

Galson Laboratories

Project Reference: L62002

Client Sample ID : TCLP BLANK

Date Sampled : 07/24/00      Order #: 396316      Sample Matrix: SOIL/SEDIMENT  
Date Received: 07/26/00      Submission #: R2003091      Analytical Run 53569

| ANALYTE              | PQL        | RESULT | UNITS |
|----------------------|------------|--------|-------|
| DATE EXTRACTED       | : 07/27/00 |        |       |
| DATE ANALYZED        | : 08/07/00 |        |       |
| ANALYTICAL DILUTION: | 10.00      |        |       |
| GAMMA-BHC (LINDANE)  | 0.50       | 5.0 U  | UG/L  |
| CHLORDANE            | 2.0        | 20 U   | UG/L  |
| ENDRIN               | 0.50       | 5.0 U  | UG/L  |
| HEPTACHLOR           | 0.50       | 5.0 U  | UG/L  |
| HEPTACHLOR EPOXIDE   | 0.50       | 5.0 U  | UG/L  |
| METHOXYCHLOR         | 2.0        | 20 U   | UG/L  |
| TOXAPHENE            | 10         | 100 U  | UG/L  |

| SURROGATE RECOVERIES           | QC LIMITS    |     |   |
|--------------------------------|--------------|-----|---|
| DECACHLOROBIPHENYL (DCB)       | (24 - 154 %) | 130 | % |
| TETRACHLORO-META-XYLENE (TCMX) | (30 - 150 %) | 102 | % |

Data Reported following TCLP Toxicity Characteristics Leaching Procedure.  
Federal Register, Part 261, Vol. 55, NO 126, June 29, 1990.

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COLUMBIA ANALYTICAL SERVICES**EXTRACTABLE ORGANICS**

METHOD 8151A

Reported: 08/09/00

Galson Laboratories

Project Reference: L62002

Client Sample ID : TCLP BLANK

Date Sampled : 07/24/00

Order #: 396316

Sample Matrix: SOIL/SEDIMENT

Date Received: 07/26/00

Submission #: R2003091

Analytical Run 53627

| ANALYTE | PQL | RESULT | UNITS |
|---------|-----|--------|-------|
|---------|-----|--------|-------|

DATE EXTRACTED : 07/28/00

DATE ANALYZED : 08/05/00

ANALYTICAL DILUTION: 100.00

|                   |      |      |      |
|-------------------|------|------|------|
| 2,4-D             | 0.50 | 50 U | UG/L |
| 2,4,5-TP (SILVEX) | 0.50 | 50 U | UG/L |

SURROGATE RECOVERIESQC LIMITS

|      |              |     |   |
|------|--------------|-----|---|
| DCAA | (18 - 152 %) | 116 | % |
|------|--------------|-----|---|

Data Reported following TCLP Toxicity Characteristics Leaching Procedure.  
Federal Register, Part 261, Vol. 55, NO 126, June 29, 1990.

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COLUMBIA ANALYTICAL SERVICES**EXTRACTABLE ORGANICS**

METHOD 8081A TCLP

Reported: 08/09/00

**Project Reference:**

Client Sample ID : METHOD BLANK

|                |                 |                              |
|----------------|-----------------|------------------------------|
| Date Sampled : | Order #: 398260 | Sample Matrix: SOIL/SEDIMENT |
| Date Received: | Submission #:   | Analytical Run 53569         |

| ANALYTE | PQL | RESULT | UNITS |
|---------|-----|--------|-------|
|---------|-----|--------|-------|

|                      |            |
|----------------------|------------|
| DATE EXTRACTED       | : 07/27/00 |
| DATE ANALYZED        | : 08/05/00 |
| ANALYTICAL DILUTION: | 10.00      |

|                     |      |       |      |
|---------------------|------|-------|------|
| GAMMA-BHC (LINDANE) | 0.50 | 5.0 U | UG/L |
| CHLORDANE           | 2.0  | 20 U  | UG/L |
| ENDRIN              | 0.50 | 5.0 U | UG/L |
| HEPTACHLOR          | 0.50 | 5.0 U | UG/L |
| HEPTACHLOR EPOXIDE  | 0.50 | 5.0 U | UG/L |
| METHOXYCHLOR        | 2.0  | 20 U  | UG/L |
| TOXAPHENE           | 10   | 100 U | UG/L |

SURROGATE RECOVERIESQC LIMITS

|                                |              |     |   |
|--------------------------------|--------------|-----|---|
| DECACHLOROBIPHENYL (DCB)       | (24 - 154 %) | 125 | % |
| TETRACHLORO-META-XYLENE (TCMX) | (30 - 150 %) | 98  | % |

Data Reported following TCLP Toxicity Characteristics Leaching Procedure.  
Federal Register, Part 261, Vol. 55, NO 126, June 29, 1990.

COLUMBIA ANALYTICAL SERVICES**EXTRACTABLE ORGANICS**

METHOD 8151A

Reported: 08/09/00

Project Reference:

Client Sample ID : METHOD BLANK

|                |                 |                              |
|----------------|-----------------|------------------------------|
| Date Sampled : | Order #: 398563 | Sample Matrix: SOIL/SEDIMENT |
| Date Received: | Submission #:   | Analytical Run 53627         |

| ANALYTE | PQL | RESULT | UNITS |
|---------|-----|--------|-------|
|---------|-----|--------|-------|

|                      |   |          |
|----------------------|---|----------|
| DATE EXTRACTED       | : | 07/28/00 |
| DATE ANALYZED        | : | 08/04/00 |
| ANALYTICAL DILUTION: |   | 100.00   |

|                   |      |      |      |
|-------------------|------|------|------|
| 2,4-D             | 0.50 | 50 U | UG/L |
| 2,4,5-TP (SILVEX) | 0.50 | 50 U | UG/L |

SURROGATE RECOVERIESQC LIMITS

|      |              |     |   |
|------|--------------|-----|---|
| DCAA | (18 - 152 %) | 110 | % |
|------|--------------|-----|---|

Data Reported following TCLP Toxicity Characteristics Leaching Procedure.  
Federal Register, Part 261, Vol. 55, NO 126, June 29, 1990.



Company Name

GALSON LABS

Project Name / Number

L62002.

### um-Around Time

☒ - Standard Service☐ - \* Rush Service

Date requested by: 8/8/00

Ph # ( ) - -

Fax # ( ) -

Page 1 of 1

## PARAMETERS FOR ANALYSIS

Send Report to: Mike B

Send Invoice to: Ham W

P.O. # 10624

[illegible]

REMARKS: TO Columbia

**Total Containers - 3**

\* Results + surrogate recoveries only.

\*TCLP rotation performed by Galsun Laboratories on 7/24/00

\* Hold time = 7/31/00

**SAMPLER'S NAME:**

**SIGNATURE:**

VOC Pres

U

P

ALL

NA

**SAMPLES RELINQUISHED BY:**

**SAMPLES RECEIVED BY:**

### Custody Seal Intact?

☐ Yes ☐ No ☒ N.A.

**Shipment Complete?**

☐ **Yes**    ☐ **No**

Temp 8 °C

TS TB TM

Airbill #

NAME: M. K. ROUSE DATE: 11/25/00  
SIGNATURE: M. K. ROUSE TIME: 4:30 PM

NAME: \_\_\_\_\_ DATE: \_\_\_\_\_  
SIGNATURE: \_\_\_\_\_ TIME: \_\_\_\_\_

|            |       |
|------------|-------|
| NAME:      | DATE: |
| SIGNATURE: | TIME: |

NAME: \_\_\_\_\_ DATE: \_\_\_\_\_  
SIGNATURE: \_\_\_\_\_ TIME: \_\_\_\_\_

Received For Laboratory By: DATE:  
(Signature) TIME:

Received For Laboratory By: *CAS* DATE: 7-26-05  
(Signature) TIME: 10:05

C-1112

**Columbia Analytical Services Inc.**  
**Cooler Receipt And Preservation Check Form**

Project/Client Gilson Submission Number R2-3091  
 Cooler received on 7-26-00 by: JE COURIER: CAS UPS FEDEX CD&L CLIENT

1. Were custody seals on outside of cooler? YES NO
2. Were custody papers properly filled out (ink, signed, etc.)? YES NO
3. Did all bottles arrive in good condition (unbroken)? YES NO
4. Did any VOA vials have significant air bubbles? YES NO N/A
5. Were Ice or Ice packs present? YES NO
6. Where did the bottles originate? CAS/ROC, CLIENT
7. Temperature of cooler(s) upon receipt: 8°

Is the temperature within 0° - 6° C?: Yes ☐ No ☒ Yes ☐ No ☐ Yes ☐ No ☐ Yes ☐ No ☐

If No, Explain Below No ☒ No ☐ No ☐ No ☐ No ☐

Date/Time Temperatures Taken: 7-26-00 @ 10:30

Thermometer ID: IR-Gun Temp Blank Sample Bottle Cooler Temp. IR. Gun

If out of Temperature, Client Approval to Run Samples \_\_\_\_\_

Cooler Breakdown: Date: 7/26/00 by: BC

1. Were all bottle labels complete (i.e. analysis, preservation, etc.)? YES NO
2. Did all bottle labels and tags agree with custody papers? YES NO
3. Were correct containers used for the tests indicated? YES NO
4. Air Samples: Cassettes / Tubes Intact Canisters Pressurized Tedlar® Bags Inflated N/A

Explain any discrepancies: \_\_\_\_\_

|      |                                | YES | NO | Sample I.D. | Reagent | Vol. Added |
|------|--------------------------------|-----|----|-------------|---------|------------|
| pH   | Reagent                        |     |    |             |         |            |
| 12   | NaOH                           |     |    |             |         |            |
| 2    | HNO <sub>3</sub>               |     |    |             |         |            |
| 2    | H <sub>2</sub> SO <sub>4</sub> |     |    |             |         |            |
| 5-9* | P/PCBs<br>(608 only)           |     |    |             |         |            |

YES = All samples OK NO = Samples were preserved at lab as listed PC OK to adjust pH  
 \*If pH adjustment is required, use NaOH and/or H<sub>2</sub>SO<sub>4</sub>

|  |  |  |  |  |
|--|--|--|--|--|
| VOC Vial pH Verification<br>(Tested after Analysis)<br>Following Samples<br>Exhibited pH > 2 |  |  |  |  |
|  |  |  |  |  |
|  |  |  |  |  |
|  |  |  |  |  |
|  |  |  |  |  |

Other Comments:





## VOLATILE ANALYTICAL REPORT

Client : Blasland, Bouck & Lee  
Account # : 10624  
Site : NMPC ONEIDA CEDAR ST

Date Received : 15-AUG-00  
Date Sampled : 15-AUG-00  
Date Extracted: NA

Matrix : Water  
Method : NYSDEC ASP 95-1  
Units : ug/l

| Galson ID: | L62754-1 | L62754-2       | WG24857-1    |
|------------|----------|----------------|--------------|
| Client ID: | PZ-1     | PZ-TRIPBLANK-2 | Method Blank |

|                            |     |     |     |
|----------------------------|-----|-----|-----|
| Chloromethane              | <10 | <10 | <10 |
| Bromomethane               | <10 | <10 | <10 |
| Vinyl Chloride             | <10 | <10 | <10 |
| Chloroethane               | <10 | <10 | <10 |
| Methylene Chloride         | <10 | <10 | <10 |
| Acetone                    | <10 | <10 | <10 |
| Carbon Disulfide           | <10 | <10 | <10 |
| 1,1-Dichloroethene         | <10 | <10 | <10 |
| 1,1-Dichloroethane         | <10 | <10 | <10 |
| 1,2-Dichloroethene (Total) | <10 | <10 | <10 |
| Chloroform                 | <10 | <10 | <10 |
| 1,2-Dichloroethane         | <10 | <10 | <10 |
| Pentanone                  | <10 | <10 | <10 |
| 1,1,1-Trichloroethane      | <10 | <10 | <10 |
| Carbon Tetrachloride       | <10 | <10 | <10 |
| Bromodichloromethane       | <10 | <10 | <10 |
| 1,2-Dichloropropane        | <10 | <10 | <10 |
| cis-1,3-Dichloropropene    | <10 | <10 | <10 |
| Trichloroethene            | <10 | <10 | <10 |
| Dibromochloromethane       | <10 | <10 | <10 |
| 1,1,2-Trichloroethane      | <10 | <10 | <10 |
| Benzene                    | <10 | <10 | <10 |
| trans-1,3-Dichloropropene  | <10 | <10 | <10 |
| Bromoform                  | <10 | <10 | <10 |
| 4-Methyl-2-Pentanone       | <10 | <10 | <10 |
| 2-Hexanone                 | <10 | <10 | <10 |
| Tetrachloroethene          | <10 | <10 | <10 |

Approved by : PJT  
Date : 30-AUG-00  
QC by : *[Signature]*  
Date : 9/6/02  
NYS DOH # : 11626  
Footnotes:





## VOLATILE ANALYTICAL REPORT

Client : Blasland, Bouck & Lee  
Account # : 10624  
Site : NMPC ONEIDA CEDAR ST

Date Received : 15-AUG-00  
Date Sampled : 15-AUG-00  
Date Extracted: NA

Matrix : Water  
Method : NYSDEC ASP 95-1  
Units : ug/l

| Galson ID:                | L62754-1 | L62754-2       | WG24857-1    |
|---------------------------|----------|----------------|--------------|
| Client ID:                | PZ-1     | PZ-TRIPBLANK-2 | Method Blank |
| 1,1,2,2-Tetrachloroethane | <10      | <10            | <10          |
| Toluene                   | <10      | <10            | <10          |
| Chlorobenzene             | <10      | <10            | <10          |
| Ethylbenzene              | <10      | <10            | <10          |
| Styrene                   | <10      | <10            | <10          |
| Xylene (total)            | <10      | <10            | <10          |
| Analysis Date             | 08/16/00 | 08/16/00       | 08/16/00     |
| Dilution Factor           | 1        | 1              | 1            |

Approved by : PJT  
Date : 30-AUG-00  
QC by : *[Signature]*  
Date : 9/6/00  
NYS DOH # : 11626  
Footnotes:



## WATER VOLATILE SURROGATE RECOVERY

Client : Blasland, Bouck & Lee

Login # : L62754

[illegible]

SMC1 (TOL) = Toluene-d8  
SMC2 (BFB) = Bromofluorobenzene  
SMC3 (DCE) = 1,2-Dichloroethane-d4

QC LIMITS  
(62-127)  
(73-120)  
(68-121)

```
# Column to be used to flag recovery values
* Values outside of QC limits
D Surrogate diluted out
```

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

PZ-1

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62754

Matrix: (soil/water) WATER

Lab Sample ID: L62754-1

Sample wt/vol: 1050 (g/mL) mL

Lab File ID: ED82104

Level: (low/med) LOW

Date Received: 08/15/00

% Moisture: decanted: (Y/N) N

Date Extracted: 08/18/00

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: 7.0

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/l

CAS NO.

COMPOUND

Q

|               |                              |    |   |
|---------------|------------------------------|----|---|
| 108-95-2----- | Phenol                       | 10 | U |
| 111-44-4----- | bis(2-Chloroethyl) ether     | 10 | U |
| 95-57-8-----  | 2-Chlorophenol               | 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 |
| 95-48-7-----  | 2-Methylphenol               | 10 | U |
| 108-60-1----- | 2,2'-oxybis(1-Chloropropane) | 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 |
| 120-82-1----- | 1,2,4-Trichlorobenzene       | 10 | U |
| 91-20-3-----  | Naphthalene                  | 10 | U |
| 106-47-8----- | 4-Chloroaniline              | 10 | U |
| 87-68-3-----  | Hexachlorobutadiene          | 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        | 24 | U |
| 91-58-7-----  | 2-Chloronaphthalene          | 10 | U |
| 88-74-4-----  | 2-Nitroaniline               | 24 | U |
| 131-11-3----- | Dimethylphthalate            | 10 | U |
| 208-96-8----- | Acenaphthylene               | 10 | U |
| 606-20-2----- | 2,6-Dinitrotoluene           | 10 | U |
| 99-09-2-----  | 3-Nitroaniline               | 24 | U |
| 83-32-9-----  | Acenaphthene                 | 10 | U |

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

PZ-1

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62754

Matrix: (soil/water) WATER

Lab Sample ID: L62754-1

Sample wt/vol: 1050 (g/mL) mL

Lab File ID: ED82104

Level: (low/med) LOW

Date Received: 08/15/00

% Moisture: decanted: (Y/N) N

Date Extracted: 08/18/00

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: 7.0

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/l

CAS NO.

COMPOUND

Q

|                |                            |    |   |
|----------------|----------------------------|----|---|
| 51-28-5-----   | 2,4-Dinitrophenol          | 24 | U |
| 100-02-7-----  | 4-Nitrophenol              | 24 | 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             | 24 | U |
| 534-52-1-----  | 4,6-Dinitro-2-methylphenol | 24 | U |
| 86-30-6-----   | N-Nitrosodiphenylamine     | 10 | U |
| 101-55-3-----  | 4-Bromophenyl-phenylether  | 10 | U |
| 118-74-1-----  | Hexachlorobenzene          | 10 | U |
| 87-86-5-----   | Pentachlorophenol          | 24 | 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 |
| 56-55-3-----   | Benzo(a)anthracene         | 10 | U |
| 91-94-1-----   | 3,3'-Dichlorobenzidine     | 10 | U |
| 218-01-9-----  | Chrysene                   | 10 | U |
| 117-81-7-----  | bis(2-Ethylhexyl)phthalate | 10 | U |
| 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-----   | Dibenzo(a,h)anthracene     | 10 | U |
| 191-24-2-----  | Benzo(g,h,i)perylene       | 10 | U |

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

SBLK WG24913

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62754

Matrix: (soil/water) WATER

Lab Sample ID: WG24913-1

Sample wt/vol: 1000 (g/mL) mL

Lab File ID: ED82103

Level: (low/med) LOW

Date Received: 00/00/00

% Moisture: decanted: (Y/N) N

Date Extracted: 08/18/00

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: 7.0

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/l

CAS NO.

COMPOUND

Q

|               |                              |    |   |
|---------------|------------------------------|----|---|
| 108-95-2----- | Phenol                       | 10 | U |
| 111-44-4----- | bis(2-Chloroethyl) ether     | 10 | U |
| 95-57-8-----  | 2-Chlorophenol               | 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 |
| 95-48-7-----  | 2-Methylphenol               | 10 | U |
| 108-60-1----- | 2,2'-oxybis(1-Chloropropane) | 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 |
| 120-82-1----- | 1,2,4-Trichlorobenzene       | 10 | U |
| 91-20-3-----  | Naphthalene                  | 10 | U |
| 106-47-8----- | 4-Chloroaniline              | 10 | U |
| 87-68-3-----  | Hexachlorobutadiene          | 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 |
| 91-58-7-----  | 2-Chloronaphthalene          | 10 | U |
| 88-74-4-----  | 2-Nitroaniline               | 25 | U |
| 131-11-3----- | Dimethylphthalate            | 10 | U |
| 208-96-8----- | Acenaphthylene               | 10 | U |
| 606-20-2----- | 2,6-Dinitrotoluene           | 10 | U |
| 99-09-2-----  | 3-Nitroaniline               | 25 | U |
| 83-32-9-----  | Acenaphthene                 | 10 | U |

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SBLK WG24913

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62754

Matrix: (soil/water) WATER

Lab Sample ID: WG24913-1

Sample wt/vol: 1000 (g/mL) mL

Lab File ID: ED82103

Level: (low/med) LOW

Date Received: 00/00/00

% Moisture: decanted: (Y/N) N

Date Extracted: 08/18/00

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: 7.0

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/l

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     | 10 | U |
| 101-55-3-----  | 4-Bromophenyl-phenylether  | 10 | U |
| 118-74-1-----  | Hexachlorobenzene          | 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 |
| 56-55-3-----   | Benzo(a)anthracene         | 10 | U |
| 91-94-1-----   | 3,3'-Dichlorobenzidine     | 10 | U |
| 218-01-9-----  | Chrysene                   | 10 | U |
| 117-81-7-----  | bis(2-Ethylhexyl)phthalate | 10 | U |
| 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-----   | Dibenzo(a,h)anthracene     | 10 | U |
| 191-24-2-----  | Benzo(g,h,i)perylene       | 10 | U |

2C

Contract :

SDG No. : L62754

[illegible]

S1 (2FP) = 2-Fluorophenol  
S2 (PHL) = Phenol-d6  
S3 (NBZ) = Nitrobenzene-d5  
S4 (FBP) = 2-Fluorobiphenyl  
S5 (TBP) = 2,4,6-Tribromophenol  
S6 (TPH) = Terphenyl-d14  
S7 (2CP) = 2-Chlorophenol-d4  
S8 (DCB) = 1,2-Dichlorobenzene-d4

(21-110)

(10-110)  
(35-114)  
(43-116)  
(10-123)  
(33-141)  
(33-110) (advisory)  
(16-110) (advisory)

```
# Column to be used to flag recovery values
* Values outside of QC limits
D Surrogate diluted out
```



1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

PZ-1

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62754

Matrix: (soil/water) WATER

Lab Sample ID: L62754-1

Sample wt/vol: 1000 (g/mL) mL

Lab File ID: HP13A\0817C020

% Moisture: decanted: (Y/N) N

Date Received: 08/15/00

Extraction: (SepF/Cont/Sonc) SEPF

Date Extracted: 08/17/00

Concentrated Extract Volume: 10000 (uL)

Date Analyzed: 08/18/00

Injection Volume: 1.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: 7.0

Sulfur Cleanup: (Y/N) N

|         |          |  |   |
|---------|----------|--|---|
| CAS NO. | COMPOUND | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) ug/l | Q |
|---------|----------|--|---|

|                                  |      |   |
|----------------------------------|------|---|
| 319-84-6-----alpha-BHC           | 0.05 | U |
| 319-85-7-----beta-BHC            | 0.05 | U |
| 319-86-8-----delta-BHC           | 0.05 | U |
| 58-89-9-----gamma-BHC (Lindane)  | 0.05 | U |
| 76-44-8-----Heptachlor           | 0.05 | U |
| 309-00-2-----Aldrin              | 0.05 | U |
| 1024-57-3-----Heptachlor epoxide | 0.05 | U |
| 959-98-8-----Endosulfan I        | 0.05 | U |
| 60-57-1-----Dieldrin             | 0.1  | U |
| 72-55-9-----4,4'-DDE             | 0.1  | U |
| 72-20-8-----Endrin               | 0.1  | U |
| 332-13-659-----Endosulfan II     | 0.1  | U |
| 72-54-8-----4,4'-DDD             | 0.1  | U |
| 103-10-78-----Endosulfan sulfate | 0.1  | U |
| 50-29-3-----4,4'-DDT             | 0.1  | U |
| 72-43-5-----Methoxychlor         | 0.5  | U |
| 53494-70-5-----Endrin ketone     | 0.1  | U |
| 7421-36-3-----Endrin aldehyde    | 0.1  | U |
| 5103-71-9-----alpha-Chlordane    | 0.05 | U |
| 5103-7-42-----gamma-Chlordane    | 0.05 | 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 |

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

PBLK WG24862

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62754

Matrix: (soil/water) WATER

Lab Sample ID: WG24862-1

Sample wt/vol: 1000 (g/mL) mL

Lab File ID: HP13A\0817C015

% Moisture: decanted: (Y/N) N

Date Received: 00/00/00

Extraction: (SepF/Cont/Sonc) SEPF

Date Extracted: 08/17/00

Concentrated Extract Volume: 10000 (uL)

Date Analyzed: 08/17/00

Injection Volume: 1.0 (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

CAS NO.

COMPOUND

Q

|                 |                     |      |   |
|-----------------|---------------------|------|---|
| 319-84-6-----   | alpha-BHC           | 0.05 | U |
| 319-85-7-----   | beta-BHC            | 0.05 | U |
| 319-86-8-----   | delta-BHC           | 0.05 | U |
| 58-89-9-----    | gamma-BHC (Lindane) | 0.05 | U |
| 76-44-8-----    | Heptachlor          | 0.05 | U |
| 309-00-2-----   | Aldrin              | 0.05 | U |
| 1024-57-3-----  | Heptachlor epoxide  | 0.05 | U |
| 959-98-8-----   | Endosulfan I        | 0.05 | U |
| 60-57-1-----    | Dieldrin            | 0.1  | U |
| 72-55-9-----    | 4,4'-DDE            | 0.1  | U |
| 72-20-8-----    | Endrin              | 0.1  | U |
| 332-13-659----- | Endosulfan II       | 0.1  | U |
| 72-54-8-----    | 4,4'-DDD            | 0.1  | U |
| 103-10-78-----  | Endosulfan sulfate  | 0.1  | U |
| 50-29-3-----    | 4,4'-DDT            | 0.1  | U |
| 72-43-5-----    | Methoxychlor        | 0.5  | U |
| 53494-70-5----- | Endrin ketone       | 0.1  | U |
| 7421-36-3-----  | Endrin aldehyde     | 0.1  | U |
| 5103-71-9-----  | alpha-Chlordane     | 0.05 | U |
| 5103-7-42-----  | gamma-Chlordane     | 0.05 | 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 |

Contract:

SDG No.: L62754

ID: .53 (mm)

[illegible]

TCX = TETRACHLORO-M-XYLENE  
DCB = DECACHLOROBIPHENYL

ADVISORY  
QC LIMITS  
(30-150)  
(30-150)

```
# Column to be used to flag recovery values
* Values outside of QC limits
D Surrogate diluted out
```



## METALS ANALYTICAL REPORT

Client : Blasland, Bouck & Lee  
Account # : 10624  
Site : NMPC ONEIDA CEDAR ST

Date Received : 15-AUG-00  
Date Sampled : 15-AUG-00

Matrix : Water  
Method : SW846 6010B/7470A; ICP/CVAA

| Galson ID:          | L62754-1     | L62754-3 | WG24971-1    |
|---------------------|--------------|----------|--------------|
| Client ID:          | PZ-1         | PZ-1     | Method Blank |
|                     | Units        |          |              |
| Mercury             | mg/l <0.0002 | NR       | NR           |
| Mercury, dissolved  | mg/l NR      | <0.00020 | NR           |
| Aluminum            | mg/l 0.21    | NR       | <0.1         |
| Antimony            | mg/l <0.01   | NR       | <0.01        |
| Arsenic             | mg/l <0.01   | NR       | <0.01        |
| Barium              | mg/l 0.036   | NR       | <0.005       |
| Beryllium           | mg/l <0.005  | NR       | <0.005       |
| Cadmium             | mg/l <0.005  | NR       | <0.005       |
| Calcium             | mg/l 450     | NR       | <0.2         |
| Chromium            | mg/l <0.01   | NR       | <0.01        |
| Cobalt              | mg/l <0.01   | NR       | <0.01        |
| Copper              | mg/l <0.01   | NR       | <0.01        |
| Iron                | mg/l 1.4     | NR       | <0.1         |
| Inorganic Lead      | mg/l 0.0093  | NR       | <0.003       |
| Magnesium           | mg/l 21      | NR       | <0.1         |
| Manganese           | mg/l 2.0     | NR       | 0.0054       |
| Nickel              | mg/l <0.02   | NR       | <0.02        |
| Potassium           | mg/l 10      | NR       | <1           |
| Aluminum, dissolved | mg/l NR      | <0.10    | NR           |
| Selenium            | mg/l <0.005  | NR       | <0.005       |
| Antimony, dissolved | mg/l NR      | <0.010   | NR           |
| Silver              | mg/l <0.01   | NR       | <0.01        |
| Arsenic, dissolved  | mg/l NR      | <0.010   | NR           |
| Sodium              | mg/l 5.6     | NR       | <2           |
| Thallium            | mg/l <0.01   | NR       | <0.01        |
| Barium, dissolved   | mg/l NR      | 0.032    | NR           |
| Vanadium            | mg/l <0.01   | NR       | <0.01        |

Approved by : JK  
Date : 06-SEP-00  
QC by : *[Signature]*  
Date : *9/6/00*  
NYS DOH # : 11626  
Footnotes:





## METALS ANALYTICAL REPORT

Client : Blasland, Bouck & Lee  
Account # : 10624  
Site : NMPC ONEIDA CEDAR ST

Date Received : 15-AUG-00  
Date Sampled : 15-AUG-00

Matrix : Water  
Method : SW846 6010B/7470A; ICP/CVAA

| Galson ID: | L62754-1 | L62754-3 | WG24971-1    |
|------------|----------|----------|--------------|
| Client ID: | PZ-1     | PZ-1     | Method Blank |
|            | Units    |          |              |

|                      |      |     |         |       |
|----------------------|------|-----|---------|-------|
| Beryllium, dissolved | mg/l | NR  | <0.0050 | NR    |
| Zinc                 | mg/l | 1.3 | NR      | <0.01 |
| Cadmium, dissolved   | mg/l | NR  | <0.0050 | NR    |
| Calcium, dissolved   | mg/l | NR  | 440     | NR    |
| Chromium, dissolved  | mg/l | NR  | <0.010  | NR    |
| Cobalt, dissolved    | mg/l | NR  | <0.010  | NR    |
| Copper, dissolved    | mg/l | NR  | <0.010  | NR    |
| Iron, dissolved      | mg/l | NR  | 0.79    | NR    |
| Lead, dissolved      | mg/l | NR  | <0.0030 | NR    |
| Magnesium, dissolved | mg/l | NR  | 20      | NR    |
| Manganese, dissolved | mg/l | NR  | 1.8     | NR    |
| Nickel, dissolved    | mg/l | NR  | <0.020  | NR    |
| Potassium, dissolved | mg/l | NR  | 12      | NR    |
| Selenium, dissolved  | mg/l | NR  | 0.019   | NR    |
| Silver, dissolved    | mg/l | NR  | <0.010  | NR    |
| Sodium, dissolved    | mg/l | NR  | 5.3     | NR    |
| Thallium, dissolved  | mg/l | NR  | <0.010  | NR    |
| Vanadium, dissolved  | mg/l | NR  | <0.010  | NR    |
| Zinc, dissolved      | mg/l | NR  | 1.1     | NR    |

Approved by : JK  
Date : 06-SEP-00  
QC by : *[Signature]*  
Date : 9/6/00  
NYS DOH # : 11626  
Footnotes:





## METALS ANALYTICAL REPORT

Client : Blasland, Bouck & Lee  
Account # : 10624  
Site : NMPC ONEIDA CEDAR ST

Date Received : 15-AUG-00  
Date Sampled : 15-AUG-00

Matrix : Water  
Method : SW846 6010B/7470A; ICP/CVAA

Galson ID: WG25084-1  
Client ID: Method Blank

## Units

|                     |      |         |
|---------------------|------|---------|
| Mercury             | mg/l | <0.0002 |
| Mercury, dissolved  | mg/l | NR      |
| Aluminum            | mg/l | NR      |
| Antimony            | mg/l | NR      |
| Arsenic             | mg/l | NR      |
| Barium              | mg/l | NR      |
| Beryllium           | mg/l | NR      |
| Cadmium             | mg/l | NR      |
| Calcium             | mg/l | NR      |
| Chromium            | mg/l | NR      |
| Cobalt              | mg/l | NR      |
| Copper              | mg/l | NR      |
| Iron                | mg/l | NR      |
| Inorganic Lead      | mg/l | NR      |
| Magnesium           | mg/l | NR      |
| Manganese           | mg/l | NR      |
| Nickel              | mg/l | NR      |
| Potassium           | mg/l | NR      |
| Aluminum, dissolved | mg/l | NR      |
| Selenium            | mg/l | NR      |
| Antimony, dissolved | mg/l | NR      |
| Silver              | mg/l | NR      |
| Arsenic, dissolved  | mg/l | NR      |
| Sodium              | mg/l | NR      |
| Thallium            | mg/l | NR      |
| Barium, dissolved   | mg/l | NR      |
| Vanadium            | mg/l | NR      |

Approved by : JK  
Date : 06-SEP-00  
QC by : *[Signature]*  
Date : 9/6/00  
NYS DOH # : 11626  
Footnotes:





## INORGANIC ANALYTICAL REPORT

Client : Blasland, Bouck & Lee  
Account # : 10624  
Site : NMPC ONEIDA CEDAR ST

Date Received : 15-AUG-00  
Date Sampled : 15-AUG-00

Matrix : Water

Galson ID: L62754-1  
Client ID: PZ-1

|                | Method | Units |       |
|----------------|--------|-------|-------|
| Cyanide, Total | CLP-M  | mg/l  | <0.01 |

Approved by : DM  
Date : 24-AUG-00  
QC by : *[Signature]*  
Date : *[Signature]*  
NYS DOH # : 11626  
Footnotes:







BLASLAND, BOUCK & LEE, INC.

LABORATORY DATA REVIEW REPORT

Project: NMPC - Cedar Street  
Analytical Laboratory: Galson Laboratories  
Laboratory Report Identification Number: L62005  
Date of Laboratory Report: August 21, 2000  
Date of Review: November 2, 2000  
Reviewer: Linda Waters  
Number of Samples: 7  
Sample Matrix: soil  
Date of Collection: 7/18/00 - 7/20/00

Sample Analysis: Volatiles

Quality Control Checks

|  |            |           |                       |
|--|------------|-----------|-----------------------|
| 1. Field Chain-of-Custody complete   | <u>yes</u> | no        | not applicable        |
| 2. Proper methods for analysis used  | <u>yes</u> | no        | not applicable        |
| 3. All documentation supplied  | <u>yes</u> | no        | not applicable        |
| 4. Samples analyzed within specified holding times                             | <u>yes</u> | no        | not applicable        |
| 5. The minimum number of field and laboratory QC samples analyzed              | <u>yes</u> | no        | not applicable        |
| 6. Laboratory accuracy maintained within established ranges for the following: |            |           |                       |
| - %RSD, initial calibration  | <u>yes</u> | no        | not applicable        |
| - %D, continuing calibration   | <u>yes</u> | no        | not applicable        |
| - %Recovery, matrix spike  | <u>yes</u> | no        | not applicable        |
| - %Recovery, blank spike   | <u>yes</u> | no        | not applicable        |
| - %Recovery, surrogate   | yes        | <u>no</u> | not applicable        |
| - %Recovery, control sample  | yes        | no        | <u>not applicable</u> |

7. Laboratory precision maintained within established ranges for the following:
- |                     |            |    |                |
|---------------------|------------|----|----------------|
| - RPD, matrix spike | <u>yes</u> | no | not applicable |
| - RPD, duplicates   | <u>yes</u> | no | not applicable |
8. Target analyte concentrations below detection limit in all blank samples
- |     |           |                |
|-----|-----------|----------------|
| yes | <u>no</u> | not applicable |
|-----|-----------|----------------|

Notes: Methylene chloride was detected in three method blanks. Data for methylene chloride have been qualified as undetected in samples SB-2(10-12), SB-2(10-12)RE, DUP-1, DUP-1RE and SB-4A(13-15) based on the blank content.

Recovery for one or more surrogates were outside control limits in samples DUP-1, DUP-1RE and SB-2(10-12)RE. All positive data for samples DUP-1RE and SB-2(10-12)RE and all data for sample DUP-1 have been qualified as estimated based on the recoveries.

The response for one or more internal standards was below established limits in samples DUP-1, DUP-1RE, SB-2(10-12) and SB-2(10-12)RE. Data for all compounds quantitated under the non-compliant standards have been qualified as estimated based on the deviations.

Data for samples DUP-1, SB-2(10-12) and their associated method blank were incorrectly calculated. The samples results and surrogate recoveries were manually recalculated and corrected.

Other than for the deviations noted in this review, all data quality parameters were within method-specified limits and the data is acceptable for use as reported by the laboratory.

Report Number: L62005  
 Sample Analysis: Semivolatiles

## Quality Control Checks

- |   |            |           |                       |
|---|------------|-----------|-----------------------|
| 1. Field Chain-of-Custody complete  | <u>yes</u> | no        | not applicable        |
| 2. Proper methods for analysis used   | <u>yes</u> | no        | not applicable        |
| 3. All documentation supplied   | <u>yes</u> | no        | not applicable        |
| 4. Samples analyzed within specified holding times                              | <u>yes</u> | no        | not applicable        |
| 5. The minimum number of field and laboratory QC samples analyzed               | <u>yes</u> | no        | not applicable        |
| 6. Laboratory accuracy maintained within established ranges for the following:  |            |           |                       |
| - %RSD, initial calibration   | <u>yes</u> | no        | not applicable        |
| - %D, continuing calibration  | <u>yes</u> | no        | not applicable        |
| - %Recovery, matrix spike   | yes        | <u>no</u> | not applicable        |
| - %Recovery, blank spike  | yes        | <u>no</u> | not applicable        |
| - %Recovery, surrogate  | <u>yes</u> | no        | not applicable        |
| - %Recovery, control sample   | yes        | no        | <u>not applicable</u> |
| 7. Laboratory precision maintained within established ranges for the following: |            |           |                       |
| - RPD, matrix spike   | yes        | <u>no</u> | not applicable        |
| - RPD, duplicates   | <u>yes</u> | no        | not applicable        |
| 8. Target analyte concentrations below detection limit in all blank samples     | <u>yes</u> | no        | not applicable        |

Notes: \_\_\_\_\_

A rinse blank was submitted and received for semivolatile analysis. No data for this sample was, however, included in the data package.

Although no target compounds were detected, several non-target compounds were detected in the method blank. When common to the blank and samples, their presence in the samples have been rejected.

The non-target compound benzo(e)pyrene was present but not reported in the

majority of the samples. The non-target sample data sheets have been corrected to reflect the presence.

The response for one or more internal standards was below established limits in samples DUP-1, DUP-1DL, SB-2(19-23), SB-1(4-6) and SB-1(4-6)DL. Data for all compounds quantitated under the non-compliant standards have been qualified as estimated based on the deviations.

Recovery for pentachlorophenol was below control limits in the matrix spike and recovery for 4-nitrophenol was above control limits in the matrix spike blank.

Since the deviations were minor and since neither of the compounds were detected in the samples, no data have been qualified based on the recoveries.

Other than for the deviations noted in this review, the data quality parameters were within method specifications and the data is acceptable for use as reported by the laboratory.

Report Number: L62005  
 Sample Analysis: Pesticides/PCB

## Quality Control Checks

- |   |            |           |                       |
|---|------------|-----------|-----------------------|
| 1. Field Chain-of-Custody complete  | <u>yes</u> | no        | not applicable        |
| 2. Proper methods for analysis used   | <u>yes</u> | no        | not applicable        |
| 3. All documentation supplied   | <u>yes</u> | no        | not applicable        |
| 4. Samples analyzed within specified holding times                              | <u>yes</u> | no        | not applicable        |
| 5. The minimum number of field and laboratory QC samples analyzed               | <u>yes</u> | no        | not applicable        |
| 6. Laboratory accuracy maintained within established ranges for the following:  |            |           |                       |
| - %RSD, initial calibration   | <u>yes</u> | no        | not applicable        |
| - %D, continuing calibration  | <u>yes</u> | no        | not applicable        |
| - %Recovery, matrix spike   | <u>yes</u> | no        | not applicable        |
| - %Recovery, blank spike  | <u>yes</u> | no        | not applicable        |
| - %Recovery, surrogate  | yes        | <u>no</u> | not applicable        |
| - %Recovery, control sample   | yes        | no        | <u>not applicable</u> |
| 7. Laboratory precision maintained within established ranges for the following: |            |           |                       |
| - RPD, matrix spike   | <u>yes</u> | no        | not applicable        |
| - RPD, duplicates   | <u>yes</u> | no        | not applicable        |
| 8. Target analyte concentrations below detection limit in all blank samples     | <u>yes</u> | no        | not applicable        |

Notes: \_\_\_\_\_  
Surrogate recoveries were outside control limits in several samples due to  
interfering peaks. No data have been qualified based on biased recoveries due  
to matrix interference.

\_\_\_\_\_  
Data for several compounds were rejected due to matrix interference. In all  
cases the interference made the determination of presence of absence of the  
rejected compound impossible.

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Other than for the deviations noted in this review, the data quality parameters were within method specifications and the data is acceptable for use as reported by the laboratory.

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Report Number: L62005  
 Sample Analysis: Metals

## Quality Control Checks

- |   |            |           |                |
|---|------------|-----------|----------------|
| 1. Field Chain-of-Custody complete  | <u>yes</u> | no        | not applicable |
| 2. Proper methods for analysis used   | <u>yes</u> | no        | not applicable |
| 3. All documentation supplied   | <u>yes</u> | no        | not applicable |
| 4. Samples analyzed within specified holding times                              | <u>yes</u> | no        | not applicable |
| 5. The minimum number of field and laboratory QC samples analyzed               | <u>yes</u> | no        | not applicable |
| 6. Laboratory accuracy maintained within established ranges for the following:  |            |           |                |
| - r2, initial calibration   | <u>yes</u> | no        | not applicable |
| - %R, continuing calibration  | <u>yes</u> | no        | not applicable |
| - %Recovery, matrix spike   | yes        | <u>no</u> | not applicable |
| - %Recovery, control sample   | <u>yes</u> | no        | not applicable |
| 7. Laboratory precision maintained within established ranges for the following: |            |           |                |
| - RPD, matrix spike   | <u>yes</u> | no        | not applicable |
| - RPD, duplicates   | yes        | <u>no</u> | not applicable |
| 8. Target analyte concentrations below detection limit in all blank samples     | <u>yes</u> | no        | not applicable |

## Notes

Recoveries were below control limits for antimony and manganese and recoveries were above control limits for lead in the soil matrix spike. All data for antimony and manganese and all positive data for lead have been qualified as estimated based on the recoveries.

The field duplicate results are unacceptable for lead. All data for this compound have been qualified as estimated based on the duplicate results.

No target analytes were detected above the CRDL in the preparation and calibration blanks; however, beryllium was detected below the CRDL. Based on

the similarity between the samples and blanks, beryllium data for samples  
DUP1, SB-1(4-6), SB-2(10-12), SB-3(12-14) and SB4(13-15) should be considered  
highly suspect.

Other than for the deviations noted in this review, all quality control parameters  
were within method specifications and the data is acceptable for use as reported  
by the laboratory.

Reviewed and Approved:

  
\_\_\_\_\_  
Quality Assurance Manager

\_\_\_\_\_  
Project Manager



## SAMPLE COMPLIANCE REPORT

## NMPC - Cedar Street

| Sample Delivery Group | Sampling Date | ASP Protocol | Sample ID    | Matrix | Compliance <sup>1</sup> |     |      |     | Noncompliance  |
|-----------------------|---------------|--------------|--------------|--------|-------------------------|-----|------|-----|--|
|                       |               |              |              |        | VOA                     | BNA | Pest | TAL |  |
| L62005                | 7/18/00       | 1995         | SB-1(4-6)    | soil   | yes                     | no  | yes  | yes | BN - ms <sup>2</sup> , msb <sup>2</sup> , int std<br>Pest - id<br>TAL - ms, f.dup                                |
| L58771                | 7/19/00       | 1995         | SB-2(10-12)  | soil   | no                      | no  | yes  | yes | VOA - blank <sup>3</sup><br>BN - ms <sup>2</sup> , msb <sup>2</sup><br>TAL - ms, f.dup                           |
| L58771                | 7/19/00       | 1995         | SB-2(19-23)  | soil   | yes                     | no  | yes  | yes | BN - ms <sup>2</sup> , msb <sup>2</sup><br>Pest - id<br>TAL - ms, f.dup  |
| L58771                | 7/19/00       | 1995         | SB-3(12-14)  | soil   | yes                     | no  | yes  | yes | BN - ms <sup>2</sup> , msb <sup>2</sup><br>TAL - ms, f.dup   |
| L58771                | 7/19/00       | 1995         | DUP-1        | soil   | no                      | no  | yes  | yes | VOA - blank <sup>3</sup> , surr, int std<br>BN - ms <sup>2</sup> , msb <sup>2</sup> , int std<br>TAL - ms, f.dup |
| L58771                | 7/20/00       | 1995         | SB-4A(13-15) | soil   | no                      | no  | yes  | yes | VOA - blank <sup>3</sup><br>BN - ms <sup>2</sup> , msb <sup>2</sup><br>TAL - ms, f.dup                           |
| L58771                | --            | 1995         | RINSE BLANK  | water  | yes                     | --  | yes  | yes | TAL - ms   |
|                       |               |              |              |        |                         |     |      |     |  |

- 1 Samples which are compliant with no added validation qualifiers are listed as "yes". Samples which are non-compliant or which have added qualifiers are listed as "no". A "no" designation does not necessarily indicate that the data have been rejected or are otherwise unusable.
- 2 The noncompliance resulted in no qualification of data.
- 3 Although the deviation resulted in the qualification of data, the laboratory was method compliant.







1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SB-1 (4-6)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) Soil

Lab Sample ID: L62005-1

Sample wt/vol: 4 (g/mL) g

Lab File ID: CD072018

Level: (low/med) MED

Date Received: 07/19/00

%Moisture: not dec. 12

Date Analyzed: 07/20/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 1

Soil Extract Volume: 10000 (uL)

Soil Aliquot Volume: 100 (uL)

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|                 |                            |       |   |
|-----------------|----------------------------|-------|---|
| 74-87-3-----    | Chloromethane              | 1400  | U |
| 74-83-9-----    | Bromomethane               | 1400  | U |
| 75-01-4-----    | Vinyl Chloride             | 1400  | U |
| 75-00-3-----    | Chloroethane               | 1400  | U |
| 75-09-2-----    | Methylene Chloride         | 460   | J |
| 67-64-1-----    | Acetone                    | 1400  | U |
| 75-15-0-----    | Carbon Disulfide           | 1400  | U |
| 75-35-4-----    | 1,1-Dichloroethene         | 1400  | U |
| 75-34-3-----    | 1,1-Dichloroethane         | 1400  | U |
| 540-59-0-----   | 1,2-Dichloroethene (Total) | 1400  | U |
| 67-66-3-----    | Chloroform                 | 1400  | U |
| 107-06-2-----   | 1,2-Dichloroethane         | 1400  | U |
| 78-93-3-----    | 2-Butanone                 | 1400  | U |
| 71-55-6-----    | 1,1,1-Trichloroethane      | 1400  | U |
| 56-23-5-----    | Carbon Tetrachloride       | 1400  | U |
| 75-27-4-----    | Bromodichloromethane       | 1400  | U |
| 78-87-5-----    | 1,2-Dichloropropane        | 1400  | U |
| 10061-01-5----- | cis-1,3-Dichloropropene    | 1400  | U |
| 79-01-6-----    | Trichloroethene            | 1400  | U |
| 124-48-1-----   | Dibromochloromethane       | 1400  | U |
| 79-00-5-----    | 1,1,2-Trichloroethane      | 1400  | U |
| 71-43-2-----    | Benzene                    | 380   | J |
| 10061-02-6----- | trans-1,3-Dichloropropene  | 1400  | U |
| 75-25-2-----    | Bromoform                  | 1400  | U |
| 108-10-1-----   | 4-Methyl-2-Pentanone       | 1400  | U |
| 591-78-6-----   | 2-Hexanone                 | 1400  | U |
| 127-18-4-----   | Tetrachloroethene          | 1400  | U |
| 79-34-5-----    | 1,1,2,2-Tetrachloroethane  | 1400  | U |
| 108-88-3-----   | Toluene                    | 2800  |   |
| 108-90-7-----   | Chlorobenzene              | 1400  | U |
| 100-41-4-----   | Ethylbenzene               | 480   | J |
| 100-42-5-----   | Styrene                    | 950   | J |
| 1330-20-7-----  | Xylene (total)             | 15000 |   |

1E  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SB-1 (4-6)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) Soil

Lab Sample ID: L62005-1

Sample wt/vol: 4 (g/mL) g

Lab File ID: CD072018

Level: (low/med) MED

Date Received: 07/19/00

%Moisture: not dec. 12

Date Analyzed: 07/20/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 1

Soil Extract Volume: 10000 (uL)

Soil Aliquot Volume: 100 (uL)

Number TICS found: 10

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg

| CAS NUMBER     | COMPOUND NAME          | RT    | EST. CONC. | Q |
|----------------|------------------------|-------|------------|---|
| 1. 000091-20-3 | Naphthalene            | 21.56 | 1400       | J |
| 2.             | C3-Substituted benzene | 27.90 | 3500       | J |
| 3.             | C3-Substituted Benzene | 28.11 | 4600       | J |
| 4.             | C3-Substituted Benzene | 28.70 | 1500       | J |
| 5.             | C3-Substituted Benzene | 29.12 | 9000       | J |
| 6. 000271-89-6 | Benzofuran             | 29.93 | 1700       | J |
| 7.             | C3-Substituted Benzene | 30.25 | 2600       | J |
| 8.             | Unknown Aromatic       | 30.81 | 3100       | J |
| 9.             | Unknown Aromatic       | 31.44 | 15000      | J |
| 10.            | C4-Substituted Benzene | 32.95 | 1700       | J |
| 11.            |                        |       |            |   |
| 12.            |                        |       |            |   |
| 13.            |                        |       |            |   |
| 14.            |                        |       |            |   |
| 15.            |                        |       |            |   |
| 16.            |                        |       |            |   |
| 17.            |                        |       |            |   |
| 18.            |                        |       |            |   |
| 19.            |                        |       |            |   |
| 20.            |                        |       |            |   |
| 21.            |                        |       |            |   |
| 22.            |                        |       |            |   |
| 23.            |                        |       |            |   |
| 24.            |                        |       |            |   |
| 25.            |                        |       |            |   |
| 26.            |                        |       |            |   |
| 27.            |                        |       |            |   |
| 28.            |                        |       |            |   |
| 29.            |                        |       |            |   |
| 30.            |                        |       |            |   |

FORM I VOA-TIC

12/91

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SB-2 (10-12)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) Soil

Lab Sample ID: L62005-2

Sample wt/vol: 5 (g/mL) g

Lab File ID: CD072005

Level: (low/med) LOW

Date Received: 07/19/00

%Moisture: not dec. 25

Date Analyzed: 07/20/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 1

Soil Extract Volume: (uL)

Soil Aliquot Volume: (uL)

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|                 |                            |    |      |
|-----------------|----------------------------|----|------|
| 74-87-3-----    | Chloromethane              | 13 | U    |
| 74-83-9-----    | Bromomethane               | 13 | U    |
| 75-01-4-----    | Vinyl Chloride             | 13 | U    |
| 75-00-3-----    | Chloroethane               | 13 | U    |
| 75-09-2-----    | Methylene Chloride         | 13 | B u  |
| 67-64-1-----    | Acetone                    | 13 | U    |
| 75-15-0-----    | Carbon Disulfide           | 13 | U    |
| 75-35-4-----    | 1,1-Dichloroethene         | 13 | U    |
| 75-34-3-----    | 1,1-Dichloroethane         | 13 | U    |
| 540-59-0-----   | 1,2-Dichloroethene (Total) | 13 | U    |
| 67-66-3-----    | Chloroform                 | 13 | U    |
| 107-06-2-----   | 1,2-Dichloroethane         | 13 | U    |
| 78-93-3-----    | 2-Butanone                 | 13 | U    |
| 71-55-6-----    | 1,1,1-Trichloroethane      | 13 | U    |
| 56-23-5-----    | Carbon Tetrachloride       | 13 | U    |
| 75-27-4-----    | Bromodichloromethane       | 13 | U    |
| 78-87-5-----    | 1,2-Dichloropropane        | 13 | U    |
| 10061-01-5----- | cis-1,3-Dichloropropene    | 13 | U    |
| 79-01-6-----    | Trichloroethene            | 13 | U    |
| 124-48-1-----   | Dibromochloromethane       | 13 | U    |
| 79-00-5-----    | 1,1,2-Trichloroethane      | 13 | U    |
| 71-43-2-----    | Benzene                    | 13 | U    |
| 10061-02-6----- | trans-1,3-Dichloropropene  | 13 | U    |
| 75-25-2-----    | Bromoform                  | 13 | U    |
| 108-10-1-----   | 4-Methyl-2-Pentanone       | 13 | U u5 |
| 591-78-6-----   | 2-Hexanone                 | 13 | U u5 |
| 127-18-4-----   | Tetrachloroethene          | 13 | U u5 |
| 79-34-5-----    | 1,1,2,2-Tetrachloroethane  | 13 | U u5 |
| 108-88-3-----   | Toluene                    | 18 | J    |
| 108-90-7-----   | Chlorobenzene              | 13 | U u5 |
| 100-41-4-----   | Ethylbenzene               | 13 | U u5 |
| 100-42-5-----   | Styrene                    | 13 | U u5 |
| 1330-20-7-----  | Xylene (total)             | 13 | U u5 |

1E  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SB-2 (10-12)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) Soil

Lab Sample ID: L62005-2

Sample wt/vol: 5 (g/mL) g

Lab File ID: CD072005

Level: (low/med) LOW

Date Received: 07/19/00

%Moisture: not dec. 25

Date Analyzed: 07/20/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 1

Soil Extract Volume: (uL)

Soil Aliquot Volume: (uL)

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg

Number TICS found: 0

| CAS NUMBER | COMPOUND NAME      | RT    | EST. CONC. | Q     |
|------------|--------------------|-------|------------|-------|
| =====      | =====              | ===== | =====      | ===== |
| 1.         | No Volatiles Found |       |            |       |
| 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

SAMPLE NO.

SB-2 (10-12) RE

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) Soil

Lab Sample ID: L62005-2RE

Sample wt/vol: 5 (g/mL) g

Lab File ID: CD072417

Level: (low/med) LOW

Date Received: 07/19/00

%Moisture: not dec. 25

Date Analyzed: 07/24/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 1

Soil Extract Volume: (uL)

Soil Aliquot Volume: (uL)

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|                 |                            |      |       |
|-----------------|----------------------------|------|-------|
| 74-87-3-----    | Chloromethane              | 13   | U     |
| 74-83-9-----    | Bromomethane               | 13   | U     |
| 75-01-4-----    | Vinyl Chloride             | 13   | U     |
| 75-00-3-----    | Chloroethane               | 13   | U     |
| 75-09-2-----    | Methylene Chloride         | 13 5 | JB u  |
| 67-64-1-----    | Acetone                    | 13   | U     |
| 75-15-0-----    | Carbon Disulfide           | 13   | U     |
| 75-35-4-----    | 1,1-Dichloroethene         | 13   | U     |
| 75-34-3-----    | 1,1-Dichloroethane         | 13   | U     |
| 540-59-0-----   | 1,2-Dichloroethene (Total) | 13   | U     |
| 67-66-3-----    | Chloroform                 | 13   | U     |
| 107-06-2-----   | 1,2-Dichloroethane         | 13   | U     |
| 78-93-3-----    | 2-Butanone                 | 13   | U     |
| 71-55-6-----    | 1,1,1-Trichloroethane      | 13   | U     |
| 56-23-5-----    | Carbon Tetrachloride       | 13   | U     |
| 75-27-4-----    | Bromodichloromethane       | 13   | U     |
| 78-87-5-----    | 1,2-Dichloropropane        | 13   | U     |
| 10061-01-5----- | cis-1,3-Dichloropropene    | 13   | U     |
| 79-01-6-----    | Trichloroethene            | 13   | U     |
| 124-48-1-----   | Dibromochloromethane       | 13   | U     |
| 79-00-5-----    | 1,1,2-Trichloroethane      | 13   | U     |
| 71-43-2-----    | Benzene                    | 2    | J J   |
| 10061-02-6----- | trans-1,3-Dichloropropene  | 13   | U     |
| 75-25-2-----    | Bromoform                  | 13   | U     |
| 108-10-1-----   | 4-Methyl-2-Pentanone       | 13   | u u s |
| 591-78-6-----   | 2-Hexanone                 | 13   | u u s |
| 127-18-4-----   | Tetrachloroethene          | 13   | u u s |
| 79-34-5-----    | 1,1,2,2-Tetrachloroethane  | 13   | u u s |
| 108-88-3-----   | Toluene                    | 6    | J J   |
| 108-90-7-----   | Chlorobenzene              | 13   | u u s |
| 100-41-4-----   | Ethylbenzene               | 13   | u u s |
| 100-42-5-----   | Styrene                    | 13   | u u s |
| 1330-20-7-----  | Xylene (total)             | 13   | u u s |

1E  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SB-2 (10-12) RE

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) Soil

Lab Sample ID: L62005-2 RE

Sample wt/vol: 5 (g/mL) g

Lab File ID: CD072417

Level: (low/med) LOW

Date Received: 07/19/00

%Moisture: not dec. 25

Date Analyzed: 07/24/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 1

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.         | No Volatiles Found |    |            |   |
| 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

SAMPLE NO.

SB-2 (19-23)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) Soil

Lab Sample ID: L62005-3

Sample wt/vol: 4 (g/mL) g

Lab File ID: CD072105

Level: (low/med) MED

Date Received: 07/19/00

%Moisture: not dec. 49

Date Analyzed: 07/21/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 2

Soil Extract Volume: 10000 (uL)

Soil Aliquot Volume: 50 (uL)

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|                 |                            |       |   |
|-----------------|----------------------------|-------|---|
| 74-87-3-----    | Chloromethane              | 4900  | U |
| 74-83-9-----    | Bromomethane               | 4900  | U |
| 75-01-4-----    | Vinyl Chloride             | 4900  | U |
| 75-00-3-----    | Chloroethane               | 4900  | U |
| 75-09-2-----    | Methylene Chloride         | 660   | J |
| 67-64-1-----    | Acetone                    | 4900  | U |
| 75-15-0-----    | Carbon Disulfide           | 4900  | U |
| 75-35-4-----    | 1,1-Dichloroethene         | 4900  | U |
| 75-34-3-----    | 1,1-Dichloroethane         | 4900  | U |
| 540-59-0-----   | 1,2-Dichloroethene (Total) | 4900  | U |
| 67-66-3-----    | Chloroform                 | 4900  | U |
| 107-06-2-----   | 1,2-Dichloroethane         | 4900  | U |
| 78-93-3-----    | 2-Butanone                 | 4900  | U |
| 71-55-6-----    | 1,1,1-Trichloroethane      | 4900  | U |
| 56-23-5-----    | Carbon Tetrachloride       | 4900  | U |
| 75-27-4-----    | Bromodichloromethane       | 4900  | U |
| 78-87-5-----    | 1,2-Dichloropropane        | 4900  | U |
| 10061-01-5----- | cis-1,3-Dichloropropene    | 4900  | U |
| 79-01-6-----    | Trichloroethene            | 4900  | U |
| 124-48-1-----   | Dibromochloromethane       | 4900  | U |
| 79-00-5-----    | 1,1,2-Trichloroethane      | 4900  | U |
| 71-43-2-----    | Benzene                    | 27000 |   |
| 10061-02-6----- | trans-1,3-Dichloropropene  | 4900  | U |
| 75-25-2-----    | Bromoform                  | 4900  | U |
| 108-10-1-----   | 4-Methyl-2-Pentanone       | 4900  | U |
| 591-78-6-----   | 2-Hexanone                 | 4900  | U |
| 127-18-4-----   | Tetrachloroethene          | 4900  | U |
| 79-34-5-----    | 1,1,2,2-Tetrachloroethane  | 4900  | U |
| 108-88-3-----   | Toluene                    | 28000 |   |
| 108-90-7-----   | Chlorobenzene              | 4900  | U |
| 100-41-4-----   | Ethylbenzene               | 18000 |   |
| 100-42-5-----   | Styrene                    | 10000 |   |
| 1330-20-7-----  | Xylene (total)             | 97000 |   |

1E  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SB-2 (19-23)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) Soil

Lab Sample ID: L62005-3

Sample wt/vol: 4 (g/mL) g

Lab File ID: CD072105

Level: (low/med) MED

Date Received: 07/19/00

%Moisture: not dec. 49

Date Analyzed: 07/21/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 2

Soil Extract Volume: 10000 (uL)

Soil Aliquot Volume: 50 (uL)

Number TICS found: 10

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg

| CAS NUMBER      | COMPOUND NAME          | RT    | EST. CONC. | Q |
|-----------------|------------------------|-------|------------|---|
| 1.              | C3-Substituted benzene | 27.92 | 14000      | J |
| 2.              | C3-Substituted benzene | 28.11 | 16000      | J |
| 3.              | C3-Substituted benzene | 29.12 | 36000      | J |
| 4. 000271-89-6  | Benzofuran             | 29.93 | 28000      | J |
| 5.              | Unknown aromatic       | 31.45 | 140000     | J |
| 6.              | Unknown aromatic       | 33.28 | 17000      | J |
| 7.              | Unknown aromatic       | 34.68 | 14000      | J |
| 8.              | Unknown aromatic       | 35.04 | 13000      | J |
| 9. 000091-20-3  | Naphthalene            | 36.61 | 350000     | J |
| 10. 000270-82-6 | Benzo[c]thiophene      | 37.05 | 14000      | J |
| 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

SAMPLE NO.

SB-4A (13-15)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) Soil

Lab Sample ID: L62041-1

Sample wt/vol: 5 (g/mL) g

Lab File ID: CD072414

Level: (low/med) LOW

Date Received: 07/20/00

%Moisture: not dec. 17

Date Analyzed: 07/24/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 1

Soil Extract Volume: (uL)

Soil Aliquot Volume: (uL)

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|                 |                            |    |   |
|-----------------|----------------------------|----|---|
| 74-87-3-----    | Chloromethane              | 12 | U |
| 74-83-9-----    | Bromomethane               | 12 | U |
| 75-01-4-----    | Vinyl Chloride             | 12 | U |
| 75-00-3-----    | Chloroethane               | 12 | U |
| 75-09-2-----    | Methylene Chloride         | 12 | U |
| 67-64-1-----    | Acetone                    | 12 | U |
| 75-15-0-----    | Carbon Disulfide           | 12 | U |
| 75-35-4-----    | 1,1-Dichloroethene         | 12 | U |
| 75-34-3-----    | 1,1-Dichloroethane         | 12 | U |
| 540-59-0-----   | 1,2-Dichloroethene (Total) | 12 | U |
| 67-66-3-----    | Chloroform                 | 12 | U |
| 107-06-2-----   | 1,2-Dichloroethane         | 12 | U |
| 78-93-3-----    | 2-Butanone                 | 12 | U |
| 71-55-6-----    | 1,1,1-Trichloroethane      | 12 | U |
| 56-23-5-----    | Carbon Tetrachloride       | 12 | U |
| 75-27-4-----    | Bromodichloromethane       | 12 | U |
| 78-87-5-----    | 1,2-Dichloropropane        | 12 | U |
| 10061-01-5----- | cis-1,3-Dichloropropene    | 12 | U |
| 79-01-6-----    | Trichloroethene            | 12 | U |
| 124-48-1-----   | Dibromochloromethane       | 12 | U |
| 79-00-5-----    | 1,1,2-Trichloroethane      | 12 | U |
| 71-43-2-----    | Benzene                    | 12 | U |
| 10061-02-6----- | trans-1,3-Dichloropropene  | 12 | U |
| 75-25-2-----    | Bromoform                  | 12 | U |
| 108-10-1-----   | 4-Methyl-2-Pentanone       | 12 | U |
| 591-78-6-----   | 2-Hexanone                 | 12 | U |
| 127-18-4-----   | Tetrachloroethene          | 12 | U |
| 79-34-5-----    | 1,1,2,2-Tetrachloroethane  | 12 | U |
| 108-88-3-----   | Toluene                    | 12 | U |
| 108-90-7-----   | Chlorobenzene              | 12 | U |
| 100-41-4-----   | Ethylbenzene               | 12 | U |
| 100-42-5-----   | Styrene                    | 12 | U |
| 1330-20-7-----  | Xylene (total)             | 12 | U |

1E  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SB-4A (13-15)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) Soil

Lab Sample ID: L62041-1

Sample wt/vol: 5 (g/mL) g

Lab File ID: CD072414

Level: (low/med) LOW

Date Received: 07/20/00

%Moisture: not dec. 17

Date Analyzed: 07/24/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 1

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.         | No Volatiles Found |       |            |       |
| 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

SAMPLE NO.

DUP-1

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) Soil

Lab Sample ID: L62005-5

Sample wt/vol: 5 (g/mL) g

Lab File ID: CD072006

Level: (low/med) LOW

Date Received: 07/19/00

%Moisture: not dec. 26

Date Analyzed: 07/20/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 1

Soil Extract Volume: (uL)

Soil Aliquot Volume: (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) ug/kg Q

CAS NO.

COMPOUND

|                 |                            |       |   |   |   |
|-----------------|----------------------------|-------|---|---|---|
| 74-87-3-----    | Chloromethane              | 14    | U | U | U |
| 74-83-9-----    | Bromomethane               | 14    | U | U | U |
| 75-01-4-----    | Vinyl Chloride             | 14    | U | U | U |
| 75-00-3-----    | Chloroethane               | 14    | U | U | U |
| 75-09-2-----    | Methylene Chloride         | 17 18 | B | U | U |
| 67-64-1-----    | Acetone                    | 14    | U | U | U |
| 75-15-0-----    | Carbon Disulfide           | 14    | U | U | U |
| 75-35-4-----    | 1,1-Dichloroethene         | 14    | U | U | U |
| 75-34-3-----    | 1,1-Dichloroethane         | 14    | U | U | U |
| 540-59-0-----   | 1,2-Dichloroethene (Total) | 14    | U | U | U |
| 67-66-3-----    | Chloroform                 | 14    | U | U | U |
| 107-06-2-----   | 1,2-Dichloroethane         | 14    | U | U | U |
| 78-93-3-----    | 2-Butanone                 | 14    | U | U | U |
| 71-55-6-----    | 1,1,1-Trichloroethane      | 14    | U | U | U |
| 56-23-5-----    | Carbon Tetrachloride       | 14    | U | U | U |
| 75-27-4-----    | Bromodichloromethane       | 14    | U | U | U |
| 78-87-5-----    | 1,2-Dichloropropane        | 14    | U | U | U |
| 10061-01-5----- | cis-1,3-Dichloropropene    | 14    | U | U | U |
| 79-01-6-----    | Trichloroethene            | 14    | U | U | U |
| 124-48-1-----   | Dibromochloromethane       | 14    | U | U | U |
| 79-00-5-----    | 1,1,2-Trichloroethane      | 14    | U | U | U |
| 71-43-2-----    | Benzene                    | 14    | U | U | U |
| 10061-02-6----- | trans-1,3-Dichloropropene  | 14    | U | U | U |
| 75-25-2-----    | Bromoform                  | 14    | U | U | U |
| 108-10-1-----   | 4-Methyl-2-Pentanone       | 14    | U | U | U |
| 591-78-6-----   | 2-Hexanone                 | 14    | U | U | U |
| 127-18-4-----   | Tetrachloroethene          | 14    | U | U | U |
| 79-34-5-----    | 1,1,2,2-Tetrachloroethane  | 14    | U | U | U |
| 108-88-3-----   | Toluene                    | 14    | U | U | U |
| 108-90-7-----   | Chlorobenzene              | 14    | U | U | U |
| 100-41-4-----   | Ethylbenzene               | 14    | U | U | U |
| 100-42-5-----   | Styrene                    | 14    | U | U | U |
| 1330-20-7-----  | Xylene (total)             | 14    | U | U | U |

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1E  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

DUP-1

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) Soil

Lab Sample ID: L62005-5

Sample wt/vol: 5 (g/mL) g

Lab File ID: CD072006

Level: (low/med) LOW

Date Received: 07/19/00

%Moisture: not dec. 26

Date Analyzed: 07/20/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 1

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.         | No Volatiles Found |    |            |   |
| 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

SAMPLE NO.

DUP-1 RE

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) Soil

Lab Sample ID: L62005-5RE

Sample wt/vol: 5 (g/mL) g

Lab File ID: CD072418

Level: (low/med) LOW

Date Received: 07/19/00

%Moisture: not dec. 26

Date Analyzed: 07/24/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 1

Soil Extract Volume: (uL)

Soil Aliquot Volume: (uL)

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|                 |                            |      |      |
|-----------------|----------------------------|------|------|
| 74-87-3-----    | Chloromethane              | 14   | U    |
| 74-83-9-----    | Bromomethane               | 14   | U    |
| 75-01-4-----    | Vinyl Chloride             | 14   | U    |
| 75-00-3-----    | Chloroethane               | 14   | U    |
| 75-09-2-----    | Methylene Chloride         | 14 6 | JB U |
| 67-64-1-----    | Acetone                    | 14   | U    |
| 75-15-0-----    | Carbon Disulfide           | 14   | U    |
| 75-35-4-----    | 1,1-Dichloroethene         | 14   | U    |
| 75-34-3-----    | 1,1-Dichloroethane         | 14   | U    |
| 540-59-0-----   | 1,2-Dichloroethene (Total) | 14   | U    |
| 67-66-3-----    | Chloroform                 | 14   | U    |
| 107-06-2-----   | 1,2-Dichloroethane         | 14   | U    |
| 78-93-3-----    | 2-Butanone                 | 14   | U    |
| 71-55-6-----    | 1,1,1-Trichloroethane      | 14   | U    |
| 56-23-5-----    | Carbon Tetrachloride       | 14   | U    |
| 75-27-4-----    | Bromodichloromethane       | 14   | U    |
| 78-87-5-----    | 1,2-Dichloropropane        | 14   | U    |
| 10061-01-5----- | cis-1,3-Dichloropropene    | 14   | U    |
| 79-01-6-----    | Trichloroethene            | 14   | U    |
| 124-48-1-----   | Dibromochloromethane       | 14   | U    |
| 79-00-5-----    | 1,1,2-Trichloroethane      | 14   | U    |
| 71-43-2-----    | Benzene                    | 2    | J J  |
| 10061-02-6----- | trans-1,3-Dichloropropene  | 14   | U    |
| 75-25-2-----    | Bromoform                  | 14   | U    |
| 108-10-1-----   | 4-Methyl-2-Pentanone       | 14   | U U  |
| 591-78-6-----   | 2-Hexanone                 | 14   | U U  |
| 127-18-4-----   | Tetrachloroethene          | 14   | U U  |
| 79-34-5-----    | 1,1,2,2-Tetrachloroethane  | 14   | U U  |
| 108-88-3-----   | Toluene                    | 3    | J J  |
| 108-90-7-----   | Chlorobenzene              | 14   | U U  |
| 100-41-4-----   | Ethylbenzene               | 14   | U U  |
| 100-42-5-----   | Styrene                    | 14   | U U  |
| 1330-20-7-----  | Xylene (total)             | 14   | U U  |

1E  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

DUP-1 RE

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) Soil

Lab Sample ID: L62005-5 RE

Sample wt/vol: 5 (g/mL) g

Lab File ID: CD072418

Level: (low/med) LOW

Date Received: 07/19/00

%Moisture: not dec. 26

Date Analyzed: 07/24/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 1

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.         | No Volatiles Found |       |            |       |
| 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

SAMPLE NO.

RINSE BLANK

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) Water

Lab Sample ID: L62041-5

Sample wt/vol: 5 (g/mL) mL

Lab File ID: CD072404

Level: (low/med) LOW

Date Received: 07/20/00

%Moisture: not dec.

Date Analyzed: 07/24/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 1

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 |
|------------|----------------------------|--|---|
| 74-87-3    | Chloromethane              | 10   | U |
| 74-83-9    | Bromomethane               | 10   | U |
| 75-01-4    | Vinyl Chloride             | 10   | U |
| 75-00-3    | Chloroethane               | 10   | U |
| 75-09-2    | Methylene Chloride         | 10   | U |
| 67-64-1    | Acetone                    | 10   | U |
| 75-15-0    | Carbon Disulfide           | 10   | U |
| 75-35-4    | 1,1-Dichloroethene         | 10   | U |
| 75-34-3    | 1,1-Dichloroethane         | 10   | U |
| 540-59-0   | 1,2-Dichloroethene (Total) | 10   | U |
| 67-66-3    | Chloroform                 | 10   | U |
| 107-06-2   | 1,2-Dichloroethane         | 10   | U |
| 78-93-3    | 2-Butanone                 | 10   | U |
| 71-55-6    | 1,1,1-Trichloroethane      | 10   | U |
| 56-23-5    | Carbon Tetrachloride       | 10   | U |
| 75-27-4    | Bromodichloromethane       | 10   | U |
| 78-87-5    | 1,2-Dichloropropane        | 10   | U |
| 10061-01-5 | cis-1,3-Dichloropropene    | 10   | U |
| 79-01-6    | Trichloroethene            | 10   | U |
| 124-48-1   | Dibromochloromethane       | 10   | U |
| 79-00-5    | 1,1,2-Trichloroethane      | 10   | U |
| 71-43-2    | Benzene                    | 10   | U |
| 10061-02-6 | trans-1,3-Dichloropropene  | 10   | U |
| 75-25-2    | Bromoform                  | 10   | U |
| 108-10-1   | 4-Methyl-2-Pentanone       | 10   | U |
| 591-78-6   | 2-Hexanone                 | 10   | U |
| 127-18-4   | Tetrachloroethene          | 10   | U |
| 79-34-5    | 1,1,2,2-Tetrachloroethane  | 10   | U |
| 108-88-3   | Toluene                    | 10   | U |
| 108-90-7   | Chlorobenzene              | 10   | U |
| 100-41-4   | Ethylbenzene               | 10   | U |
| 100-42-5   | Styrene                    | 10   | U |
| 1330-20-7  | Xylene (total)             | 10   | U |

1E  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

RINSE BLANK

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) Water

Lab Sample ID: L62041-5

Sample wt/vol: 5 (g/mL) mL

Lab File ID: CD072404

Level: (low/med) LOW

Date Received: 07/20/00

%Moisture: not dec.

Date Analyzed: 07/24/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 1

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.         | No Volatiles Found |       |            |       |
| 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.        |                    |       |            |       |

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

SB-1 (4-6)

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-1

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED73139

Level: (low/med) LOW

Date Received: 07/19/00

% Moisture: 12 decanted: (Y/N) N

Date Extracted: 07/24/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/01/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg Q

|               |                              |                        |                  |
|---------------|------------------------------|------------------------|------------------|
| 108-95-2----- | Phenol                       | 380                    | U                |
| 111-44-4----- | bis(2-Chloroethyl) ether     | 380                    | U                |
| 95-57-8-----  | 2-Chlorophenol               | 380                    | U                |
| 541-73-1----- | 1,3-Dichlorobenzene          | 380                    | U                |
| 106-46-7----- | 1,4-Dichlorobenzene          | 380                    | U                |
| 95-50-1-----  | 1,2-Dichlorobenzene          | 380                    | U                |
| 95-48-7-----  | 2-Methylphenol               | 380                    | U                |
| 108-60-1----- | 2,2'-oxybis(1-Chloropropane) | 380                    | U                |
| 106-44-5----- | 4-Methylphenol               | 380                    | U                |
| 621-64-7----- | N-Nitroso-di-n-propylamine   | 380                    | U                |
| 67-72-1-----  | Hexachloroethane             | 380                    | U                |
| 98-95-3-----  | Nitrobenzene                 | 380                    | U W              |
| 78-59-1-----  | Isophorone                   | 380                    | U W              |
| 88-75-5-----  | 2-Nitrophenol                | 380                    | U W              |
| 105-67-9----- | 2,4-Dimethylphenol           | 380                    | U W              |
| 111-91-1----- | bis(2-Chloroethoxy) methane  | 380                    | U W              |
| 120-83-2----- | 2,4-Dichlorophenol           | 380                    | U W              |
| 120-82-1----- | 1,2,4-Trichlorobenzene       | 380                    | U W              |
| 91-20-3-----  | Naphthalene                  | <del>23000</del> 24000 | <del>E</del> J D |
| 106-47-8----- | 4-Chloroaniline              | 380                    | U W              |
| 87-68-3-----  | Hexachlorobutadiene          | 380                    | U W              |
| 59-50-7-----  | 4-Chloro-3-methylphenol      | 380                    | U W              |
| 91-57-6-----  | 2-Methylnaphthalene          | <del>6200</del> 12000  | <del>E</del> J D |
| 77-47-4-----  | Hexachlorocyclopentadiene    | 380                    | U W              |
| 88-06-2-----  | 2,4,6-Trichlorophenol        | 380                    | U W              |
| 95-95-4-----  | 2,4,5-Trichlorophenol        | 940                    | U W              |
| 91-58-7-----  | 2-Chloronaphthalene          | 380                    | U W              |
| 88-74-4-----  | 2-Nitroaniline               | 940                    | U W              |
| 131-11-3----- | Dimethylphthalate            | 380                    | U W              |
| 208-96-8----- | Acenaphthylene               | <del>18000</del> 24000 | <del>E</del> J D |
| 606-20-2----- | 2,6-Dinitrotoluene           | 380                    | U W              |
| 99-09-2-----  | 3-Nitroaniline               | 940                    | U W              |
| 83-32-9-----  | Acenaphthene                 | <del>4700</del> 5700   | <del>E</del> J D |

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

SB-1 (4-6)

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-1

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED73139

Level: (low/med) LOW

Date Received: 07/19/00

% Moisture: 12 decanted: (Y/N) N

Date Extracted: 07/24/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/01/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|                |                            |        |       |   |     |
|----------------|----------------------------|--------|-------|---|-----|
| 51-28-5-----   | 2,4-Dinitrophenol          | 940    | U     | U | U   |
| 100-02-7-----  | 4-Nitrophenol              | 940    | U     | U | U   |
| 132-64-9-----  | Dibenzofuran               | 11000  | 15000 | E | J D |
| 121-14-2-----  | 2,4-Dinitrotoluene         | 380    | U     | U | U   |
| 84-66-2-----   | Diethylphthalate           | 380    | U     | U | U   |
| 86-73-7-----   | Fluorene                   | 27000  | 20000 | E | J D |
| 7005-72-3----- | 4-Chlorophenyl-phenylether | 380    | U     | U | U   |
| 100-01-6-----  | 4-Nitroaniline             | 940    | U     | U | U   |
| 534-52-1-----  | 4,6-Dinitro-2-methylphenol | 940    | U     | U | U   |
| 86-30-6-----   | N-Nitrosodiphenylamine     | 380    | U     | U | U   |
| 101-55-3-----  | 4-Bromophenyl-phenylether  | 380    | U     | U | U   |
| 118-74-1-----  | Hexachlorobenzene          | 380    | U     | U | U   |
| 87-86-5-----   | Pentachlorophenol          | 940    | U     | U | U   |
| 85-01-8-----   | Phenanthrene               | 120000 | E     | J |     |
| 120-12-7-----  | Anthracene                 | 62000  | E     | J |     |
| 86-74-8-----   | Carbazole                  | 5400   | E     | J |     |
| 84-74-2-----   | Di-n-butylphthalate        | 380    | U     | U | U   |
| 206-44-0-----  | Fluoranthene               | 180000 | E     | J |     |
| 129-00-0-----  | Pyrene                     | 73000  | E     |   |     |
| 85-68-7-----   | Butylbenzylphthalate       | 380    | U     |   |     |
| 56-55-3-----   | Benzo(a)anthracene         | 66000  | E     |   |     |
| 91-94-1-----   | 3,3'-Dichlorobenzidine     | 380    | U     |   |     |
| 218-01-9-----  | Chrysene                   | 20000  | E     |   |     |
| 117-81-7-----  | bis(2-Ethylhexyl)phthalate | 380    | U     |   |     |
| 117-84-0-----  | Di-n-octylphthalate        | 380    | U     |   |     |
| 205-99-2-----  | Benzo(b)fluoranthene       | 44000  | E     |   |     |
| 207-08-9-----  | Benzo(k)fluoranthene       | 9200   | E     |   |     |
| 50-32-8-----   | Benzo(a)pyrene             | 36000  | E     |   |     |
| 193-39-5-----  | Indeno(1,2,3-cd)pyrene     | 24000  | E     |   |     |
| 53-70-3-----   | Dibenzo(a,h)anthracene     | 31000  | E     |   |     |
| 191-24-2-----  | Benzo(g,h,i)perylene       | 26000  | E     |   |     |

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SB-1 (4-6)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-1

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED73139

Level: (low/med) LOW

Date Received: 07/19/00

% Moisture: 12 decanted: (Y/N) N

Date Extracted: 07/24/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/01/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

Number TICS found: 22

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg

| CAS NUMBER     | COMPOUND NAME                | RT    | EST. CONC. | Q   |
|----------------|------------------------------|-------|------------|-----|
| 1.             | Unknown                      | 1.42  | 1700       | JB  |
| 2. 010574-37-5 | 2-Pentene, 2,3-dimethyl-     | 1.91  | 48000      | JBA |
| 3. 000123-42-2 | 2-Pentanone, 4-hydroxy-4-met | 2.49  | 5300       | JA  |
| 4. 000100-42-5 | Styrene                      | 3.42  | 650        | J   |
| 5.             | Unknown                      | 5.13  | 1000       | J   |
| 6. 000095-13-6 | Indene                       | 6.52  | 3000       | J   |
| 7.             | Substituted Naphthalene      | 11.32 | 1100       | J   |
| 8.             | Substituted Naphthalene      | 11.47 | 1700       | J   |
| 9.             | Substituted Naphthalene      | 11.52 | 650        | J   |
| 10.            | Substituted Naphthalene      | 11.68 | 930        | J   |
| 11.            | Substituted Naphthalene      | 12.64 | 1000       | J   |
| 12.            | Substituted Naphthalene      | 12.78 | 770        | J   |
| 13.            | Substituted Naphthalene      | 12.97 | 990        | J   |
| 14.            | Unknown PNA                  | 13.52 | 1300       | J   |
| 15.            | Unknown PNA                  | 15.24 | 400        | J   |
| 16.            | Unknown PNA                  | 17.73 | 5800       | J   |
| 17.            | Unknown PNA                  | 18.15 | 1700       | J   |
| 18.            | Unknown PNA                  | 18.26 | 3600       | J   |
| 19.            | Unknown                      | 19.24 | 710        | J   |
| 20.            | Unknown                      | 19.29 | 430        | J   |
| 21.            | Unknown                      | 19.98 | 760        | J   |
| 22.            | Unknown                      | 20.37 | 8100       | J   |
| 23.            | benzo(e)pyrene               | 22.85 | 8100       | J E |
| 24.            |                              |       |            |     |
| 25.            |                              |       |            |     |
| 26.            |                              |       |            |     |
| 27.            |                              |       |            |     |
| 28.            |                              |       |            |     |
| 29.            |                              |       |            |     |
| 30.            |                              |       |            |     |

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

SB-1 (4-6) DL

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-1DL

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED80414

Level: (low/med) LOW

Date Received: 07/19/00

% Moisture: 12 decanted: (Y/N) N

Date Extracted: 07/24/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/04/00

Injection Volume: 2.0 (uL)

Dilution Factor: 100.0

GPC Cleanup: (Y/N) Y pH: 7.0

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg Q

|               |                              |       |    |
|---------------|------------------------------|-------|----|
| 108-95-2----- | Phenol                       | 38000 | U  |
| 111-44-4----- | bis(2-Chloroethyl) ether     | 38000 | U  |
| 95-57-8-----  | 2-Chlorophenol               | 38000 | U  |
| 541-73-1----- | 1,3-Dichlorobenzene          | 38000 | U  |
| 106-46-7----- | 1,4-Dichlorobenzene          | 38000 | U  |
| 95-50-1-----  | 1,2-Dichlorobenzene          | 38000 | U  |
| 95-48-7-----  | 2-Methylphenol               | 38000 | U  |
| 108-60-1----- | 2,2'-oxybis(1-Chloropropane) | 38000 | U  |
| 106-44-5----- | 4-Methylphenol               | 38000 | U  |
| 621-64-7----- | N-Nitroso-di-n-propylamine   | 38000 | U  |
| 67-72-1-----  | Hexachloroethane             | 38000 | U  |
| 98-95-3-----  | Nitrobenzene                 | 38000 | U  |
| 78-59-1-----  | Isophorone                   | 38000 | U  |
| 88-75-5-----  | 2-Nitrophenol                | 38000 | U  |
| 105-67-9----- | 2,4-Dimethylphenol           | 38000 | U  |
| 111-91-1----- | bis(2-Chloroethoxy) methane  | 38000 | U  |
| 120-83-2----- | 2,4-Dichlorophenol           | 38000 | U  |
| 120-82-1----- | 1,2,4-Trichlorobenzene       | 38000 | U  |
| 91-20-3-----  | Naphthalene                  | 24000 | JD |
| 106-47-8----- | 4-Chloroaniline              | 38000 | U  |
| 87-68-3-----  | Hexachlorobutadiene          | 38000 | U  |
| 59-50-7-----  | 4-Chloro-3-methylphenol      | 38000 | U  |
| 91-57-6-----  | 2-Methylnaphthalene          | 12000 | JD |
| 77-47-4-----  | Hexachlorocyclopentadiene    | 38000 | U  |
| 88-06-2-----  | 2,4,6-Trichlorophenol        | 38000 | U  |
| 95-95-4-----  | 2,4,5-Trichlorophenol        | 94000 | U  |
| 91-58-7-----  | 2-Chloronaphthalene          | 38000 | U  |
| 88-74-4-----  | 2-Nitroaniline               | 94000 | U  |
| 131-11-3----- | Dimethylphthalate            | 38000 | U  |
| 208-96-8----- | Acenaphthylene               | 24000 | JD |
| 606-20-2----- | 2,6-Dinitrotoluene           | 38000 | U  |
| 99-09-2-----  | 3-Nitroaniline               | 94000 | U  |
| 83-32-9-----  | Acenaphthene                 | 5700  | JD |



1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SB-1 (4-6) DL

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-1DL

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED80414

Level: (low/med) LOW

Date Received: 07/19/00

% Moisture: 12 decanted: (Y/N) N

Date Extracted: 07/24/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/04/00

Injection Volume: 2.0 (uL)

Dilution Factor: 100.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|                |                            |        |    |
|----------------|----------------------------|--------|----|
| 51-28-5-----   | 2,4-Dinitrophenol          | 94000  | U  |
| 100-02-7-----  | 4-Nitrophenol              | 94000  | U  |
| 132-64-9-----  | Dibenzofuran               | 15000  | JD |
| 121-14-2-----  | 2,4-Dinitrotoluene         | 38000  | U  |
| 84-66-2-----   | Diethylphthalate           | 38000  | U  |
| 86-73-7-----   | Fluorene                   | 26000  | JD |
| 7005-72-3----- | 4-Chlorophenyl-phenylether | 38000  | U  |
| 100-01-6-----  | 4-Nitroaniline             | 94000  | U  |
| 534-52-1-----  | 4,6-Dinitro-2-methylphenol | 94000  | U  |
| 86-30-6-----   | N-Nitrosodiphenylamine     | 38000  | U  |
| 101-55-3-----  | 4-Bromophenyl-phenylether  | 38000  | U  |
| 118-74-1-----  | Hexachlorobenzene          | 38000  | U  |
| 87-86-5-----   | Pentachlorophenol          | 94000  | U  |
| 85-01-8-----   | Phenanthrene               | 95000  | D  |
| 120-12-7-----  | Anthracene                 | 39000  | D  |
| 86-74-8-----   | Carbazole                  | 5000   | JD |
| 84-74-2-----   | Di-n-butylphthalate        | 38000  | U  |
| 206-44-0-----  | Fluoranthene               | 200000 | D  |
| 129-00-0-----  | Pyrene                     | 210000 | D  |
| 85-68-7-----   | Butylbenzylphthalate       | 38000  | U  |
| 56-55-3-----   | Benzo (a) anthracene       | 120000 | D  |
| 91-94-1-----   | 3,3'-Dichlorobenzidine     | 38000  | U  |
| 218-01-9-----  | Chrysene                   | 80000  | D  |
| 117-81-7-----  | bis(2-Ethylhexyl)phthalate | 38000  | U  |
| 117-84-0-----  | Di-n-octylphthalate        | 38000  | U  |
| 205-99-2-----  | Benzo (b) fluoranthene     | 98000  | D  |
| 207-08-9-----  | Benzo (k) fluoranthene     | 37000  | JD |
| 50-32-8-----   | Benzo (a) pyrene           | 76000  | D  |
| 193-39-5-----  | Indeno (1,2,3-cd) pyrene   | 22000  | JD |
| 53-70-3-----   | Dibenzo (a,h) anthracene   | 6400   | JD |
| 191-24-2-----  | Benzo (g,h,i) perylene     | 22000  | JD |

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SB-1 (4-6) DL

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-1 DL

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED80414

Level: (low/med) LOW

Date Received: 07/19/00

% Moisture: 12 decanted: (Y/N) N

Date Extracted: 07/24/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/04/00

Injection Volume: 2.0 (uL)

Dilution Factor: 100.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:

(ug/L or ug/Kg) ug/kg

Number TICS found: 22

| CAS NUMBER     | COMPOUND NAME             | RT    | EST. CONC. | Q   |
|----------------|---------------------------|-------|------------|-----|
| 1. 000141-79-7 | 3-Penten-2-one, 4-methyl- | 1.45  | 53000      | JBA |
| 2.             | Substituted Naphthalene   | 12.05 | 18000      | J   |
| 3.             | Unknown PAH               | 12.71 | 21000      | J   |
| 4.             | Unknown PAH               | 13.57 | 18000      | J   |
| 5.             | Unknown PAH               | 13.61 | 22000      | J   |
| 6.             | Unknown PAH               | 13.77 | 23000      | J   |
| 7.             | Unknown                   | 14.57 | 31000      | J   |
| 8.             | Unknown PAH               | 14.91 | 30000      | J   |
| 9.             | Unknown PAH               | 14.96 | 36000      | J   |
| 10.            | Unknown PAH               | 15.06 | 77000      | J   |
| 11.            | Unknown PAH               | 15.10 | 24000      | J   |
| 12.            | Unknown PAH               | 15.14 | 30000      | J   |
| 13.            | Substituted Naphthalene   | 15.51 | 22000      | J   |
| 14.            | Unknown PAH               | 15.97 | 63000      | J   |
| 15.            | Unknown PAH               | 16.09 | 47000      | J   |
| 16.            | Unknown PAH               | 16.36 | 42000      | J   |
| 17.            | Unknown PAH               | 16.87 | 18000      | J   |
| 18.            | Unknown PAH               | 17.27 | 26000      | J   |
| 19.            | Unknown PAH               | 18.22 | 21000      | J   |
| 20.            | Unknown PAH               | 18.44 | 20000      | J   |
| 21.            | Unknown PAH               | 20.49 | 18000      | J   |
| 22.            | Unknown PAH               | 20.88 | 36000      | J   |
| 23.            | benzo (k) pyrene          | 21.13 | 34000      | J   |
| 24.            |                           |       |            |     |
| 25.            |                           |       |            |     |
| 26.            |                           |       |            |     |
| 27.            |                           |       |            |     |
| 28.            |                           |       |            |     |
| 29.            |                           |       |            |     |
| 30.            |                           |       |            |     |

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SB-2 (10-12)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-2

Sample wt/vol: 30.3 (g/mL) g

Lab File ID: ED73129

Level: (low/med) LOW

Date Received: 07/19/00

% Moisture: 25 decanted: (Y/N) N

Date Extracted: 07/24/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/01/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|               |                              |      |   |
|---------------|------------------------------|------|---|
| 108-95-2----- | Phenol                       | 440  | U |
| 111-44-4----- | bis(2-Chloroethyl) ether     | 440  | U |
| 95-57-8-----  | 2-Chlorophenol               | 440  | U |
| 541-73-1----- | 1,3-Dichlorobenzene          | 440  | U |
| 106-46-7----- | 1,4-Dichlorobenzene          | 440  | U |
| 95-50-1-----  | 1,2-Dichlorobenzene          | 440  | U |
| 95-48-7-----  | 2-Methylphenol               | 440  | U |
| 108-60-1----- | 2,2'-oxybis(1-Chloropropane) | 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 |
| 120-82-1----- | 1,2,4-Trichlorobenzene       | 440  | U |
| 91-20-3-----  | Naphthalene                  | 680  |   |
| 106-47-8----- | 4-Chloroaniline              | 440  | U |
| 87-68-3-----  | Hexachlorobutadiene          | 440  | U |
| 59-50-7-----  | 4-Chloro-3-methylphenol      | 440  | U |
| 91-57-6-----  | 2-Methylnaphthalene          | 260  | J |
| 77-47-4-----  | Hexachlorocyclopentadiene    | 440  | U |
| 88-06-2-----  | 2,4,6-Trichlorophenol        | 440  | U |
| 95-95-4-----  | 2,4,5-Trichlorophenol        | 1100 | U |
| 91-58-7-----  | 2-Chloronaphthalene          | 440  | U |
| 88-74-4-----  | 2-Nitroaniline               | 1100 | U |
| 131-11-3----- | Dimethylphthalate            | 440  | U |
| 208-96-8----- | Acenaphthylene               | 440  |   |
| 606-20-2----- | 2,6-Dinitrotoluene           | 440  | U |
| 99-09-2-----  | 3-Nitroaniline               | 1100 | U |
| 83-32-9-----  | Acenaphthene                 | 330  | J |

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SB-2 (10-12)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-2

Sample wt/vol: 30.3 (g/mL) g

Lab File ID: ED73129

Level: (low/med) LOW

Date Received: 07/19/00

% Moisture: 25 decanted: (Y/N) N

Date Extracted: 07/24/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/01/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

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               | 610                  |              |   |
| 121-14-2-----  | 2,4-Dinitrotoluene         | 440                  | U            |   |
| 84-66-2-----   | Diethylphthalate           | 440                  | U            |   |
| 86-73-7-----   | Fluorene                   | 490                  |              |   |
| 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     | 440                  | U            |   |
| 101-55-3-----  | 4-Bromophenyl-phenylether  | 440                  | U            |   |
| 118-74-1-----  | Hexachlorobenzene          | 440                  | U            |   |
| 87-86-5-----   | Pentachlorophenol          | 1100                 | U            |   |
| 85-01-8-----   | Phenanthrene               | <del>4600</del> 6000 | <del>U</del> | D |
| 120-12-7-----  | Anthracene                 | 1100                 |              |   |
| 86-74-8-----   | Carbazole                  | 470                  |              |   |
| 84-74-2-----   | Di-n-butylphthalate        | 440                  | U            |   |
| 206-44-0-----  | Fluoranthene               | <del>4100</del> 5400 | <del>U</del> | D |
| 129-00-0-----  | Pyrene                     | 2800                 |              |   |
| 85-68-7-----   | Butylbenzylphthalate       | 440                  | U            |   |
| 56-55-3-----   | Benzo(a)anthracene         | 1600                 |              |   |
| 91-94-1-----   | 3,3'-Dichlorobenzidine     | 440                  | U            |   |
| 218-01-9-----  | Chrysene                   | 1400                 |              |   |
| 117-81-7-----  | bis(2-Ethylhexyl)phthalate | 440                  | U            |   |
| 117-84-0-----  | Di-n-octylphthalate        | 440                  | U            |   |
| 205-99-2-----  | Benzo(b)fluoranthene       | 1600                 |              |   |
| 207-08-9-----  | Benzo(k)fluoranthene       | 250                  | J            |   |
| 50-32-8-----   | Benzo(a)pyrene             | 1000                 |              |   |
| 193-39-5-----  | Indeno(1,2,3-cd)pyrene     | 560                  |              |   |
| 53-70-3-----   | Dibenzo(a,h)anthracene     | 170                  | J            |   |
| 191-24-2-----  | Benzo(g,h,i)perylene       | 680                  |              |   |

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SB-2 (10-12)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-2

Sample wt/vol: 30.3 (g/mL) g

Lab File ID: ED73129

Level: (low/med) LOW

Date Received: 07/19/00

% Moisture: 25 decanted: (Y/N) N

Date Extracted: 07/24/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/01/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

Number TICS found: 8

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg

| CAS NUMBER     | COMPOUND NAME                | RT    | EST. CONC. | Q   |
|----------------|------------------------------|-------|------------|-----|
| 1.             | Unknown                      | 1.08  | 2500       | J   |
| 2.             | Unknown                      | 1.43  | 3300       | JB  |
| 3. 000141-79-7 | 3-Penten-2-one, 4-methyl     | 1.95  | 81000      | JBA |
| 4.             | Unknown                      | 2.25  | 1100       | J   |
| 5. 000123-42-2 | 2-Pentanone, 4-hydroxy-4-met | 2.52  | 8400       | JA  |
| 6.             | Unknown                      | 4.89  | 3500       | J   |
| 7.             | Unknown                      | 5.76  | 1300       | J   |
| 8.             | Unknown                      | 6.81  | 1400       | J   |
| 9.             | benzo(e) pyrene              | 22.52 | 750        | J   |
| 10.            |                              |       |            |     |
| 11.            |                              |       |            |     |
| 12.            |                              |       |            |     |
| 13.            |                              |       |            |     |
| 14.            |                              |       |            |     |
| 15.            |                              |       |            |     |
| 16.            |                              |       |            |     |
| 17.            |                              |       |            |     |
| 18.            |                              |       |            |     |
| 19.            |                              |       |            |     |
| 20.            |                              |       |            |     |
| 21.            |                              |       |            |     |
| 22.            |                              |       |            |     |
| 23.            |                              |       |            |     |
| 24.            |                              |       |            |     |
| 25.            |                              |       |            |     |
| 26.            |                              |       |            |     |
| 27.            |                              |       |            |     |
| 28.            |                              |       |            |     |
| 29.            |                              |       |            |     |
| 30.            |                              |       |            |     |

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SB-2 (10-12) DL

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-2DL

Sample wt/vol: 30.3 (g/mL) g

Lab File ID: ED80416

Level: (low/med) LOW

Date Received: 07/19/00

% Moisture: 25 decanted: (Y/N) N

Date Extracted: 07/24/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/04/00

Injection Volume: 2.0 (uL)

Dilution Factor: 2.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:

(ug/L or ug/Kg) ug/kg Q

CAS NO.

COMPOUND

|               |                              |      |    |
|---------------|------------------------------|------|----|
| 108-95-2----- | Phenol                       | 880  | U  |
| 111-44-4----- | bis(2-Chloroethyl) ether     | 880  | U  |
| 95-57-8-----  | 2-Chlorophenol               | 880  | U  |
| 541-73-1----- | 1,3-Dichlorobenzene          | 880  | U  |
| 106-46-7----- | 1,4-Dichlorobenzene          | 880  | U  |
| 95-50-1-----  | 1,2-Dichlorobenzene          | 880  | U  |
| 95-48-7-----  | 2-Methylphenol               | 880  | U  |
| 108-60-1----- | 2,2'-oxybis(1-Chloropropane) | 880  | U  |
| 106-44-5----- | 4-Methylphenol               | 880  | U  |
| 621-64-7----- | N-Nitroso-di-n-propylamine   | 880  | U  |
| 67-72-1-----  | Hexachloroethane             | 880  | U  |
| 98-95-3-----  | Nitrobenzene                 | 880  | U  |
| 78-59-1-----  | Isophorone                   | 880  | U  |
| 88-75-5-----  | 2-Nitrophenol                | 880  | U  |
| 105-67-9----- | 2,4-Dimethylphenol           | 880  | U  |
| 111-91-1----- | bis(2-Chloroethoxy) methane  | 880  | U  |
| 120-83-2----- | 2,4-Dichlorophenol           | 880  | U  |
| 120-82-1----- | 1,2,4-Trichlorobenzene       | 880  | U  |
| 91-20-3-----  | Naphthalene                  | 610  | JD |
| 106-47-8----- | 4-Chloroaniline              | 880  | U  |
| 87-68-3-----  | Hexachlorobutadiene          | 880  | U  |
| 59-50-7-----  | 4-Chloro-3-methylphenol      | 880  | U  |
| 91-57-6-----  | 2-Methylnaphthalene          | 330  | JD |
| 77-47-4-----  | Hexachlorocyclopentadiene    | 880  | U  |
| 88-06-2-----  | 2,4,6-Trichlorophenol        | 880  | U  |
| 95-95-4-----  | 2,4,5-Trichlorophenol        | 2200 | U  |
| 91-58-7-----  | 2-Chloronaphthalene          | 880  | U  |
| 88-74-4-----  | 2-Nitroaniline               | 2200 | U  |
| 131-11-3----- | Dimethylphthalate            | 880  | U  |
| 208-96-8----- | Acenaphthylene               | 350  | JD |
| 606-20-2----- | 2,6-Dinitrotoluene           | 880  | U  |
| 99-09-2-----  | 3-Nitroaniline               | 2200 | U  |
| 83-32-9-----  | Acenaphthene                 | 250  | JD |

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SB-2 (10-12) DL

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-2DL

Sample wt/vol: 30.3 (g/mL) g

Lab File ID: ED80416

Level: (low/med) LOW

Date Received: 07/19/00

% Moisture: 25 decanted: (Y/N) N

Date Extracted: 07/24/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/04/00

Injection Volume: 2.0 (uL)

Dilution Factor: 2.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:

(ug/L or ug/Kg) ug/kg

Q

CAS NO.

COMPOUND

|                |                            |      |    |
|----------------|----------------------------|------|----|
| 51-28-5-----   | 2,4-Dinitrophenol          | 2200 | U  |
| 100-02-7-----  | 4-Nitrophenol              | 2200 | U  |
| 132-64-9-----  | Dibenzofuran               | 500  | JD |
| 121-14-2-----  | 2,4-Dinitrotoluene         | 880  | U  |
| 84-66-2-----   | Diethylphthalate           | 880  | U  |
| 86-73-7-----   | Fluorene                   | 430  | JD |
| 7005-72-3----- | 4-Chlorophenyl-phenylether | 880  | U  |
| 100-01-6-----  | 4-Nitroaniline             | 2200 | U  |
| 534-52-1-----  | 4,6-Dinitro-2-methylphenol | 2200 | U  |
| 86-30-6-----   | N-Nitrosodiphenylamine     | 880  | U  |
| 101-55-3-----  | 4-Bromophenyl-phenylether  | 880  | U  |
| 118-74-1-----  | Hexachlorobenzene          | 880  | U  |
| 87-86-5-----   | Pentachlorophenol          | 2200 | U  |
| 85-01-8-----   | Phenanthrene               | 6000 | D  |
| 120-12-7-----  | Anthracene                 | 1100 | D  |
| 86-74-8-----   | Carbazole                  | 480  | JD |
| 84-74-2-----   | Di-n-butylphthalate        | 880  | U  |
| 206-44-0-----  | Fluoranthene               | 5400 | D  |
| 129-00-0-----  | Pyrene                     | 6200 | D  |
| 85-68-7-----   | Butylbenzylphthalate       | 880  | U  |
| 56-55-3-----   | Benzo(a)anthracene         | 2700 | D  |
| 91-94-1-----   | 3,3'-Dichlorobenzidine     | 880  | U  |
| 218-01-9-----  | Chrysene                   | 2300 | D  |
| 117-81-7-----  | bis(2-Ethylhexyl)phthalate | 880  | U  |
| 117-84-0-----  | Di-n-octylphthalate        | 880  | U  |
| 205-99-2-----  | Benzo(b)fluoranthene       | 4300 | D  |
| 207-08-9-----  | Benzo(k)fluoranthene       | 1100 | D  |
| 50-32-8-----   | Benzo(a)pyrene             | 2500 | D  |
| 193-39-5-----  | Indeno(1,2,3-cd)pyrene     | 1400 | D  |
| 53-70-3-----   | Dibenzo(a,h)anthracene     | 300  | JD |
| 191-24-2-----  | Benzo(g,h,i)perylene       | 1800 | D  |

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

SB-2 (10-12) DI

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-2 DL

Sample wt/vol: 30.3 (g/mL) g

Lab File ID: ED80416

Level: (low/med) LOW

Date Received: 07/19/00

% Moisture: 25 decanted: (Y/N) N

Date Extracted: 07/24/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/04/00

Injection Volume: 2.0 (uL)

Dilution Factor: 2.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg

Number TICS found: 9

| CAS NUMBER     | COMPOUND NAME                | RT    | EST. CONC. | Q   |
|----------------|------------------------------|-------|------------|-----|
| 1.             | Unknown                      | 0.84  | 7100       | J   |
| 2.             | Unknown                      | 1.14  | 4100       | J   |
| 3. 000141-79-7 | 3-Penten-2-one, 4-methyl-    | 1.54  | 98000      | JBA |
| 4. 000123-42-2 | 2-Pentanone, 4-hydroxy-4-met | 1.98  | 9900       | JBA |
| 5.             | Unknown                      | 4.20  | 3900       | J   |
| 6.             | Unknown                      | 5.14  | 1400       | J   |
| 7.             | Unknown                      | 6.20  | 1600       | J   |
| 8.             | Unknown                      | 13.29 | 1400       | J   |
| 9. 000084-65-1 | 9,10-Anthracenedione         | 15.54 | 1500       | J   |
| 10.            | benzo(e) pyrene              | 21.21 | 1800       | J   |
| 11.            |                              |       |            |     |
| 12.            |                              |       |            |     |
| 13.            |                              |       |            |     |
| 14.            |                              |       |            |     |
| 15.            |                              |       |            |     |
| 16.            |                              |       |            |     |
| 17.            |                              |       |            |     |
| 18.            |                              |       |            |     |
| 19.            |                              |       |            |     |
| 20.            |                              |       |            |     |
| 21.            |                              |       |            |     |
| 22.            |                              |       |            |     |
| 23.            |                              |       |            |     |
| 24.            |                              |       |            |     |
| 25.            |                              |       |            |     |
| 26.            |                              |       |            |     |
| 27.            |                              |       |            |     |
| 28.            |                              |       |            |     |
| 29.            |                              |       |            |     |
| 30.            |                              |       |            |     |



1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SB-2 (19-23)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-3

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED73131

Level: (low/med) LOW

Date Received: 07/19/00

% Moisture: 49 decanted: (Y/N) N

Date Extracted: 07/24/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/01/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|               |                              |        |    |
|---------------|------------------------------|--------|----|
| 108-95-2----- | Phenol                       | 650    | U  |
| 111-44-4----- | bis(2-Chloroethyl) ether     | 650    | U  |
| 95-57-8-----  | 2-Chlorophenol               | 650    | U  |
| 541-73-1----- | 1,3-Dichlorobenzene          | 650    | U  |
| 106-46-7----- | 1,4-Dichlorobenzene          | 650    | U  |
| 95-50-1-----  | 1,2-Dichlorobenzene          | 650    | U  |
| 95-48-7-----  | 2-Methylphenol               | 600    | J  |
| 108-60-1----- | 2,2'-oxybis(1-Chloropropane) | 650    | U  |
| 106-44-5----- | 4-Methylphenol               | 1500   |    |
| 621-64-7----- | N-Nitroso-di-n-propylamine   | 650    | U  |
| 67-72-1-----  | Hexachloroethane             | 650    | U  |
| 98-95-3-----  | Nitrobenzene                 | 650    | U  |
| 78-59-1-----  | Isophorone                   | 650    | U  |
| 88-75-5-----  | 2-Nitrophenol                | 650    | U  |
| 105-67-9----- | 2,4-Dimethylphenol           | 20000  | EJ |
| 111-91-1----- | bis(2-Chloroethoxy) methane  | 650    | U  |
| 120-83-2----- | 2,4-Dichlorophenol           | 650    | U  |
| 120-82-1----- | 1,2,4-Trichlorobenzene       | 650    | U  |
| 91-20-3-----  | Naphthalene                  | 400000 | ED |
| 106-47-8----- | 4-Chloroaniline              | 650    | U  |
| 87-68-3-----  | Hexachlorobutadiene          | 650    | U  |
| 59-50-7-----  | 4-Chloro-3-methylphenol      | 260    | J  |
| 91-57-6-----  | 2-Methylnaphthalene          | 80000  | ED |
| 77-47-4-----  | Hexachlorocyclopentadiene    | 650    | U  |
| 88-06-2-----  | 2,4,6-Trichlorophenol        | 650    | U  |
| 95-95-4-----  | 2,4,5-Trichlorophenol        | 1600   | U  |
| 91-58-7-----  | 2-Chloronaphthalene          | 650    | U  |
| 88-74-4-----  | 2-Nitroaniline               | 1600   | U  |
| 131-11-3----- | Dimethylphthalate            | 650    | U  |
| 208-96-8----- | Acenaphthylene               | 150000 | ED |
| 606-20-2----- | 2,6-Dinitrotoluene           | 650    | U  |
| 99-09-2-----  | 3-Nitroaniline               | 1600   | U  |
| 83-32-9-----  | Acenaphthene                 | 24000  | ED |

FORM I SV-I

451

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

SB-2 (19-23)

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-3

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED73131

Level: (low/med) LOW

Date Received: 07/19/00

% Moisture: 49 decanted: (Y/N) N

Date Extracted: 07/24/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/01/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|                |                            |        |        |   |
|----------------|----------------------------|--------|--------|---|
| 51-28-5-----   | 2,4-Dinitrophenol          | 1600   |        | Y |
| 100-02-7-----  | 4-Nitrophenol              | 1600   |        | Y |
| 132-64-9-----  | Dibenzofuran               | 79000  |        | E |
| 121-14-2-----  | 2,4-Dinitrotoluene         | 650    |        | Y |
| 84-66-2-----   | Diethylphthalate           | 650    |        | Y |
| 86-73-7-----   | Fluorene                   | 110000 |        | E |
| 7005-72-3----- | 4-Chlorophenyl-phenylether | 650    |        | Y |
| 100-01-6-----  | 4-Nitroaniline             | 1600   |        | Y |
| 534-52-1-----  | 4,6-Dinitro-2-methylphenol | 1600   |        | U |
| 86-30-6-----   | N-Nitrosodiphenylamine     | 650    |        | U |
| 101-55-3-----  | 4-Bromophenyl-phenylether  | 650    |        | U |
| 118-74-1-----  | Hexachlorobenzene          | 650    |        | U |
| 87-86-5-----   | Pentachlorophenol          | 1600   |        | U |
| 85-01-8-----   | Phenanthrene               | 210000 | 230000 | E |
| 120-12-7-----  | Anthracene                 | 76000  |        | E |
| 86-74-8-----   | Carbazole                  | 26000  |        | E |
| 84-74-2-----   | Di-n-butylphthalate        | 650    |        | U |
| 206-44-0-----  | Fluoranthene               | 98000  | 120000 | E |
| 129-00-0-----  | Pyrene                     | 84000  | 95000  | E |
| 85-68-7-----   | Butylbenzylphthalate       | 650    |        | Y |
| 56-55-3-----   | Benzo (a) anthracene       | 43000  | 55000  | E |
| 91-94-1-----   | 3,3'-Dichlorobenzidine     | 650    |        | Y |
| 218-01-9-----  | Chrysene                   | 22000  | 40000  | E |
| 117-81-7-----  | bis(2-Ethylhexyl)phthalate | 920    |        | Y |
| 117-84-0-----  | Di-n-octylphthalate        | 650    |        | U |
| 205-99-2-----  | Benzo (b) fluoranthene     | 24000  | 42000  | E |
| 207-08-9-----  | Benzo (k) fluoranthene     | 5000   |        |   |
| 50-32-8-----   | Benzo (a) pyrene           | 19000  | 40000  | E |
| 193-39-5-----  | Indeno (1,2,3-cd) pyrene   | 8700   |        | E |
| 53-70-3-----   | Dibenzo (a,h) anthracene   | 4900   |        |   |
| 191-24-2-----  | Benzo (g,h,i) perylene     | 9200   |        | E |

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SB-2 (19-23)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-3

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED73131

Level: (low/med) LOW

Date Received: 07/19/00

% Moisture: 49 decanted: (Y/N) N

Date Extracted: 07/24/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/01/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg

Number TICS found: 22

| CAS NUMBER      | COMPOUND NAME                | RT    | EST. CONC. | Q |
|-----------------|------------------------------|-------|------------|---|
| 1.              | Unknown                      | 1.43  | 1300       | J |
| 2. 000141-79-7  | 3-Penten-2-one, 4-methyl-    | 1.94  | 34000      | J |
| 3. 000123-42-2  | 2-Pentanone, 4-hydroxy-4-met | 2.51  | 3600       | J |
| 4. 000108-38-3  | Benzene, 1,3-dimethyl-       | 2.96  | 3700       | J |
| 5. 000106-42-3  | p-Xylene                     | 3.44  | 2600       | J |
| 6. 000620-14-4  | Benzene, 1-ethyl-3-methyl-   | 5.03  | 1300       | J |
| 7. 000526-73-8  | Benzene, 1,2,3-trimethyl-    | 5.21  | 1700       | J |
| 8. 000095-36-3  | 1,2,4-Trimethylbenzene       | 6.16  | 1300       | J |
| 9. 000095-13-6  | Indene                       | 6.61  | 20000      | J |
| 10. 017059-52-8 | Benzofuran, 7-methyl-        | 7.39  | 680        | J |
| 11. 017059-52-8 | Benzofuran, 7-methyl-        | 7.52  | 1900       | J |
| 12. 000767-59-9 | 1H-Indene, 1-methyl-         | 8.21  | 1400       | J |
| 13.             | Unknown                      | 9.02  | 390        | J |
| 14.             | Substituted Naphthalene      | 10.36 | 8800       | J |
| 15.             | Substituted Naphthalene      | 11.03 | 760        | J |
| 16.             | Substituted Naphthalene      | 11.34 | 840        | J |
| 17.             | Substituted Naphthalene      | 11.52 | 1400       | J |
| 18.             | Unknown                      | 12.97 | 450        | J |
| 19. 007320-53-8 | Dibenzofuran, 4-methyl-      | 13.53 | 660        | J |
| 20.             | Unknown PAH                  | 18.51 | 540        | J |
| 21. 000243-17-4 | 11H-Benzo[b]fluorene         | 18.65 | 670        | J |
| 22.             | Unknown PAH                  | 22.30 | 450        | J |
| 23.             | Benzo[a]pyrene               | 22.55 | 7700       | J |
| 24.             |                              |       |            |   |
| 25.             |                              |       |            |   |
| 26.             |                              |       |            |   |
| 27.             |                              |       |            |   |
| 28.             |                              |       |            |   |
| 29.             |                              |       |            |   |
| 30.             |                              |       |            |   |

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

SB-2 (19-23) DL

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-3DL

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED80412

Level: (low/med) LOW

Date Received: 07/19/00

% Moisture: 49 decanted: (Y/N) N

Date Extracted: 07/24/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/04/00

Injection Volume: 2.0 (uL)

Dilution Factor: 200.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|               |                              |        |    |
|---------------|------------------------------|--------|----|
| 108-95-2----- | Phenol                       | 130000 | U  |
| 111-44-4----- | bis(2-Chloroethyl) ether     | 130000 | U  |
| 95-57-8-----  | 2-Chlorophenol               | 130000 | U  |
| 541-73-1----- | 1,3-Dichlorobenzene          | 130000 | U  |
| 106-46-7----- | 1,4-Dichlorobenzene          | 130000 | U  |
| 95-50-1-----  | 1,2-Dichlorobenzene          | 130000 | U  |
| 95-48-7-----  | 2-Methylphenol               | 130000 | U  |
| 108-60-1----- | 2,2'-oxybis(1-Chloropropane) | 130000 | U  |
| 106-44-5----- | 4-Methylphenol               | 130000 | U  |
| 621-64-7----- | N-Nitroso-di-n-propylamine   | 130000 | U  |
| 67-72-1-----  | Hexachloroethane             | 130000 | U  |
| 98-95-3-----  | Nitrobenzene                 | 130000 | U  |
| 78-59-1-----  | Isophorone                   | 130000 | U  |
| 88-75-5-----  | 2-Nitrophenol                | 130000 | U  |
| 105-67-9----- | 2,4-Dimethylphenol           | 18000  | JD |
| 111-91-1----- | bis(2-Chloroethoxy) methane  | 130000 | U  |
| 120-83-2----- | 2,4-Dichlorophenol           | 130000 | U  |
| 120-82-1----- | 1,2,4-Trichlorobenzene       | 130000 | U  |
| 91-20-3-----  | Naphthalene                  | 610000 | D  |
| 106-47-8----- | 4-Chloroaniline              | 130000 | U  |
| 87-68-3-----  | Hexachlorobutadiene          | 130000 | U  |
| 59-50-7-----  | 4-Chloro-3-methylphenol      | 130000 | U  |
| 91-57-6-----  | 2-Methylnaphthalene          | 110000 | JD |
| 77-47-4-----  | Hexachlorocyclopentadiene    | 130000 | U  |
| 88-06-2-----  | 2,4,6-Trichlorophenol        | 130000 | U  |
| 95-95-4-----  | 2,4,5-Trichlorophenol        | 320000 | U  |
| 91-58-7-----  | 2-Chloronaphthalene          | 130000 | U  |
| 88-74-4-----  | 2-Nitroaniline               | 320000 | U  |
| 131-11-3----- | Dimethylphthalate            | 130000 | U  |
| 208-96-8----- | Acenaphthylene               | 120000 | JD |
| 606-20-2----- | 2,6-Dinitrotoluene           | 130000 | U  |
| 99-09-2-----  | 3-Nitroaniline               | 320000 | U  |
| 83-32-9-----  | Acenaphthene                 | 19000  | JD |

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SB-2 (19-23) DL

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-3DL

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED80412

Level: (low/med) LOW

Date Received: 07/19/00

% Moisture: 49 decanted: (Y/N) N

Date Extracted: 07/24/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/04/00

Injection Volume: 2.0 (uL)

Dilution Factor: 200.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg Q

| CAS NO.        | COMPOUND                   | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) ug/kg | Q  |
|----------------|----------------------------|---|----|
| 51-28-5-----   | 2,4-Dinitrophenol          | 320000  | U  |
| 100-02-7-----  | 4-Nitrophenol              | 320000  | U  |
| 132-64-9-----  | Dibenzofuran               | 65000   | JD |
| 121-14-2-----  | 2,4-Dinitrotoluene         | 130000  | U  |
| 84-66-2-----   | Diethylphthalate           | 130000  | U  |
| 86-73-7-----   | Fluorene                   | 82000   | JD |
| 7005-72-3----- | 4-Chlorophenyl-phenylether | 130000  | U  |
| 100-01-6-----  | 4-Nitroaniline             | 320000  | U  |
| 534-52-1-----  | 4,6-Dinitro-2-methylphenol | 320000  | U  |
| 86-30-6-----   | N-Nitrosodiphenylamine     | 130000  | U  |
| 101-55-3-----  | 4-Bromophenyl-phenylether  | 130000  | U  |
| 118-74-1-----  | Hexachlorobenzene          | 130000  | U  |
| 87-86-5-----   | Pentachlorophenol          | 320000  | U  |
| 85-01-8-----   | Phenanthrene               | 230000  | D  |
| 120-12-7-----  | Anthracene                 | 81000   | JD |
| 86-74-8-----   | Carbazole                  | 30000   | JD |
| 84-74-2-----   | Di-n-butylphthalate        | 130000  | U  |
| 206-44-0-----  | Fluoranthene               | 120000  | JD |
| 129-00-0-----  | Pyrene                     | 95000   | JD |
| 85-68-7-----   | Butylbenzylphthalate       | 130000  | U  |
| 56-55-3-----   | Benzo(a)anthracene         | 55000   | JD |
| 91-94-1-----   | 3,3'-Dichlorobenzidine     | 130000  | U  |
| 218-01-9-----  | Chrysene                   | 40000   | JD |
| 117-81-7-----  | bis(2-Ethylhexyl)phthalate | 130000  | U  |
| 117-84-0-----  | Di-n-octylphthalate        | 130000  | U  |
| 205-99-2-----  | Benzo(b)fluoranthene       | 42000   | JD |
| 207-08-9-----  | Benzo(k)fluoranthene       | 17000   | JD |
| 50-32-8-----   | Benzo(a)pyrene             | 40000   | JD |
| 193-39-5-----  | Indeno(1,2,3-cd)pyrene     | 130000  | U  |
| 53-70-3-----   | Dibenzo(a,h)anthracene     | 130000  | U  |
| 191-24-2-----  | Benzo(g,h,i)perylene       | 130000  | U  |

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SB-2 (19-23) DL

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-3 DL

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED80412

Level: (low/med) LOW

Date Received: 07/19/00

% Moisture: 49 decanted: (Y/N) N

Date Extracted: 07/24/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/04/00

Injection Volume: 2.0 (uL)

Dilution Factor: 200.0

GPC Cleanup: (Y/N) Y pH: 7.0

Number TICS found: 7

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg

| CAS NUMBER     | COMPOUND NAME            | RT    | EST. CONC. | Q   |
|----------------|--------------------------|-------|------------|-----|
| 1. 000141-79-7 | 3-Penten-2-one, 4-methyl | 1.45  | 120000     | JBA |
| 2. 000271-89-6 | Benzofuran               | 4.99  | 30000      | J   |
| 3. 000673-32-5 | Benzene, 1-propynyl-     | 5.85  | 63000      | J   |
| 4.             | Substituted Naphthalene  | 9.45  | 79000      | J   |
| 5.             | Substituted Naphthalene  | 10.48 | 28000      | J   |
| 6.             | Substituted Naphthalene  | 10.62 | 31000      | J   |
| 7.             | Unknown PNA              | 15.04 | 48000      | J   |
| 8.             | benzo (e) pyrene         | 21.10 | 17000      | J   |
| 9.             |                          |       |            |     |
| 10.            |                          |       |            |     |
| 11.            |                          |       |            |     |
| 12.            |                          |       |            |     |
| 13.            |                          |       |            |     |
| 14.            |                          |       |            |     |
| 15.            |                          |       |            |     |
| 16.            |                          |       |            |     |
| 17.            |                          |       |            |     |
| 18.            |                          |       |            |     |
| 19.            |                          |       |            |     |
| 20.            |                          |       |            |     |
| 21.            |                          |       |            |     |
| 22.            |                          |       |            |     |
| 23.            |                          |       |            |     |
| 24.            |                          |       |            |     |
| 25.            |                          |       |            |     |
| 26.            |                          |       |            |     |
| 27.            |                          |       |            |     |
| 28.            |                          |       |            |     |
| 29.            |                          |       |            |     |
| 30.            |                          |       |            |     |

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SB-3 (12-14)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-4

Sample wt/vol: 30 (g/mL) g

Lab File ID: ED73128

Level: (low/med) LOW

Date Received: 07/19/00

% Moisture: 17 decanted: (Y/N) N

Date Extracted: 07/24/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 07/31/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg Q

|               |                              |      |   |
|---------------|------------------------------|------|---|
| 108-95-2----- | Phenol                       | 400  | U |
| 111-44-4----- | bis(2-Chloroethyl) ether     | 400  | U |
| 95-57-8-----  | 2-Chlorophenol               | 400  | U |
| 541-73-1----- | 1,3-Dichlorobenzene          | 400  | U |
| 106-46-7----- | 1,4-Dichlorobenzene          | 400  | U |
| 95-50-1-----  | 1,2-Dichlorobenzene          | 400  | U |
| 95-48-7-----  | 2-Methylphenol               | 400  | U |
| 108-60-1----- | 2,2'-oxybis(1-Chloropropane) | 400  | U |
| 106-44-5----- | 4-Methylphenol               | 400  | U |
| 621-64-7----- | N-Nitroso-di-n-propylamine   | 400  | U |
| 67-72-1-----  | Hexachloroethane             | 400  | U |
| 98-95-3-----  | Nitrobenzene                 | 400  | U |
| 78-59-1-----  | Isophorone                   | 400  | U |
| 88-75-5-----  | 2-Nitrophenol                | 400  | U |
| 105-67-9----- | 2,4-Dimethylphenol           | 400  | U |
| 111-91-1----- | bis(2-Chloroethoxy) methane  | 400  | U |
| 120-83-2----- | 2,4-Dichlorophenol           | 400  | U |
| 120-82-1----- | 1,2,4-Trichlorobenzene       | 400  | U |
| 91-20-3-----  | Naphthalene                  | 400  | U |
| 106-47-8----- | 4-Chloroaniline              | 400  | U |
| 87-68-3-----  | Hexachlorobutadiene          | 400  | U |
| 59-50-7-----  | 4-Chloro-3-methylphenol      | 400  | U |
| 91-57-6-----  | 2-Methylnaphthalene          | 400  | U |
| 77-47-4-----  | Hexachlorocyclopentadiene    | 400  | U |
| 88-06-2-----  | 2,4,6-Trichlorophenol        | 400  | U |
| 95-95-4-----  | 2,4,5-Trichlorophenol        | 1000 | U |
| 91-58-7-----  | 2-Chloronaphthalene          | 400  | U |
| 88-74-4-----  | 2-Nitroaniline               | 1000 | U |
| 131-11-3----- | Dimethylphthalate            | 400  | U |
| 208-96-8----- | Acenaphthylene               | 400  | U |
| 606-20-2----- | 2,6-Dinitrotoluene           | 400  | U |
| 99-09-2-----  | 3-Nitroaniline               | 1000 | U |
| 83-32-9-----  | Acenaphthene                 | 400  | U |

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

SB-3 (12-14)

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-4

Sample wt/vol: 30 (g/mL) g

Lab File ID: ED73128

Level: (low/med) LOW

Date Received: 07/19/00

% Moisture: 17 decanted: (Y/N) N

Date Extracted: 07/24/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 07/31/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg Q

|                |                            |      |   |
|----------------|----------------------------|------|---|
| 51-28-5-----   | 2,4-Dinitrophenol          | 1000 | U |
| 100-02-7-----  | 4-Nitrophenol              | 1000 | U |
| 132-64-9-----  | Dibenzofuran               | 400  | U |
| 121-14-2-----  | 2,4-Dinitrotoluene         | 400  | U |
| 84-66-2-----   | Diethylphthalate           | 400  | U |
| 86-73-7-----   | Fluorene                   | 400  | U |
| 7005-72-3----- | 4-Chlorophenyl-phenylether | 400  | U |
| 100-01-6-----  | 4-Nitroaniline             | 1000 | U |
| 534-52-1-----  | 4,6-Dinitro-2-methylphenol | 1000 | U |
| 86-30-6-----   | N-Nitrosodiphenylamine     | 400  | U |
| 101-55-3-----  | 4-Bromophenyl-phenylether  | 400  | U |
| 118-74-1-----  | Hexachlorobenzene          | 400  | U |
| 87-86-5-----   | Pentachlorophenol          | 1000 | U |
| 85-01-8-----   | Phenanthrene               | 62   | J |
| 120-12-7-----  | Anthracene                 | 400  | U |
| 86-74-8-----   | Carbazole                  | 400  | U |
| 84-74-2-----   | Di-n-butylphthalate        | 400  | U |
| 206-44-0-----  | Fluoranthene               | 290  | J |
| 129-00-0-----  | Pyrene                     | 400  |   |
| 85-68-7-----   | Butylbenzylphthalate       | 400  | U |
| 56-55-3-----   | Benzo(a)anthracene         | 250  | J |
| 91-94-1-----   | 3,3'-Dichlorobenzidine     | 400  | U |
| 218-01-9-----  | Chrysene                   | 200  | J |
| 117-81-7-----  | bis(2-Ethylhexyl)phthalate | 400  | U |
| 117-84-0-----  | Di-n-octylphthalate        | 400  | U |
| 205-99-2-----  | Benzo(b)fluoranthene       | 320  | J |
| 207-08-9-----  | Benzo(k)fluoranthene       | 140  | J |
| 50-32-8-----   | Benzo(a)pyrene             | 280  | J |
| 193-39-5-----  | Indeno(1,2,3-cd)pyrene     | 130  | J |
| 53-70-3-----   | Dibenzo(a,h)anthracene     | 400  | U |
| 191-24-2-----  | Benzo(g,h,i)perylene       | 140  | J |



1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SB-3 (12-14)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-4

Sample wt/vol: 30 (g/mL) g

Lab File ID: ED73128

Level: (low/med) LOW

Date Received: 07/19/00

% Moisture: 17 decanted: (Y/N) N

Date Extracted: 07/24/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 07/31/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

Number TICS found: 9

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg

| CAS NUMBER     | COMPOUND NAME                | RT    | EST. CONC. | Q    |
|----------------|------------------------------|-------|------------|------|
| 1.             | Unknown                      | 1.08  | 720        | J    |
| 2.             | Unknown                      | 1.42  | 2800       | JB   |
| 3. 010574-37-5 | 2-Pentene, 2,3-dimethyl      | 1.94  | 71000      | JBA- |
| 4.             | Unknown                      | 2.25  | 1800       | J    |
| 5. 000123-42-2 | 2-Pentanone, 4-hydroxy-4-met | 2.51  | 7100       | JA-  |
| 6.             | Unknown                      | 4.85  | 1600       | J    |
| 7.             | Unknown                      | 5.76  | 1600       | J    |
| 8.             | Unknown                      | 6.81  | 1500       | J    |
| 9.             | Unknown Alkane               | 14.11 | 840        | J    |
| 10.            | benzo(e) pyrene              | 22.38 | 160        | J    |
| 11.            |                              |       |            |      |
| 12.            |                              |       |            |      |
| 13.            |                              |       |            |      |
| 14.            |                              |       |            |      |
| 15.            |                              |       |            |      |
| 16.            |                              |       |            |      |
| 17.            |                              |       |            |      |
| 18.            |                              |       |            |      |
| 19.            |                              |       |            |      |
| 20.            |                              |       |            |      |
| 21.            |                              |       |            |      |
| 22.            |                              |       |            |      |
| 23.            |                              |       |            |      |
| 24.            |                              |       |            |      |
| 25.            |                              |       |            |      |
| 26.            |                              |       |            |      |
| 27.            |                              |       |            |      |
| 28.            |                              |       |            |      |
| 29.            |                              |       |            |      |
| 30.            |                              |       |            |      |

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

SB-4A (13-15)

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62041-1

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED73125

Level: (low/med) LOW

Date Received: 07/20/00

% Moisture: 17 decanted: (Y/N) N

Date Extracted: 07/24/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 07/31/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|               |                              |      |   |
|---------------|------------------------------|------|---|
| 108-95-2----- | Phenol                       | 400  | U |
| 111-44-4----- | bis(2-Chloroethyl) ether     | 400  | U |
| 95-57-8-----  | 2-Chlorophenol               | 400  | U |
| 541-73-1----- | 1,3-Dichlorobenzene          | 400  | U |
| 106-46-7----- | 1,4-Dichlorobenzene          | 400  | U |
| 95-50-1-----  | 1,2-Dichlorobenzene          | 400  | U |
| 95-48-7-----  | 2-Methylphenol               | 400  | U |
| 108-60-1----- | 2,2'-oxybis(1-Chloropropane) | 400  | U |
| 106-44-5----- | 4-Methylphenol               | 400  | U |
| 621-64-7----- | N-Nitroso-di-n-propylamine   | 400  | U |
| 67-72-1-----  | Hexachloroethane             | 400  | U |
| 98-95-3-----  | Nitrobenzene                 | 400  | U |
| 78-59-1-----  | Isophorone                   | 400  | U |
| 88-75-5-----  | 2-Nitrophenol                | 400  | U |
| 105-67-9----- | 2,4-Dimethylphenol           | 400  | U |
| 111-91-1----- | bis(2-Chloroethoxy) methane  | 400  | U |
| 120-83-2----- | 2,4-Dichlorophenol           | 400  | U |
| 120-82-1----- | 1,2,4-Trichlorobenzene       | 400  | U |
| 91-20-3-----  | Naphthalene                  | 400  | U |
| 106-47-8----- | 4-Chloroaniline              | 400  | U |
| 87-68-3-----  | Hexachlorobutadiene          | 400  | U |
| 59-50-7-----  | 4-Chloro-3-methylphenol      | 400  | U |
| 91-57-6-----  | 2-Methylnaphthalene          | 400  | U |
| 77-47-4-----  | Hexachlorocyclopentadiene    | 400  | U |
| 88-06-2-----  | 2,4,6-Trichlorophenol        | 400  | U |
| 95-95-4-----  | 2,4,5-Trichlorophenol        | 1000 | U |
| 91-58-7-----  | 2-Chloronaphthalene          | 400  | U |
| 88-74-4-----  | 2-Nitroaniline               | 1000 | U |
| 131-11-3----- | Dimethylphthalate            | 400  | U |
| 208-96-8----- | Acenaphthylene               | 400  | U |
| 606-20-2----- | 2,6-Dinitrotoluene           | 400  | U |
| 99-09-2-----  | 3-Nitroaniline               | 1000 | U |
| 83-32-9-----  | Acenaphthene                 | 400  | U |

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SB-4A (13-15)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62041-1

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED73125

Level: (low/med) LOW

Date Received: 07/20/00

% Moisture: 17 decanted: (Y/N) N

Date Extracted: 07/24/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 07/31/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg Q

|                |                            |      |   |
|----------------|----------------------------|------|---|
| 51-28-5-----   | 2,4-Dinitrophenol          | 1000 | U |
| 100-02-7-----  | 4-Nitrophenol              | 1000 | U |
| 132-64-9-----  | Dibenzofuran               | 400  | U |
| 121-14-2-----  | 2,4-Dinitrotoluene         | 400  | U |
| 84-66-2-----   | Diethylphthalate           | 400  | U |
| 86-73-7-----   | Fluorene                   | 400  | U |
| 7005-72-3----- | 4-Chlorophenyl-phenylether | 400  | U |
| 100-01-6-----  | 4-Nitroaniline             | 1000 | U |
| 534-52-1-----  | 4,6-Dinitro-2-methylphenol | 1000 | U |
| 86-30-6-----   | N-Nitrosodiphenylamine     | 400  | U |
| 101-55-3-----  | 4-Bromophenyl-phenylether  | 400  | U |
| 118-74-1-----  | Hexachlorobenzene          | 400  | U |
| 87-86-5-----   | Pentachlorophenol          | 1000 | U |
| 85-01-8-----   | Phenanthrene               | 64   | J |
| 120-12-7-----  | Anthracene                 | 400  | U |
| 86-74-8-----   | Carbazole                  | 400  | U |
| 84-74-2-----   | Di-n-butylphthalate        | 400  | U |
| 206-44-0-----  | Fluoranthene               | 46   | J |
| 129-00-0-----  | Pyrene                     | 49   | J |
| 85-68-7-----   | Butylbenzylphthalate       | 400  | U |
| 56-55-3-----   | Benzo(a)anthracene         | 400  | U |
| 91-94-1-----   | 3,3'-Dichlorobenzidine     | 400  | U |
| 218-01-9-----  | Chrysene                   | 400  | U |
| 117-81-7-----  | bis(2-Ethylhexyl)phthalate | 400  | U |
| 117-84-0-----  | Di-n-octylphthalate        | 400  | U |
| 205-99-2-----  | Benzo(b)fluoranthene       | 400  | U |
| 207-08-9-----  | Benzo(k)fluoranthene       | 400  | U |
| 50-32-8-----   | Benzo(a)pyrene             | 400  | U |
| 193-39-5-----  | Indeno(1,2,3-cd)pyrene     | 400  | U |
| 53-70-3-----   | Dibenzo(a,h)anthracene     | 400  | U |
| 191-24-2-----  | Benzo(g,h,i)perylene       | 400  | U |

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SB-4A (13-15)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62041-1

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED73125

Level: (low/med) LOW

Date Received: 07/20/00

% Moisture: 17 decanted: (Y/N) N

Date Extracted: 07/24/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 07/31/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

Number TICS found: 7

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg

| CAS NUMBER     | COMPOUND NAME                | RT    | EST. CONC. | Q   |
|----------------|------------------------------|-------|------------|-----|
| 1.             | Unknown                      | 1.43  | 2800       | JB  |
| 2. 010574-37-5 | 2-Pentene, 2,3-dimethyl      | 1.94  | 68000      | JBA |
| 3. 000123-42-2 | 2-Pentanone, 4-hydroxy-4-met | 2.50  | 7600       | JA  |
| 4.             | Unknown                      | 4.86  | 1800       | J   |
| 5.             | Unknown                      | 5.74  | 750        | J   |
| 6.             | Unknown Hydrocarbon          | 6.59  | 690        | J   |
| 7.             | Unknown Hydrocarbon          | 14.11 | 680        | J   |
| 8.             |                              |       |            |     |
| 9.             |                              |       |            |     |
| 10.            |                              |       |            |     |
| 11.            |                              |       |            |     |
| 12.            |                              |       |            |     |
| 13.            |                              |       |            |     |
| 14.            |                              |       |            |     |
| 15.            |                              |       |            |     |
| 16.            |                              |       |            |     |
| 17.            |                              |       |            |     |
| 18.            |                              |       |            |     |
| 19.            |                              |       |            |     |
| 20.            |                              |       |            |     |
| 21.            |                              |       |            |     |
| 22.            |                              |       |            |     |
| 23.            |                              |       |            |     |
| 24.            |                              |       |            |     |
| 25.            |                              |       |            |     |
| 26.            |                              |       |            |     |
| 27.            |                              |       |            |     |
| 28.            |                              |       |            |     |
| 29.            |                              |       |            |     |
| 30.            |                              |       |            |     |

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

DUP-1

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-5

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED73130

Level: (low/med) LOW

Date Received: 07/19/00

% Moisture: 26 decanted: (Y/N) N

Date Extracted: 07/24/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/01/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

|         |          |   |   |
|---------|----------|---|---|
| CAS NO. | COMPOUND | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) ug/kg | Q |
|---------|----------|---|---|

|   |      |   |
|---|------|---|
| 108-95-2-----Phenol                       | 450  | U |
| 111-44-4-----bis(2-Chloroethyl) ether     | 450  | U |
| 95-57-8-----2-Chlorophenol                | 450  | U |
| 541-73-1-----1,3-Dichlorobenzene          | 450  | U |
| 106-46-7-----1,4-Dichlorobenzene          | 450  | U |
| 95-50-1-----1,2-Dichlorobenzene           | 450  | U |
| 95-48-7-----2-Methylphenol                | 450  | U |
| 108-60-1-----2,2'-oxybis(1-Chloropropane) | 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 |
| 120-82-1-----1,2,4-Trichlorobenzene       | 450  | U |
| 91-20-3-----Naphthalene                   | 1600 |   |
| 106-47-8-----4-Chloroaniline              | 450  | U |
| 87-68-3-----Hexachlorobutadiene           | 450  | U |
| 59-50-7-----4-Chloro-3-methylphenol       | 450  | U |
| 91-57-6-----2-Methylnaphthalene           | 620  |   |
| 77-47-4-----Hexachlorocyclopentadiene     | 450  | U |
| 88-06-2-----2,4,6-Trichlorophenol         | 450  | U |
| 95-95-4-----2,4,5-Trichlorophenol         | 1100 | U |
| 91-58-7-----2-Chloronaphthalene           | 450  | U |
| 88-74-4-----2-Nitroaniline                | 1100 | U |
| 131-11-3-----Dimethylphthalate            | 450  | U |
| 208-96-8-----Acenaphthylene               | 810  |   |
| 606-20-2-----2,6-Dinitrotoluene           | 450  | U |
| 99-09-2-----3-Nitroaniline                | 1100 | U |
| 83-32-9-----Acenaphthene                  | 770  |   |

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

DUP-1

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-5

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED73130

Level: (low/med) LOW

Date Received: 07/19/00

% Moisture: 26 decanted: (Y/N) N

Date Extracted: 07/24/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/01/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

| CAS NO.   | COMPOUND                   | CONCENTRATION UNITS:<br>(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               | 1800  |     |
| 121-14-2  | 2,4-Dinitrotoluene         | 450   | U   |
| 84-66-2   | Diethylphthalate           | 450   | U   |
| 86-73-7   | Fluorene                   | 1200  |     |
| 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 U |
| 86-30-6   | N-Nitrosodiphenylamine     | 450   | U U |
| 101-55-3  | 4-Bromophenyl-phenylether  | 450   | U U |
| 118-74-1  | Hexachlorobenzene          | 450   | U U |
| 87-86-5   | Pentachlorophenol          | 1100  | U U |
| 85-01-8   | Phenanthrene               | 16000   | E J |
| 120-12-7  | Anthracene                 | 3300  | J   |
| 86-74-8   | Carbazole                  | 1400  | J   |
| 84-74-2   | Di-n-butylphthalate        | 450   | U U |
| 206-44-0  | Fluoranthene               | 12000   | E J |
| 129-00-0  | Pyrene                     | 10000   | E J |
| 85-68-7   | Butylbenzylphthalate       | 450   | U U |
| 56-55-3   | Benzo(a)anthracene         | 5800  | E J |
| 91-94-1   | 3,3'-Dichlorobenzidine     | 450   | U U |
| 218-01-9  | Chrysene                   | 4500  | E J |
| 117-81-7  | bis(2-Ethylhexyl)phthalate | 450   | U U |
| 117-84-0  | Di-n-octylphthalate        | 450   | U U |
| 205-99-2  | Benzo(b)fluoranthene       | 6600  | E J |
| 207-08-9  | Benzo(k)fluoranthene       | 1000  | J   |
| 50-32-8   | Benzo(a)pyrene             | 3600  | E J |
| 193-39-5  | Indeno(1,2,3-cd)pyrene     | 2300  | J   |
| 53-70-3   | Dibenzo(a,h)anthracene     | 820   | J   |
| 191-24-2  | Benzo(g,h,i)perylene       | 2600  | J   |

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

DUP-1

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-5

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED73130

Level: (low/med) LOW

Date Received: 07/19/00

% Moisture: 26 decanted: (Y/N) N

Date Extracted: 07/24/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/01/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg

Number TICS found: 10

| CAS NUMBER     | COMPOUND NAME                | RT    | EST. CONC. | Q   |
|----------------|------------------------------|-------|------------|-----|
| 1.             | Unknown                      | 1.08  | 1800       | J   |
| 2.             | Unknown                      | 1.43  | 2200       | JB  |
| 3. 000141-79-7 | 3-Penten-2-one, 4-methyl-    | 1.96  | 63000      | JBA |
| 4. 000123-42-2 | 2-Pentanone, 4-hydroxy-4-met | 2.52  | 5200       | JA  |
| 5.             | Unknown                      | 4.90  | 2300       | J   |
| 6.             | Unknown                      | 5.77  | 900        | J   |
| 7.             | Unknown                      | 6.82  | 650        | J   |
| 8.             | Substituted Naphthalene      | 11.39 | 1000       | J   |
| 9.             | Unknown PNA                  | 14.55 | 91         | J   |
| 10.            | Unknown PNA                  | 15.95 | 100        | J   |
| 11.            | benzo(e)pyrene               | 22.57 | 2700       | J   |
| 12.            |                              |       |            |     |
| 13.            |                              |       |            |     |
| 14.            |                              |       |            |     |
| 15.            |                              |       |            |     |
| 16.            |                              |       |            |     |
| 17.            |                              |       |            |     |
| 18.            |                              |       |            |     |
| 19.            |                              |       |            |     |
| 20.            |                              |       |            |     |
| 21.            |                              |       |            |     |
| 22.            |                              |       |            |     |
| 23.            |                              |       |            |     |
| 24.            |                              |       |            |     |
| 25.            |                              |       |            |     |
| 26.            |                              |       |            |     |
| 27.            |                              |       |            |     |
| 28.            |                              |       |            |     |
| 29.            |                              |       |            |     |
| 30.            |                              |       |            |     |

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

DUP-1 DL

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-5DL

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED80413

Level: (low/med) LOW

Date Received: 07/19/00

% Moisture: 26 decanted: (Y/N) N

Date Extracted: 07/24/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/04/00

Injection Volume: 2.0 (uL)

Dilution Factor: 10.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|               |                              |       |    |
|---------------|------------------------------|-------|----|
| 108-95-2----- | Phenol                       | 4500  | U  |
| 111-44-4----- | bis(2-Chloroethyl) ether     | 4500  | U  |
| 95-57-8-----  | 2-Chlorophenol               | 4500  | U  |
| 541-73-1----- | 1,3-Dichlorobenzene          | 4500  | U  |
| 106-46-7----- | 1,4-Dichlorobenzene          | 4500  | U  |
| 95-50-1-----  | 1,2-Dichlorobenzene          | 4500  | U  |
| 95-48-7-----  | 2-Methylphenol               | 4500  | U  |
| 108-60-1----- | 2,2'-oxybis(1-Chloropropane) | 4500  | U  |
| 106-44-5----- | 4-Methylphenol               | 4500  | U  |
| 621-64-7----- | N-Nitroso-di-n-propylamine   | 4500  | U  |
| 67-72-1-----  | Hexachloroethane             | 4500  | U  |
| 98-95-3-----  | Nitrobenzene                 | 4500  | U  |
| 78-59-1-----  | Isophorone                   | 4500  | U  |
| 88-75-5-----  | 2-Nitrophenol                | 4500  | U  |
| 105-67-9----- | 2,4-Dimethylphenol           | 4500  | U  |
| 111-91-1----- | bis(2-Chloroethoxy) methane  | 4500  | U  |
| 120-83-2----- | 2,4-Dichlorophenol           | 4500  | U  |
| 120-82-1----- | 1,2,4-Trichlorobenzene       | 4500  | U  |
| 91-20-3-----  | Naphthalene                  | 1400  | JD |
| 106-47-8----- | 4-Chloroaniline              | 4500  | U  |
| 87-68-3-----  | Hexachlorobutadiene          | 4500  | U  |
| 59-50-7-----  | 4-Chloro-3-methylphenol      | 4500  | U  |
| 91-57-6-----  | 2-Methylnaphthalene          | 740   | JD |
| 77-47-4-----  | Hexachlorocyclopentadiene    | 4500  | U  |
| 88-06-2-----  | 2,4,6-Trichlorophenol        | 4500  | U  |
| 95-95-4-----  | 2,4,5-Trichlorophenol        | 11000 | U  |
| 91-58-7-----  | 2-Chloronaphthalene          | 4500  | U  |
| 88-74-4-----  | 2-Nitroaniline               | 11000 | U  |
| 131-11-3----- | Dimethylphthalate            | 4500  | U  |
| 208-96-8----- | Acenaphthylene               | 610   | JD |
| 606-20-2----- | 2,6-Dinitrotoluene           | 4500  | U  |
| 99-09-2-----  | 3-Nitroaniline               | 11000 | U  |
| 83-32-9-----  | Acenaphthene                 | 630   | JD |



1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

DUP-1 DL

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-5DL

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED80413

Level: (low/med) LOW

Date Received: 07/19/00

% Moisture: 26 decanted: (Y/N) N

Date Extracted: 07/24/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/04/00

Injection Volume: 2.0 (uL)

Dilution Factor: 10.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|                |                            |       |    |
|----------------|----------------------------|-------|----|
| 51-28-5-----   | 2,4-Dinitrophenol          | 11000 | U  |
| 100-02-7-----  | 4-Nitrophenol              | 11000 | U  |
| 132-64-9-----  | Dibenzofuran               | 1600  | JD |
| 121-14-2-----  | 2,4-Dinitrotoluene         | 4500  | U  |
| 84-66-2-----   | Diethylphthalate           | 4500  | U  |
| 86-73-7-----   | Fluorene                   | 950   | JD |
| 7005-72-3----- | 4-Chlorophenyl-phenylether | 4500  | U  |
| 100-01-6-----  | 4-Nitroaniline             | 11000 | U  |
| 534-52-1-----  | 4,6-Dinitro-2-methylphenol | 11000 | U  |
| 86-30-6-----   | N-Nitrosodiphenylamine     | 4500  | U  |
| 101-55-3-----  | 4-Bromophenyl-phenylether  | 4500  | U  |
| 118-74-1-----  | Hexachlorobenzene          | 4500  | U  |
| 87-86-5-----   | Pentachlorophenol          | 11000 | U  |
| 85-01-8-----   | Phenanthrene               | 16000 | D  |
| 120-12-7-----  | Anthracene                 | 3000  | JD |
| 86-74-8-----   | Carbazole                  | 1200  | JD |
| 84-74-2-----   | Di-n-butylphthalate        | 4500  | U  |
| 206-44-0-----  | Fluoranthene               | 15000 | D  |
| 129-00-0-----  | Pyrene                     | 17000 | D  |
| 85-68-7-----   | Butylbenzylphthalate       | 4500  | U  |
| 56-55-3-----   | Benzo(a)anthracene         | 7100  | D  |
| 91-94-1-----   | 3,3'-Dichlorobenzidine     | 4500  | U  |
| 218-01-9-----  | Chrysene                   | 5900  | D  |
| 117-81-7-----  | bis(2-Ethylhexyl)phthalate | 4500  | U  |
| 117-84-0-----  | Di-n-octylphthalate        | 4500  | U  |
| 205-99-2-----  | Benzo(b)fluoranthene       | 7900  | D  |
| 207-08-9-----  | Benzo(k)fluoranthene       | 1800  | JD |
| 50-32-8-----   | Benzo(a)pyrene             | 4200  | JD |
| 193-39-5-----  | Indeno(1,2,3-cd)pyrene     | 2000  | JD |
| 53-70-3-----   | Dibenzo(a,h)anthracene     | 4500  | U  |
| 191-24-2-----  | Benzo(g,h,i)perylene       | 2200  | JD |

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

DUP-1 DL

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-5 DL

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED80413

Level: (low/med) LOW

Date Received: 07/19/00

% Moisture: 26 decanted: (Y/N) N

Date Extracted: 07/24/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/04/00

Injection Volume: 2.0 (uL)

Dilution Factor: 10.0

GPC Cleanup: (Y/N) Y pH: 7.0

Number TICS found: 16

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg

| CAS NUMBER      | COMPOUND NAME                | RT    | EST. CONC. | Q   |
|-----------------|------------------------------|-------|------------|-----|
| 1.              | Unknown                      | 0.82  | 7000       | J   |
| 2.              | Unknown                      | 1.12  | 5500       | J   |
| 3.              | Unknown                      | 1.47  | 110000     | JB  |
| 4. 000123-42-2  | 2-Pentanone, 4-hydroxy-4-met | 1.92  | 10000      | JBA |
| 5.              | Unknown                      | 4.13  | 4100       | J   |
| 6.              | Unknown                      | 5.09  | 1600       | J   |
| 7.              | Substituted Naphthalene      | 9.45  | 910        | J   |
| 8.              | Substituted Naphthalene      | 10.48 | 1300       | J   |
| 9.              | Substituted Naphthalene      | 10.62 | 1000       | J   |
| 10.             | Unknown PAH                  | 13.62 | 1700       | J   |
| 11. 000132-65-0 | Dibenzothiophene             | 13.72 | 1100       | J   |
| 12.             | Unknown PAH                  | 14.90 | 1200       | J   |
| 13.             | Unknown PAH                  | 14.95 | 1500       | J   |
| 14. 000084-65-1 | 9,10-Anthracenedione         | 15.52 | 2900       | J   |
| 15.             | Unknown PAH                  | 16.06 | 980        | J   |
| 16.             | Unknown                      | 21.00 | 9000       | J   |
| 17.             | benzo(e) pyrene              | 21.15 | 2900       | J   |
| 18.             |                              |       |            |     |
| 19.             |                              |       |            |     |
| 20.             |                              |       |            |     |
| 21.             |                              |       |            |     |
| 22.             |                              |       |            |     |
| 23.             |                              |       |            |     |
| 24.             |                              |       |            |     |
| 25.             |                              |       |            |     |
| 26.             |                              |       |            |     |
| 27.             |                              |       |            |     |
| 28.             |                              |       |            |     |
| 29.             |                              |       |            |     |
| 30.             |                              |       |            |     |

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

SB-1 (4-6)

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-1

Sample wt/vol: 30.3 (g/mL) g

Lab File ID: HP13A\0806C027

% Moisture: 12 decanted: (Y/N) N

Date Received: 07/19/00

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 07/24/00

Concentrated Extract Volume: 5000 (uL)

Date Analyzed: 08/07/00

Injection Volume: 1.0 (uL)

Dilution Factor: 10.0

GPC Cleanup: (Y/N) Y pH: 7.0

Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|                 |                     |      |   |
|-----------------|---------------------|------|---|
| 319-84-6-----   | alpha-BHC           | 19   | U |
| 319-85-7-----   | beta-BHC            | 19   | U |
| 319-86-8-----   | delta-BHC           | 19   | U |
| 58-89-9-----    | gamma-BHC (Lindane) | 130  |   |
| 76-44-8-----    | Heptachlor          | 66   | P |
| 309-00-2-----   | Aldrin              | 19   | U |
| 1024-57-3-----  | Heptachlor epoxide  | 240  | P |
| 959-98-8-----   | Endosulfan I        | 19   | U |
| 60-57-1-----    | Dieldrin            | 38   | U |
| 72-55-9-----    | 4,4'-DDE            | 430  | P |
| 72-20-8-----    | Endrin              | 38   | U |
| 332-13-659----- | Endosulfan II       | 38   | U |
| 72-54-8-----    | 4,4'-DDD            | 170  | P |
| 103-10-78-----  | Endosulfan sulfate  | 38   | U |
| 50-29-3-----    | 4,4'-DDT            | 38   | U |
| 72-43-5-----    | Methoxychlor        | 190  | U |
| 53494-70-5----- | Endrin ketone       | 38   | U |
| 7421-36-3-----  | Endrin aldehyde     | 38   | U |
| 5103-71-9-----  | alpha-Chlordane     | 19   | U |
| 5103-7-42-----  | gamma-Chlordane     | 19   | U |
| 8001-35-2-----  | Toxaphene           | 1900 | U |
| 12674-11-2----- | Aroclor-1016        | 380  | U |
| 11104-28-2----- | Aroclor-1221        | 750  | U |
| 11141-16-5----- | Aroclor-1232        | 380  | U |
| 53469-21-9----- | Aroclor-1242        | 380  | U |
| 12672-29-6----- | Aroclor-1248        | 380  | U |
| 11097-69-1----- | Aroclor-1254        | 380  | U |
| 11096-82-5----- | Aroclor-1260        | 380  | U |

FORM I PEST

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

SB-1 (4-6) DL

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-1DL

Sample wt/vol: 30.3 (g/mL) g

Lab File ID: HP13A\0806C029

% Moisture: 12 decanted: (Y/N) N

Date Received: 07/19/00

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 07/24/00

Concentrated Extract Volume: 5000 (uL)

Date Analyzed: 08/07/00

Injection Volume: 1.0 (uL)

Dilution Factor: 100.0

GPC Cleanup: (Y/N) Y pH: 7.0

Sulfur Cleanup: (Y/N) N

CAS NO.

COMPOUND

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg Q

|                 |                     |       |    |
|-----------------|---------------------|-------|----|
| 319-84-6-----   | alpha-BHC           | 160   | UD |
| 319-85-7-----   | beta-BHC            | 160   | UD |
| 319-86-8-----   | delta-BHC           | 160   | UD |
| 58-89-9-----    | gamma-BHC (Lindane) | 150   | JD |
| 76-44-8-----    | Heptachlor          | 160   | UD |
| 309-00-2-----   | Aldrin              | 160   | UD |
| 1024-57-3-----  | Heptachlor epoxide  | 230   | PD |
| 959-98-8-----   | Endosulfan I        | 160   | UD |
| 60-57-1-----    | Dieldrin            | 330   | UD |
| 72-55-9-----    | 4,4'-DDE            | 440   | PD |
| 72-20-8-----    | Endrin              | 330   | UD |
| 332-13-659----- | Endosulfan II       | 330   | UD |
| 72-54-8-----    | 4,4'-DDD            | 330   | UD |
| 103-10-78-----  | Endosulfan sulfate  | 330   | UD |
| 50-29-3-----    | 4,4'-DDT            | 330   | UD |
| 72-43-5-----    | Methoxychlor        | 1600  | UD |
| 53494-70-5----- | Endrin ketone       | 330   | UD |
| 7421-36-3-----  | Endrin aldehyde     | 330   | UD |
| 5103-71-9-----  | alpha-Chlordane     | 160   | UD |
| 5103-7-42-----  | gamma-Chlordane     | 160   | UD |
| 8001-35-2-----  | Toxaphene           | 16000 | UD |
| 12674-11-2----- | Aroclor-1016        | 3300  | UD |
| 11104-28-2----- | Aroclor-1221        | 6600  | UD |
| 11141-16-5----- | Aroclor-1232        | 3300  | UD |
| 53469-21-9----- | Aroclor-1242        | 3300  | UD |
| 12672-29-6----- | Aroclor-1248        | 3300  | UD |
| 11097-69-1----- | Aroclor-1254        | 3300  | UD |
| 11096-82-5----- | Aroclor-1260        | 3300  | UD |

FORM I PEST

\* Sample diluted due to matrix interference.

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

SB-2 (10-12)

Lab Code: Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-2

Sample wt/vol: 30.2 (g/mL) g

Lab File ID: HP13A\0806c031

% Moisture: 25 decanted: (Y/N) N

Date Received: 07/19/00

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 07/24/00

Concentrated Extract Volume: 5000 (uL)

Date Analyzed: 08/07/00

Injection Volume: 1.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) ug/kg Q

|            |                     |     |   |
|------------|---------------------|-----|---|
| 319-84-6   | alpha-BHC           | 1.6 | U |
| 319-85-7   | beta-BHC            | 1.6 | U |
| 319-86-8   | delta-BHC           | 1.6 | U |
| 58-89-9    | gamma-BHC (Lindane) | 1.6 | U |
| 76-44-8    | Heptachlor          | 1.6 | U |
| 309-00-2   | Aldrin              | 1.6 | U |
| 1024-57-3  | Heptachlor epoxide  | 1.6 | U |
| 959-98-8   | Endosulfan I        | 1.6 | U |
| 60-57-1    | Dieldrin            | 3.3 | U |
| 72-55-9    | 4,4'-DDE            | 3.3 | U |
| 72-20-8    | Endrin              | 3.3 | U |
| 332-13-659 | Endosulfan II       | 3.3 | U |
| 72-54-8    | 4,4'-DDD            | 3.3 | U |
| 103-10-78  | Endosulfan sulfate  | 3.3 | U |
| 50-29-3    | 4,4'-DDT            | 3.3 | U |
| 72-43-5    | Methoxychlor        | 16  | U |
| 53494-70-5 | Endrin ketone       | 3.3 | U |
| 7421-36-3  | Endrin aldehyde     | 3.3 | U |
| 5103-71-9  | alpha-Chlordane     | 1.6 | U |
| 5103-7-42  | gamma-Chlordane     | 1.6 | U |
| 8001-35-2  | Toxaphene           | 160 | U |
| 12674-11-2 | Aroclor-1016        | 33  | U |
| 11104-28-2 | Aroclor-1221        | 66  | U |
| 11141-16-5 | Aroclor-1232        | 33  | U |
| 53469-21-9 | Aroclor-1242        | 33  | U |
| 12672-29-6 | Aroclor-1248        | 33  | U |
| 11097-69-1 | Aroclor-1254        | 33  | U |
| 11096-82-5 | Aroclor-1260        | 33  | U |

FORM I PEST

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

SB-2 (10-12) DL

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-2DL

Sample wt/vol: 30.2 (g/mL) g

Lab File ID: HP13A\0806c033

% Moisture: 25 decanted: (Y/N) N

Date Received: 07/19/00

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 07/24/00

Concentrated Extract Volume: 5000 (uL)

Date Analyzed: 08/07/00

Injection Volume: 1.0 (uL)

Dilution Factor: 10.0

GPC Cleanup: (Y/N) Y pH: 7.0

Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|                 |                     |      |    |
|-----------------|---------------------|------|----|
| 319-84-6-----   | alpha-BHC           | 16   | UD |
| 319-85-7-----   | beta-BHC            | 16   | UD |
| 319-86-8-----   | delta-BHC           | 16   | UD |
| 58-89-9-----    | gamma-BHC (Lindane) | 16   | UD |
| 76-44-8-----    | Heptachlor          | 16   | UD |
| 309-00-2-----   | Aldrin              | 16   | UD |
| 1024-57-3-----  | Heptachlor epoxide  | 16   | UD |
| 959-98-8-----   | Endosulfan I        | 16   | UD |
| 60-57-1-----    | Dieldrin            | 33   | UD |
| 72-55-9-----    | 4,4'-DDE            | 33   | UD |
| 72-20-8-----    | Endrin              | 33   | UD |
| 332-13-659----- | Endosulfan II       | 33   | UD |
| 72-54-8-----    | 4,4'-DDD            | 33   | UD |
| 103-10-78-----  | Endosulfan sulfate  | 33   | UD |
| 50-29-3-----    | 4,4'-DDT            | 33   | UD |
| 72-43-5-----    | Methoxychlor        | 160  | UD |
| 53494-70-5----- | Endrin ketone       | 33   | UD |
| 7421-36-3-----  | Endrin aldehyde     | 33   | UD |
| 5103-71-9-----  | alpha-Chlordane     | 16   | UD |
| 5103-7-42-----  | gamma-Chlordane     | 16   | UD |
| 8001-35-2-----  | Toxaphene           | 1600 | UD |
| 12674-11-2----- | Aroclor-1016        | 330  | UD |
| 11104-28-2----- | Aroclor-1221        | 660  | UD |
| 11141-16-5----- | Aroclor-1232        | 330  | UD |
| 53469-21-9----- | Aroclor-1242        | 330  | UD |
| 12672-29-6----- | Aroclor-1248        | 330  | UD |
| 11097-69-1----- | Aroclor-1254        | 330  | UD |
| 11096-82-5----- | Aroclor-1260        | 330  | UD |

FORM I PEST

*\* Sample diluted due to matrix interference*

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SB-2 (19-23)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-3

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: HP13A\0806C035

Moisture: 49 decanted: (Y/N) N

Date Received: 07/19/00

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 07/24/00

Concentrated Extract Volume: 5000 (uL)

Date Analyzed: 08/07/00

Injection Volume: 1.0 (uL)

Dilution Factor: 1.0

GC Cleanup: (Y/N) Y pH: 7.0

Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|            |                               |     |    |   |
|------------|-------------------------------|-----|----|---|
| 319-84-6   | <del>alpha-BHC</del>          | 8.7 | P  | R |
| 319-85-7   | beta-BHC                      | 1.7 | U  |   |
| 319-86-8   | <del>delta-BHC</del>          | 44  | P  | R |
| 58-89-9    | gamma-BHC (Lindane)           | 1.7 | U  |   |
| 76-44-8    | <del>Heptachlor</del>         | 23  | P  | R |
| 309-00-2   | Aldrin                        | 1.7 | U  |   |
| 1024-57-3  | Heptachlor epoxide            | 1.7 | U  |   |
| 959-98-8   | Endosulfan I                  | 1.7 | U  |   |
| 60-57-1    | <del>Dieldrin</del>           | 21  | P  | R |
| 72-55-9    | <del>4,4'-DDE</del>           | 99  | P  | R |
| 72-20-8    | Endrin                        | 3.3 | U  |   |
| 332-13-659 | Endosulfan II                 | 3.3 | U  |   |
| 72-54-8    | 4,4'-DDD                      | 3.3 | U  |   |
| 103-10-78  | <del>Endosulfan sulfate</del> | 22  | P  | R |
| 50-29-3    | <del>4,4'-DDT</del>           | 68  |    |   |
| 72-43-5    | <del>Methoxychlor</del>       | 530 | EP | R |
| 53494-70-5 | Endrin ketone                 | 3.3 | U  |   |
| 7421-36-3  | Endrin aldehyde               | 3.3 | U  |   |
| 5103-71-9  | <del>alpha-Chlordane</del>    | 26  | P  | R |
| 5103-7-42  | gamma-Chlordane               | 1.7 | U  |   |
| 8001-35-2  | Toxaphene                     | 170 | U  |   |
| 12674-11-2 | Aroclor-1016                  | 33  | U  |   |
| 11104-28-2 | Aroclor-1221                  | 66  | U  |   |
| 11141-16-5 | Aroclor-1232                  | 33  | U  |   |
| 53469-21-9 | Aroclor-1242                  | 33  | U  |   |
| 12672-29-6 | Aroclor-1248                  | 33  | U  |   |
| 11097-69-1 | Aroclor-1254                  | 33  | U  |   |
| 11096-82-5 | Aroclor-1260                  | 33  | U  |   |

FORM I PEST

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SB-2 (19-23) DL

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-3DL

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: HP13A\0806C040

% Moisture: 49 decanted: (Y/N) N

Date Received: 07/19/00

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 07/24/00

Concentrated Extract Volume: 5000 (uL)

Date Analyzed: 08/07/00

Injection Volume: 1.0 (uL)

Dilution Factor: 10.0

GFC Cleanup: (Y/N) Y pH: 7.0

Sulfur Cleanup: (Y/N) N

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg Q

|                 |                     |      |    |
|-----------------|---------------------|------|----|
| 319-84-6-----   | alpha-BHC           | 17   | UD |
| 319-85-7-----   | beta-BHC            | 17   | UD |
| 319-86-8-----   | delta-BHC           | 17   | UD |
| 58-89-9-----    | gamma-BHC (Lindane) | 17   | UD |
| 76-44-8-----    | Heptachlor          | 35   | UD |
| 309-00-2-----   | Aldrin              | 17   | UD |
| 1024-57-3-----  | Heptachlor epoxide  | 61   | UD |
| 959-98-8-----   | Endosulfan I        | 17   | UD |
| 60-57-1-----    | Dieldrin            | 33   | UD |
| 72-55-9-----    | 4,4'-DDE            | 100  | UD |
| 72-20-8-----    | Endrin              | 33   | UD |
| 332-13-659----- | Endosulfan II       | 33   | UD |
| 72-54-8-----    | 4,4'-DDD            | 33   | UD |
| 103-10-78-----  | Endosulfan sulfate  | 33   | UD |
| 50-29-3-----    | 4,4'-DDT            | 33   | UD |
| 72-43-5-----    | Methoxychlor        | 630  | UD |
| 53494-70-5----- | Endrin ketone       | 33   | UD |
| 7421-36-3-----  | Endrin aldehyde     | 33   | UD |
| 5103-71-9-----  | alpha-Chlordane     | 17   | UD |
| 5103-7-42-----  | gamma-Chlordane     | 17   | UD |
| 8001-35-2-----  | Toxaphene           | 1700 | UD |
| 12674-11-2----- | Aroclor-1016        | 330  | UD |
| 11104-28-2----- | Aroclor-1221        | 660  | UD |
| 11141-16-5----- | Aroclor-1232        | 330  | UD |
| 53469-21-9----- | Aroclor-1242        | 330  | UD |
| 12672-29-6----- | Aroclor-1248        | 330  | UD |
| 11097-69-1----- | Aroclor-1254        | 330  | UD |
| 11096-82-5----- | Aroclor-1260        | 330  | UD |

FORM I PEST



1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

SB-3 (12-14)

Lab Code: Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-4

Sample wt/vol: 30.2 (g/mL) g

Lab File ID: HP13A\0806C042

% Moisture: 17 decanted: (Y/N) N

Date Received: 07/19/00

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 07/24/00

Concentrated Extract Volume: 5000 (uL)

Date Analyzed: 08/07/00

Injection Volume: 1.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

Sulfur Cleanup: (Y/N) N

| CAS NO. | COMPOUND | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) ug/kg | Q |
|---------|----------|---|---|
|---------|----------|---|---|

|                 |                     |     |   |
|-----------------|---------------------|-----|---|
| 319-84-6-----   | alpha-BHC           | 1.6 | U |
| 319-85-7-----   | beta-BHC            | 1.6 | U |
| 319-86-8-----   | delta-BHC           | 1.6 | U |
| 58-89-9-----    | gamma-BHC (Lindane) | 1.6 | U |
| 76-44-8-----    | Heptachlor          | 1.6 | U |
| 309-00-2-----   | Aldrin              | 1.6 | U |
| 1024-57-3-----  | Heptachlor epoxide  | 1.6 | U |
| 959-98-8-----   | Endosulfan I        | 1.6 | U |
| 60-57-1-----    | Dieldrin            | 3.3 | U |
| 72-55-9-----    | 4,4'-DDE            | 3.3 | U |
| 72-20-8-----    | Endrin              | 3.3 | U |
| 332-13-659----  | Endosulfan II       | 3.3 | U |
| 72-54-8-----    | 4,4'-DDD            | 3.3 | U |
| 103-10-78-----  | Endosulfan sulfate  | 3.3 | U |
| 50-29-3-----    | 4,4'-DDT            | 3.3 | U |
| 72-43-5-----    | Methoxychlor        | 16  | U |
| 53494-70-5----- | Endrin ketone       | 3.3 | U |
| 7421-36-3-----  | Endrin aldehyde     | 3.3 | U |
| 5103-71-9-----  | alpha-Chlordane     | 1.6 | U |
| 5103-7-42-----  | gamma-Chlordane     | 1.6 | U |
| 8001-35-2-----  | Toxaphene           | 160 | U |
| 12674-11-2----- | Aroclor-1016        | 33  | U |
| 11104-28-2----- | Aroclor-1221        | 66  | U |
| 11141-16-5----- | Aroclor-1232        | 33  | U |
| 53469-21-9----- | Aroclor-1242        | 33  | U |
| 12672-29-6----- | Aroclor-1248        | 33  | U |
| 11097-69-1----- | Aroclor-1254        | 33  | U |
| 11096-82-5----- | Aroclor-1260        | 33  | U |

FORM I PEST

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SB-4A (13-15)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62041-1

Sample wt/vol: 30.3 (g/mL) g

Lab File ID: HP13A\0806C048

% Moisture: 17 decanted: (Y/N) N

Date Received: 07/20/00

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 07/24/00

Concentrated Extract Volume: 5000 (uL)

Date Analyzed: 08/07/00

Injection Volume: 1.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|                 |                     |     |   |
|-----------------|---------------------|-----|---|
| 319-84-6-----   | alpha-BHC           | 1.6 | U |
| 319-85-7-----   | beta-BHC            | 1.6 | U |
| 319-86-8-----   | delta-BHC           | 1.6 | U |
| 58-89-9-----    | gamma-BHC (Lindane) | 1.6 | U |
| 76-44-8-----    | Heptachlor          | 1.6 | U |
| 309-00-2-----   | Aldrin              | 1.6 | U |
| 1024-57-3-----  | Heptachlor epoxide  | 1.6 | U |
| 959-98-8-----   | Endosulfan I        | 1.6 | U |
| 60-57-1-----    | Dieldrin            | 3.3 | U |
| 72-55-9-----    | 4,4'-DDE            | 3.3 | U |
| 72-20-8-----    | Endrin              | 3.3 | U |
| 332-13-659----- | Endosulfan II       | 3.3 | U |
| 72-54-8-----    | 4,4'-DDD            | 3.3 | U |
| 103-10-78-----  | Endosulfan sulfate  | 3.3 | U |
| 50-29-3-----    | 4,4'-DDT            | 3.3 | U |
| 72-43-5-----    | Methoxychlor        | 16  | U |
| 53494-70-5----- | Endrin ketone       | 3.3 | U |
| 7421-36-3-----  | Endrin aldehyde     | 3.3 | U |
| 5103-71-9-----  | alpha-Chlordane     | 1.6 | U |
| 5103-7-42-----  | gamma-Chlordane     | 1.6 | U |
| 8001-35-2-----  | Toxaphene           | 160 | U |
| 12674-11-2----- | Aroclor-1016        | 33  | U |
| 11104-28-2----- | Aroclor-1221        | 66  | U |
| 11141-16-5----- | Aroclor-1232        | 33  | U |
| 53469-21-9----- | Aroclor-1242        | 33  | U |
| 12672-29-6----- | Aroclor-1248        | 33  | U |
| 11097-69-1----- | Aroclor-1254        | 33  | U |
| 11096-82-5----- | Aroclor-1260        | 33  | U |

FORM I PEST

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

DUP-1

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-5

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: HP13A\0806C044

% Moisture: 26 decanted: (Y/N) N

Date Received: 07/19/00

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 07/24/00

Concentrated Extract Volume: 5000 (uL)

Date Analyzed: 08/07/00

Injection Volume: 1.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

Sulfur Cleanup: (Y/N) N

| CAS NO. | COMPOUND | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) ug/kg | Q |
|---------|----------|---|---|
|---------|----------|---|---|

|                 |                               |     |                 |
|-----------------|-------------------------------|-----|-----------------|
| 319-84-6-----   | <del>alpha-BHC</del>          | 8.1 | P <del>82</del> |
| 319-85-7-----   | beta-BHC                      | 1.7 | U               |
| 319-86-8-----   | delta-BHC                     | 1.7 | U               |
| 58-89-9-----    | gamma-BHC (Lindane)           | 1.7 | U               |
| 76-44-8-----    | <del>Heptachlor</del>         | 9.6 | P <del>82</del> |
| 309-00-2-----   | Aldrin                        | 1.7 | U               |
| 1024-57-3-----  | <del>Heptachlor epoxide</del> | 12  | P <del>82</del> |
| 959-98-8-----   | Endosulfan I                  | 1.7 | U               |
| 60-57-1-----    | Dieldrin                      | 3.3 | U               |
| 72-55-9-----    | 4,4'-DDE                      | 3.3 | U               |
| 72-20-8-----    | Endrin                        | 3.3 | U               |
| 332-13-659----- | Endosulfan II                 | 3.3 | U               |
| 72-54-8-----    | 4,4'-DDD                      | 3.3 | U               |
| 103-10-78-----  | Endosulfan sulfate            | 3.3 | U               |
| 50-29-3-----    | 4,4'-DDT                      | 3.3 | U               |
| 72-43-5-----    | Methoxychlor                  | 17  | U               |
| 53494-70-5----- | Endrin ketone                 | 3.3 | U               |
| 7421-36-3-----  | Endrin aldehyde               | 3.3 | U               |
| 5103-71-9-----  | alpha-Chlordane               | 1.7 | U               |
| 5103-7-42-----  | gamma-Chlordane               | 1.7 | U               |
| 8001-35-2-----  | Toxaphene                     | 170 | U               |
| 12674-11-2----- | Aroclor-1016                  | 33  | U               |
| 11104-28-2----- | Aroclor-1221                  | 66  | U               |
| 11141-16-5----- | Aroclor-1232                  | 33  | U               |
| 53469-21-9----- | Aroclor-1242                  | 33  | U               |
| 12672-29-6----- | Aroclor-1248                  | 33  | U               |
| 11097-69-1----- | Aroclor-1254                  | 33  | U               |
| 11096-82-5----- | Aroclor-1260                  | 33  | U               |

FORM I PEST

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

DUP-1 DL

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) SOIL

Lab Sample ID: L62005-5DL

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: HP13A\0806C046

% Moisture: 26 decanted: (Y/N) N

Date Received: 07/19/00

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 07/24/00

Concentrated Extract Volume: 5000 (uL)

Date Analyzed: 08/07/00

Injection Volume: 1.0 (uL)

Dilution Factor: 10.0

GPC Cleanup: (Y/N) Y pH: 7.0

Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|                 |                     |      |    |
|-----------------|---------------------|------|----|
| 319-84-6-----   | alpha-BHC           | 17   | UD |
| 319-85-7-----   | beta-BHC            | 17   | UD |
| 319-86-8-----   | delta-BHC           | 17   | UD |
| 58-89-9-----    | gamma-BHC (Lindane) | 17   | UD |
| 76-44-8-----    | Heptachlor          | 17   | UD |
| 309-00-2-----   | Aldrin              | 17   | UD |
| 1024-57-3-----  | Heptachlor epoxide  | 17   | UD |
| 959-98-8-----   | Endosulfan I        | 17   | UD |
| 60-57-1-----    | Dieldrin            | 33   | UD |
| 72-55-9-----    | 4,4'-DDE            | 33   | UD |
| 72-20-8-----    | Endrin              | 33   | UD |
| 332-13-659----- | Endosulfan II       | 33   | UD |
| 72-54-8-----    | 4,4'-DDD            | 33   | UD |
| 103-10-78-----  | Endosulfan sulfate  | 33   | UD |
| 50-29-3-----    | 4,4'-DDT            | 33   | UD |
| 72-43-5-----    | Methoxychlor        | 170  | UD |
| 53494-70-5----- | Endrin ketone       | 33   | UD |
| 7421-36-3-----  | Endrin aldehyde     | 33   | UD |
| 5103-71-9-----  | alpha-Chlordane     | 17   | UD |
| 5103-7-42-----  | gamma-Chlordane     | 17   | UD |
| 8001-35-2-----  | Toxaphene           | 1700 | UD |
| 12674-11-2----- | Aroclor-1016        | 330  | UD |
| 11104-28-2----- | Aroclor-1221        | 660  | UD |
| 11141-16-5----- | Aroclor-1232        | 330  | UD |
| 53469-21-9----- | Aroclor-1242        | 330  | UD |
| 12672-29-6----- | Aroclor-1248        | 330  | UD |
| 11097-69-1----- | Aroclor-1254        | 330  | UD |
| 11096-82-5----- | Aroclor-1260        | 330  | UD |

FORM I PEST

\* Sample diluted due to matrix interference.

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

RINSE BLANK

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L62005

Matrix: (soil/water) WATER

Lab Sample ID: L62041-5

Sample wt/vol: 1050 (g/mL) mL

Lab File ID: HP13A\0806C026

% Moisture: decanted: (Y/N) N

Date Received: 07/20/00

Extraction: (SepF/Cont/Sonc) SEPF

Date Extracted: 07/25/00

Concentrated Extract Volume: 10000 (uL)

Date Analyzed: 08/07/00

Injection Volume: 1.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: 7.0

Sulfur Cleanup: (Y/N) N

| CAS NO. | COMPOUND | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) ug/l | Q |
|---------|----------|--|---|
|---------|----------|--|---|

|                 |                     |      |   |
|-----------------|---------------------|------|---|
| 319-84-6-----   | alpha-BHC           | 0.05 | U |
| 319-85-7-----   | beta-BHC            | 0.05 | U |
| 319-86-8-----   | delta-BHC           | 0.05 | U |
| 58-89-9-----    | gamma-BHC (Lindane) | 0.05 | U |
| 76-44-8-----    | Heptachlor          | 0.05 | U |
| 309-00-2-----   | Aldrin              | 0.05 | U |
| 1024-57-3-----  | Heptachlor epoxide  | 0.05 | U |
| 959-98-8-----   | Endosulfan I        | 0.05 | U |
| 60-57-1-----    | Dieldrin            | 0.1  | U |
| 72-55-9-----    | 4,4'-DDE            | 0.1  | U |
| 72-20-8-----    | Endrin              | 0.1  | U |
| 332-13-659----  | Endosulfan II       | 0.1  | U |
| 72-54-8-----    | 4,4'-DDD            | 0.1  | U |
| 103-10-78-----  | Endosulfan sulfate  | 0.1  | U |
| 50-29-3-----    | 4,4'-DDT            | 0.1  | U |
| 72-43-5-----    | Methoxychlor        | 0.48 | U |
| 53494-70-5----- | Endrin ketone       | 0.1  | U |
| 7421-36-3-----  | Endrin aldehyde     | 0.1  | U |
| 5103-71-9-----  | alpha-Chlordane     | 0.05 | U |
| 5103-7-42-----  | gamma-Chlordane     | 0.05 | U |
| 8001-35-2-----  | Toxaphene           | 4.8  | U |
| 12674-11-2----- | Aroclor-1016        | 0.95 | U |
| 11104-28-2----- | Aroclor-1221        | 1.9  | U |
| 11141-16-5----- | Aroclor-1232        | 0.95 | U |
| 53469-21-9----- | Aroclor-1242        | 0.95 | U |
| 12672-29-6----- | Aroclor-1248        | 0.95 | U |
| 11097-69-1----- | Aroclor-1254        | 0.95 | U |
| 11096-82-5----- | Aroclor-1260        | 0.95 | U |

FORM I PEST

## NYSDEC ASP

1  
INORGANIC ANALYSIS DATA SHEET

NYSDEC SAMPLE NO.

SB-1(4-6)

Lab Name: Galson Laboratories

Contract: BBL

Lab Code: 11626

Case No.:

SAS No.:

SDG No.: L62005

Matrix (soil/water): Soil

Lab Sample ID: L62005-1

Level (low/med): LOW

Date Received: 07/19/00

% Solids: 88.3

Concentration Units (ug/L or mg/kg dry weight): mg/kg

| CAS No.   | Analyte   | Concentration | C | Q    | M  |
|-----------|-----------|---------------|---|------|----|
| 7429-90-5 | Aluminum  | 5110          | — |      | P  |
| 7440-36-0 | Antimony  | 1.1           | U | N W  | P  |
| 7440-38-2 | Arsenic   | 3.9           |   |      | P  |
| 7440-39-3 | Barium    | 20.5          | B |      | P  |
| 7440-41-7 | Beryllium | 0.52          | B | u    | P  |
| 7440-43-9 | Cadmium   | 0.45          | U |      | P  |
| 7440-70-2 | Calcium   | 761           | B | *    | P  |
| 7440-47-3 | Chromium  | 7.5           |   |      | P  |
| 7440-48-4 | Cobalt    | 4.4           | B |      | P  |
| 7440-50-8 | Copper    | 27.5          |   |      | P  |
| 7439-89-6 | Iron      | 12800         |   |      | P  |
| 7439-92-1 | Lead      | 4.7           |   | N J  | P  |
| 7439-95-4 | Magnesium | 2090          |   | *    | P  |
| 7439-96-5 | Manganese | 216           |   | *N J | P  |
| 7439-97-6 | Mercury   | 0.057         | U |      | AV |
| 7440-02-0 | Nickel    | 11.0          |   |      | P  |
| 7440-09-7 | Potassium | 892           | B |      | P  |
| 7782-49-2 | Selenium  | 1.1           | B |      | P  |
| 7440-22-4 | Silver    | 0.45          | U |      | P  |
| 7440-23-5 | Sodium    | 94.2          | U |      | P  |
| 7440-28-0 | Thallium  | 1.4           | U |      | P  |
| 7440-62-2 | Vanadium  | 9.7           | B |      | P  |
| 7440-66-6 | Zinc      | 21.1          |   |      | P  |
| 57-12-5   | Cyanide   | 0.55          | U |      | C  |

Color Before: brown

Clarity Before:

Texture: med

Color After: yellow

Clarity After: clear

Artifacts: rocks

Comments:

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## NYSDEC ASP

1  
INORGANIC ANALYSIS DATA SHEET

NYSDEC SAMPLE NO.

SB-2(10-12)

Lab Name: Galson Laboratories

Contract: BBL

Lab Code: 11626

Case No.:

SAS No.:

SDG No.: L62005

Matrix (soil/water): Soil

Lab Sample ID: L62005-2

Level (low/med): LOW

Date Received: 07/19/00

% Solids: 75.0

Concentration Units (ug/L or mg/kg dry weight): mg/kg

| CAS No.   | Analyte   | Concentration | C | Q    | M  |
|-----------|-----------|---------------|---|------|----|
| 7429-90-5 | Aluminum  | 4790          |   |      | P  |
| 7440-36-0 | Antimony  | 47.3          |   | N J  | P  |
| 7440-38-2 | Arsenic   | 19.5          |   |      | P  |
| 7440-39-3 | Barium    | 688           |   |      | P  |
| 7440-41-7 | Beryllium | 0.54          | B | U    | P  |
| 7440-43-9 | Cadmium   | 4.9           |   |      | P  |
| 7440-70-2 | Calcium   | 12700         |   | *    | P  |
| 7440-47-3 | Chromium  | 17.7          |   |      | P  |
| 7440-48-4 | Cobalt    | 11.5          | B |      | P  |
| 7440-50-8 | Copper    | 2490          |   |      | P  |
| 7439-89-6 | Iron      | 62000         |   |      | P  |
| 7439-92-1 | Lead      | 1700          |   | N J  | P  |
| 7439-95-4 | Magnesium | 10300         |   | *    | P  |
| 7439-96-5 | Manganese | 3410          |   | *N J | P  |
| 7439-97-6 | Mercury   | 0.58          |   |      | AV |
| 7440-02-0 | Nickel    | 88.3          |   |      | P  |
| 7440-09-7 | Potassium | 470           | B |      | P  |
| 7782-49-2 | Selenium  | 6.1           |   |      | P  |
| 7440-22-4 | Silver    | 0.53          | U |      | P  |
| 7440-23-5 | Sodium    | 409           | B |      | P  |
| 7440-28-0 | Thallium  | 7.5           |   |      | P  |
| 7440-62-2 | Vanadium  | 33.5          |   |      | P  |
| 7440-66-6 | Zinc      | 3430          |   |      | P  |
| 57-12-5   | Cyanide   | 0.63          | U |      | C  |

Color Before: brown

Clarity Before:

Texture: course

Color After: yellow

Clarity After: clear

Artifacts: rocks

Comments:

## NYSDEC ASP

1  
INORGANIC ANALYSIS DATA SHEET

NYSDEC SAMPLE NO.

SB-2(19-23)

Lab Name: Galson Laboratories

Contract: BBL

Lab Code: 11626

Case No.:

SAS No.:

SDG No.: L62005

Matrix (soil/water): Soil

Lab Sample ID: L62005-3

Level (low/med): LOW

Date Received: 07/19/00

% Solids: 51.1

Concentration Units (ug/L or mg/kg dry weight): mg/kg

| CAS No.   | Analyte   | Concentration | C | Q    | M  |
|-----------|-----------|---------------|---|------|----|
| 7429-90-5 | Aluminum  | 8460          |   |      | P  |
| 7440-36-0 | Antimony  | 2.5           | B | N J  | P  |
| 7440-38-2 | Arsenic   | 34.7          |   |      | P  |
| 7440-39-3 | Barium    | 91.8          |   |      | P  |
| 7440-41-7 | Beryllium | 0.39          | U |      | P  |
| 7440-43-9 | Cadmium   | 0.79          | B |      | P  |
| 7440-70-2 | Calcium   | 82700         |   | *    | P  |
| 7440-47-3 | Chromium  | 14.7          |   |      | P  |
| 7440-48-4 | Cobalt    | 8.9           | B |      | P  |
| 7440-50-8 | Copper    | 53.4          |   |      | P  |
| 7439-89-6 | Iron      | 27900         |   |      | P  |
| 7439-92-1 | Lead      | 67.2          |   | N J  | P  |
| 7439-95-4 | Magnesium | 24700         |   | *    | P  |
| 7439-96-5 | Manganese | 1150          |   | *N J | P  |
| 7439-97-6 | Mercury   | 0.26          |   |      | AV |
| 7440-02-0 | Nickel    | 18.6          |   |      | P  |
| 7440-09-7 | Potassium | 1050          | B |      | P  |
| 7782-49-2 | Selenium  | 3.7           |   |      | P  |
| 7440-22-4 | Silver    | 0.78          | U |      | P  |
| 7440-23-5 | Sodium    | 170           | B |      | P  |
| 7440-28-0 | Thallium  | 2.3           | U |      | P  |
| 7440-62-2 | Vanadium  | 12.2          | B |      | P  |
| 7440-66-6 | Zinc      | 78.0          |   |      | P  |
| 57-12-5   | Cyanide   | 14.9          |   |      | C  |

Color Before: black

Clarity Before:

Texture: med

Color After: yellow

Clarity After: clear

Artifacts: rocks

Comments:

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## NYSDEC ASP

1  
INORGANIC ANALYSIS DATA SHEET

NYSDEC SAMPLE NO.

SB-3(12-14)

Lab Name: Galson Laboratories

Contract: BBL

Lab Code: 11626

Case No.:

SAS No.:

SDG No.: L62005

Matrix (soil/water): Soil

Lab Sample ID: L62005-4

Level (low/med): LOW

Date Received: 07/19/00

% Solids: 83.3

Concentration Units (ug/L or mg/kg dry weight): mg/kg

| CAS No.   | Analyte   | Concentration | C | Q    | M  |
|-----------|-----------|---------------|---|------|----|
| 7429-90-5 | Aluminum  | 5320          | - |      | P  |
| 7440-36-0 | Antimony  | 1.2           | U | N J  | P  |
| 7440-38-2 | Arsenic   | 3.3           |   |      | P  |
| 7440-39-3 | Barium    | 24.5          | B |      | P  |
| 7440-41-7 | Beryllium | 0.29          | B | u    | P  |
| 7440-43-9 | Cadmium   | 0.48          | U |      | P  |
| 7440-70-2 | Calcium   | 28600         |   | *    | P  |
| 7440-47-3 | Chromium  | 7.5           |   |      | P  |
| 7440-48-4 | Cobalt    | 5.0           | B |      | P  |
| 7440-50-8 | Copper    | 23.3          |   |      | P  |
| 7439-89-6 | Iron      | 11000         |   |      | P  |
| 7439-92-1 | Lead      | 10.3          |   | N J  | P  |
| 7439-95-4 | Magnesium | 11300         |   | *    | P  |
| 7439-96-5 | Manganese | 219           |   | *N J | P  |
| 7439-97-6 | Mercury   | 0.068         | B |      | AV |
| 7440-02-0 | Nickel    | 10            |   |      | P  |
| 7440-09-7 | Potassium | 1060          | B |      | P  |
| 7782-49-2 | Selenium  | 2.2           |   |      | P  |
| 7440-22-4 | Silver    | 0.48          | U |      | P  |
| 7440-23-5 | Sodium    | 99.9          | U |      | P  |
| 7440-28-0 | Thallium  | 1.4           | U |      | P  |
| 7440-62-2 | Vanadium  | 9.7           | B |      | P  |
| 7440-66-6 | Zinc      | 23.9          |   |      | P  |
| 57-12-5   | Cyanide   | 22.4          |   |      | C  |

Color Before: brown

Clarity Before:

Texture: med

Color After: yellow

Clarity After: clear

Artifacts: rocks

Comments:

## NYSDEC ASP

1  
INORGANIC ANALYSIS DATA SHEET

NYSDEC SAMPLE NO.

SB-4A(13-15)

Lab Name: Galson Laboratories

Contract: BBL

Lab Code: 11626

Case No.:

SAS No.:

SDG No.: L62005

Matrix (soil/water): Soil

Lab Sample ID: L62041-1

Level (low/med): LOW

Date Received: 07/20/00

% Solids: 83.0

Concentration Units (ug/L or mg/kg dry weight): mg/kg

| CAS No.   | Analyte   | Concentration | C | Q    | M  |
|-----------|-----------|---------------|---|------|----|
| 7429-90-5 | Aluminum  | 5090          |   |      | P  |
| 7440-36-0 | Antimony  | 1.2           | U | N 45 | P  |
| 7440-38-2 | Arsenic   | 2.0           | B |      | P  |
| 7440-39-3 | Barium    | 27.1          | B |      | P  |
| 7440-41-7 | Beryllium | 0.30          | B | U    | P  |
| 7440-43-9 | Cadmium   | 0.48          | U |      | P  |
| 7440-70-2 | Calcium   | 30100         |   | *    | P  |
| 7440-47-3 | Chromium  | 7.9           |   |      | P  |
| 7440-48-4 | Cobalt    | 4.7           | B |      | P  |
| 7440-50-8 | Copper    | 15.8          |   |      | P  |
| 7439-89-6 | Iron      | 11200         |   |      | P  |
| 7439-92-1 | Lead      | 6.3           |   | N J  | P  |
| 7439-95-4 | Magnesium | 10600         |   | *    | P  |
| 7439-96-5 | Manganese | 287           |   | *N J | P  |
| 7439-97-6 | Mercury   | 0.060         | U |      | AV |
| 7440-02-0 | Nickel    | 9.1           | B |      | P  |
| 7440-09-7 | Potassium | 898           | B |      | P  |
| 7782-49-2 | Selenium  | 0.96          | U |      | P  |
| 7440-22-4 | Silver    | 0.48          | U |      | P  |
| 7440-23-5 | Sodium    | 100           | U |      | P  |
| 7440-28-0 | Thallium  | 1.4           | U |      | P  |
| 7440-62-2 | Vanadium  | 9.2           | B |      | P  |
| 7440-66-6 | Zinc      | 20.0          |   |      | P  |
| 57-12-5   | Cyanide   | 0.51          | U |      | C  |

Color Before: brown

Clarity Before:

Texture: med

Color After: yellow

Clarity After: clear

Artifacts: rocks

Comments:

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## NYSDEC ASP

1  
INORGANIC ANALYSIS DATA SHEET

NYSDEC SAMPLE NO.

DUP-1

Lab Name: Galson Laboratories

Contract: BBL

Lab Code: 11626

Case No.:

SAS No.:

SDG No.: L62005

Matrix (soil/water): Soil

Lab Sample ID: L62005-5

Level (low/med): LOW

Date Received: 07/19/00

% Solids: 73.6

Concentration Units (ug/L or mg/kg dry weight): mg/kg

| CAS No.   | Analyte   | Concentration | C   | Q    | M  |
|-----------|-----------|---------------|-----|------|----|
| 7429-90-5 | Aluminum  | 4470          |     |      | P  |
| 7440-36-0 | Antimony  | 70.5          |     | N J  | P  |
| 7440-38-2 | Arsenic   | 47.3          |     |      | P  |
| 7440-39-3 | Barium    | 469           |     |      | P  |
| 7440-41-7 | Beryllium | 0.54          | B U |      | P  |
| 7440-43-9 | Cadmium   | 2.6           |     |      | P  |
| 7440-70-2 | Calcium   | 17300         |     | *    | P  |
| 7440-47-3 | Chromium  | 19.0          |     |      | P  |
| 7440-48-4 | Cobalt    | 10.3          | B   |      | P  |
| 7440-50-8 | Copper    | 1020          |     |      | P  |
| 7439-89-6 | Iron      | 79000         |     |      | P  |
| 7439-92-1 | Lead      | 14600         |     | N J  | P  |
| 7439-95-4 | Magnesium | 3480          |     | *    | P  |
| 7439-96-5 | Manganese | 3220          |     | *N J | P  |
| 7439-97-6 | Mercury   | 0.55          |     |      | AV |
| 7440-02-0 | Nickel    | 37.4          |     |      | P  |
| 7440-09-7 | Potassium | 596           | B   |      | P  |
| 7782-49-2 | Selenium  | 1.1           | U   |      | P  |
| 7440-22-4 | Silver    | 0.54          | U   |      | P  |
| 7440-23-5 | Sodium    | 226           | B   |      | P  |
| 7440-28-0 | Thallium  | 1.6           | U   |      | P  |
| 7440-62-2 | Vanadium  | 24.9          |     |      | P  |
| 7440-66-6 | Zinc      | 1450          |     |      | P  |
| 57-12-5   | Cyanide   | 114           |     |      | C  |

Color Before: brown

Clarity Before:

Texture: course

Color After: yellow

Clarity After: clear

Artifacts: rocks

Comments:

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## NYSDEC ASP

1  
INORGANIC ANALYSIS DATA SHEET

NYSDEC SAMPLE NO.

RINSEBLANK

Lab Name: Galson Laboratories

Contract: BBL

Lab Code: 11626

Case No.:

SAS No.:

SDG No.: L62005

Matrix (soil/water): Water

Lab Sample ID: L62041-5

Level (low/med): LOW

Date Received: 07/20/00

% Solids:

Concentration Units (ug/L or mg/kg dry weight): ug/l

| CAS No.   | Analyte   | Concentration | C | Q             | M  |
|-----------|-----------|---------------|---|---------------|----|
| 7429-90-5 | Aluminum  | 87.0          | B | NJ            | P  |
| 7440-36-0 | Antimony  | 5.0           | U | <del>NJ</del> | P  |
| 7440-38-2 | Arsenic   | 4.0           | U | N             | P  |
| 7440-39-3 | Barium    | 1.1           | B |               | P  |
| 7440-41-7 | Beryllium | 1.3           | B |               | P  |
| 7440-43-9 | Cadmium   | 2.0           | U | N             | P  |
| 7440-70-2 | Calcium   | 23.0          | B |               | P  |
| 7440-47-3 | Chromium  | 3.0           | U |               | P  |
| 7440-48-4 | Cobalt    | 2.0           | U |               | P  |
| 7440-50-8 | Copper    | 3.0           | U |               | P  |
| 7439-89-6 | Iron      | 78.0          | B | NJ            | P  |
| 7439-92-1 | Lead      | 2.0           | U |               | P  |
| 7439-95-4 | Magnesium | 32.6          | B |               | P  |
| 7439-96-5 | Manganese | 2.0           | U | <del>NJ</del> | P  |
| 7439-97-6 | Mercury   | 0.10          | U |               | AV |
| 7440-02-0 | Nickel    | 3.0           | U |               | P  |
| 7440-09-7 | Potassium | 843           | B |               | P  |
| 7782-49-2 | Selenium  | 4.0           | U |               | P  |
| 7440-22-4 | Silver    | 2.0           | U |               | P  |
| 7440-23-5 | Sodium    | 416           | U |               | P  |
| 7440-28-0 | Thallium  | 6.0           | U |               | P  |
| 7440-62-2 | Vanadium  | 2.0           | U | N             | P  |
| 7440-66-6 | Zinc      | 2.0           | U | N             | P  |
| 57-12-5   | Cyanide   | 10.0          | U |               | C  |

Color Before: colorless

Clarity Before: clear

Texture:

Color After: colorless

Clarity After: clear

Artifacts:

Comments:

BLASLAND, BOUCK & LEE, INC.

LABORATORY DATA REVIEW REPORT

Project: NMPC - Cedar Street  
Analytical Laboratory: Galson Laboratories  
Laboratory Report Identification Number: L62574  
Date of Laboratory Report: September 12, 2000  
Date of Review: November 4, 2000  
Reviewer: Linda Waters  
Number of Samples: 14  
Sample Matrix: soil  
Date of Collection: 8/8/00 - 8/11/00

Sample Analysis: Volatiles

Quality Control Checks

|  |            |    |                       |
|--|------------|----|-----------------------|
| 1. Field Chain-of-Custody complete   | <u>yes</u> | no | not applicable        |
| 2. Proper methods for analysis used  | <u>yes</u> | no | not applicable        |
| 3. All documentation supplied  | <u>yes</u> | no | not applicable        |
| 4. Samples analyzed within specified holding times                             | <u>yes</u> | no | not applicable        |
| 5. The minimum number of field and laboratory QC samples analyzed              | <u>yes</u> | no | not applicable        |
| 6. Laboratory accuracy maintained within established ranges for the following: |            |    |                       |
| - %RSD, initial calibration  | <u>yes</u> | no | not applicable        |
| - %D, continuing calibration   | <u>yes</u> | no | not applicable        |
| - %Recovery, matrix spike  | <u>yes</u> | no | not applicable        |
| - %Recovery, blank spike   | <u>yes</u> | no | not applicable        |
| - %Recovery, surrogate   | <u>yes</u> | no | not applicable        |
| - %Recovery, control sample  | yes        | no | <u>not applicable</u> |

7. Laboratory precision maintained within established ranges for the following:
- |                     |            |    |                |
|---------------------|------------|----|----------------|
| - RPD, matrix spike | <u>yes</u> | no | not applicable |
| - RPD, duplicates   | <u>yes</u> | no | not applicable |
8. Target analyte concentrations below detection limit in all blank samples
- |  |            |    |                |
|--|------------|----|----------------|
|  | <u>yes</u> | no | not applicable |
|--|------------|----|----------------|

Notes: Methylene chloride was detected in three method blanks. Data for methylene chloride have been qualified as undetected in samples SB-6(16-18), SB-6(18-20), DUP-2, SB-13(4-6) and SB-15(4-6) based on the blank content.

The %D for chloromethane and vinyl chloride were outside acceptable limits in one continuing calibration standard. Data for chloromethane and vinyl chloride have been qualified as estimated in the associated sample SB-8(6.5-8) based on the %D.

The response for one or more internal standards was below established limits in samples SB-14(5-7) and SB-14(5-7)RE. Data for all compounds quantitated under the non-compliant standards have been qualified as estimated based on the deviations.

Other than for the deviations noted in this review, all data quality parameters were within method-specified limits and the data is acceptable for use as reported by the laboratory.

Report Number: L62574  
 Sample Analysis: Semivolatiles

#### Quality Control Checks

- |   |            |           |                       |
|---|------------|-----------|-----------------------|
| 1. Field Chain-of-Custody complete  | <u>yes</u> | no        | not applicable        |
| 2. Proper methods for analysis used   | <u>yes</u> | no        | not applicable        |
| 3. All documentation supplied   | <u>yes</u> | no        | not applicable        |
| 4. Samples analyzed within specified holding times                              | <u>yes</u> | no        | not applicable        |
| 5. The minimum number of field and laboratory QC samples analyzed               | <u>yes</u> | no        | not applicable        |
| 6. Laboratory accuracy maintained within established ranges for the following:  |            |           |                       |
| - %RSD, initial calibration   | <u>yes</u> | no        | not applicable        |
| - %D, continuing calibration  | <u>yes</u> | no        | not applicable        |
| - %Recovery, matrix spike   | <u>yes</u> | no        | not applicable        |
| - %Recovery, blank spike  | yes        | <u>no</u> | not applicable        |
| - %Recovery, surrogate  | yes        | <u>no</u> | not applicable        |
| - %Recovery, control sample   | yes        | no        | <u>not applicable</u> |
| 7. Laboratory precision maintained within established ranges for the following: |            |           |                       |
| - RPD, matrix spike   | <u>yes</u> | no        | not applicable        |
| - RPD, duplicates   | <u>yes</u> | no        | not applicable        |
| 8. Target analyte concentrations below detection limit in all blank samples     | <u>yes</u> | no        | not applicable        |

Notes: Although no target compounds were detected, several non-target compounds were detected in the method blank. When common to the blank and samples, their presence in the samples have been rejected.

Recovery for one surrogate was above control limits in samples SB-8(6.5-8), SB-8(6.5-8)DL, SB-15(4-6), SB-15(4-6)RE, SB-10(4-7), SB-10(4-7)DL, SB-13(4-6), SB-13(4-6)DL, SB-14(5-7), SB-14(5-7)DL and DUP-2. Since recoveries for the remaining surrogates were within control limits, no data have been qualified based on the deviations.

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The non-target compounds benzo(e)pyrene was detected but not reported in the majority of the samples. The non-target data sheets have been corrected to reflect the presence.

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The response for one or more internal standards was below established limits in samples SB-8(6.5-8), SB-8(6.5-8)DL, DUP-2, SB-15(4-6), SB-15(4-6)RE, SB-10(4-7), SB-10(4-7)DL, SB-13(4-6), SB13(4-6)DL, SB-14(5-7) and SB-14(5-7)DL. Data for all compounds quantitated under the non-compliant standards have been qualified as estimated based on the deviations.

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Recovery for several compounds was below control limits in the matrix spike blank, however since all matrix spike and matrix spike duplicate recoveries were control limits, no data have been qualified based on the recoveries.

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Other than for the deviations noted in this review, the data quality parameters were within method specifications and the data is acceptable for use as reported by the laboratory.

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Report Number: L62005  
 Sample Analysis: Pesticides/PCB

#### Quality Control Checks

- |   |            |           |                       |
|---|------------|-----------|-----------------------|
| 1. Field Chain-of-Custody complete  | <u>yes</u> | no        | not applicable        |
| 2. Proper methods for analysis used   | <u>yes</u> | no        | not applicable        |
| 3. All documentation supplied   | <u>yes</u> | no        | not applicable        |
| 4. Samples analyzed within specified holding times                              | <u>yes</u> | no        | not applicable        |
| 5. The minimum number of field and laboratory QC samples analyzed               | <u>yes</u> | no        | not applicable        |
| 6. Laboratory accuracy maintained within established ranges for the following:  |            |           |                       |
| - %RSD, initial calibration   | <u>yes</u> | no        | not applicable        |
| - %D, continuing calibration  | <u>yes</u> | no        | not applicable        |
| - %Recovery, matrix spike   | <u>yes</u> | no        | not applicable        |
| - %Recovery, blank spike  | <u>yes</u> | no        | not applicable        |
| - %Recovery, surrogate  | yes        | <u>no</u> | not applicable        |
| - %Recovery, control sample   | yes        | no        | <u>not applicable</u> |
| 7. Laboratory precision maintained within established ranges for the following: |            |           |                       |
| - RPD, matrix spike   | <u>yes</u> | no        | not applicable        |
| - RPD, duplicates   | <u>yes</u> | no        | not applicable        |
| 8. Target analyte concentrations below detection limit in all blank samples     | <u>yes</u> | no        | not applicable        |

Notes: Recovery for one surrogates was above control limits in several samples due to interfering peaks. No data have been qualified based on biased recoveries due to matrix interference.

Data for several compounds were rejected due to matrix interference. In all cases the interference made the determination of presence of absence of the rejected compound impossible.

Other than for the deviations noted in this review, the data quality parameters

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were within method specifications and the data is acceptable for use as reported  
by the laboratory.

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Report Number: L62005  
 Sample Analysis: Inorganics

## Quality Control Checks

- |   |            |           |                |
|---|------------|-----------|----------------|
| 1. Field Chain-of-Custody complete  | <u>yes</u> | no        | not applicable |
| 2. Proper methods for analysis used   | <u>yes</u> | no        | not applicable |
| 3. All documentation supplied   | <u>yes</u> | no        | not applicable |
| 4. Samples analyzed within specified holding times                              | <u>yes</u> | no        | not applicable |
| 5. The minimum number of field and laboratory QC samples analyzed               | <u>yes</u> | no        | not applicable |
| 6. Laboratory accuracy maintained within established ranges for the following:  |            |           |                |
| - r2, initial calibration   | <u>yes</u> | no        | not applicable |
| - %R, continuing calibration  | <u>yes</u> | no        | not applicable |
| - %Recovery, matrix spike   | yes        | <u>no</u> | not applicable |
| - %Recovery, control sample   | <u>yes</u> | no        | not applicable |
| 7. Laboratory precision maintained within established ranges for the following: |            |           |                |
| - RPD, matrix spike   | yes        | <u>no</u> | not applicable |
| - RPD, duplicates   | yes        | <u>no</u> | not applicable |
| 8. Target analyte concentrations below detection limit in all blank samples     | <u>yes</u> | no        | not applicable |

Notes: \_\_\_\_\_  
Recoveries for antimony and manganese were outside control limits in the  
matrix spikes. Data for antimony have been qualified as estimated in samples  
samples SB-10(4-7), SB-13(4-6), SB-14(5-7), SB-15(4-6), SB-8(6.5-8), SB-6(16-18),  
SB-6(18-20) and DUP-2 and data for manganese have been qualified as estimated  
in samples SB-10(4-7), SB-13(4-6), SB-14(5-7), SB-15(4-6), SB-8(6.5-8) and  
DUP-2 based on the recoveries.

\_\_\_\_\_  
Data for zinc was incorrectly calculated for sample SB-13(4-6). Data for the  
samples was manually recalculated and data corrected on the sample analysis  
data sheet.

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The laboratory duplicate results are unacceptable for zinc. Data for zinc have been qualified as estimated in samples SB-10(4-7), SB-13(4-6), SB-14(5-7), SB-15(4-6), SB-8(6.5-8) and DUP-2 based on the duplicate results.

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Other than for the deviations noted in this review, all quality control parameters were within method specifications and the data is acceptable for use as reported by the laboratory.

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Reviewed and Approved:



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Quality Assurance Manager

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Project Manager

## SAMPLE COMPLIANCE REPORT

NMPC - Cedar Street

| Sample Delivery Group | Sampling Date | ASP Protocol | Sample ID   | Matrix | Compliance <sup>1</sup> |     |      |     |     | Noncompliance   |
|-----------------------|---------------|--------------|-------------|--------|-------------------------|-----|------|-----|-----|---|
|                       |               |              |             |        | VOA                     | BNA | Pest | PCB | TAL |   |
| L62574                | 8/8/00        | 1995         | SB-6(16-18) | soil   | no                      | no  | yes  | --  | no  | VOA - blank <sup>3</sup><br>BN - msb <sup>2</sup><br>TAL - ms   |
| L62574                | 8/8/00        | 1995         | SB-6(18-20) | soil   | no                      | no  | yes  | --  | no  | VOA - blank <sup>3</sup><br>BN -msb <sup>2</sup><br>TAL - ms  |
| L62574                | 8/10/00       | 1995         | SB-8(6.5-8) | soil   | yes                     | no  | yes  | --  | no  | BN - msb <sup>2</sup> , int std, surr <sup>2</sup><br>TAL - ms, dup   |
| L62574                | 8/10/00       | 1995         | SB-10(4-7)  | soil   | yes                     | no  | no   | --  | no  | BN - msb <sup>2</sup> , int std, surr <sup>2</sup><br>Pest - surr, id<br>TAL - ms, dup                                      |
| L62574                | 8/10/00       | 1995         | SB-13(4-6)  | soil   | no                      | no  | no   | --  | no  | VOA - blank <sup>3</sup><br>BN - msb <sup>2</sup> , int std, surr <sup>2</sup><br>Pest - surr <sup>2</sup><br>TAL - ms, dup |
| L62574                | 8/10/00       | 1995         | SB-15(4-6)  | soil   | no                      | no  | yes  | --  | no  | VOA - blank <sup>3</sup><br>BN - msb <sup>2</sup> , int std, surr <sup>2</sup><br>TAL - ms, dup                             |
| L62574                | 8/10/00       | 1995         | SB-14(5-7)  | soil   | no                      | no  | yes  | --  | no  | VOA - int std<br>BN - msb <sup>2</sup> , int std, surr <sup>2</sup><br>Pest - id<br>TAL - ms, dup                           |
| L62574                | 8/10/00       | 1995         | DUP-2       | soil   | no                      | no  | yes  | --  | no  | VOA - blank <sup>3</sup><br>BN - msb <sup>2</sup> , surr <sup>2</sup><br>Pest - id<br>TAL - ms, dup                         |
| L62574                | 8/11/00       | 1995         | SS-101      | soil   | --                      | --  | --   | yes | --  |   |
| L62574                | 8/11/00       | 1995         | SS-102      | soil   | --                      | --  | --   | no  | --  | PCB - surr <sup>2</sup>   |
| L62574                | 8/11/00       | 1995         | SS-103      | soil   | --                      | --  | --   | yes | --  |   |

| Sample Delivery Group | Sampling Date | ASP Protocol | Sample ID | Matrix | Compliance <sup>1</sup> |     |      |     |     | Noncompliance           |
|-----------------------|---------------|--------------|-----------|--------|-------------------------|-----|------|-----|-----|-------------------------|
|                       |               |              |           |        | VOA                     | BNA | Pest | PCB | TAL |                         |
| L62574                | 8/11/00       | 1995         | SS-104    | soil   | --                      | --  | --   | yes | --  |                         |
| L62574                | 8/11/00       | 1995         | SS-DUP    | soil   | --                      | --  | --   | no  | --  | PCB - surr <sup>2</sup> |
| L62574                | --            | 1995         | RB-PCB    | water  | --                      | --  | --   | yes | --  |                         |
|                       |               |              |           |        |                         |     |      |     |     |                         |

- 1 Samples which are compliant with no added validation qualifiers are listed as "yes". Samples which are non-compliant or which have added qualifiers are listed as "no". A "no" designation does not necessarily indicate that the data have been rejected or are otherwise unusable.
- 2 The noncompliance resulted in no qualification of data.
- 3 Although the deviation resulted in the qualification of data, the laboratory was method compliant.

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SB-6 (16-18)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) Soil

Lab Sample ID: L62574-1

Sample wt/vol: 5 (g/mL) g

Lab File ID: CD081007

Level: (low/med) LOW

Date Received: 08/08/00

%Moisture: not dec. 17

Date Analyzed: 08/10/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 1

Soil Extract Volume: (uL)

Soil Aliquot Volume: (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) ug/kg

Q

CAS NO.

COMPOUND

|                 |                            |      |     |
|-----------------|----------------------------|------|-----|
| 74-87-3-----    | Chloromethane              | 12   | U   |
| 74-83-9-----    | Bromomethane               | 12   | U   |
| 75-01-4-----    | Vinyl Chloride             | 12   | U   |
| 75-00-3-----    | Chloroethane               | 12   | U   |
| 75-09-2-----    | Methylene Chloride         | 12 2 | JBU |
| 67-64-1-----    | Acetone                    | 12   | U   |
| 75-15-0-----    | Carbon Disulfide           | 12   | U   |
| 75-35-4-----    | 1,1-Dichloroethene         | 12   | U   |
| 75-34-3-----    | 1,1-Dichloroethane         | 12   | U   |
| 540-59-0-----   | 1,2-Dichloroethene (Total) | 12   | U   |
| 67-66-3-----    | Chloroform                 | 12   | U   |
| 107-06-2-----   | 1,2-Dichloroethane         | 12   | U   |
| 78-93-3-----    | 2-Butanone                 | 12   | U   |
| 71-55-6-----    | 1,1,1-Trichloroethane      | 12   | U   |
| 56-23-5-----    | Carbon Tetrachloride       | 12   | U   |
| 75-27-4-----    | Bromodichloromethane       | 12   | U   |
| 78-87-5-----    | 1,2-Dichloropropane        | 12   | U   |
| 10061-01-5----- | cis-1,3-Dichloropropene    | 12   | U   |
| 79-01-6-----    | Trichloroethene            | 12   | U   |
| 124-48-1-----   | Dibromochloromethane       | 12   | U   |
| 79-00-5-----    | 1,1,2-Trichloroethane      | 12   | U   |
| 71-43-2-----    | Benzene                    | 12   | U   |
| 10061-02-6----- | trans-1,3-Dichloropropene  | 12   | U   |
| 75-25-2-----    | Bromoform                  | 12   | U   |
| 108-10-1-----   | 4-Methyl-2-Pentanone       | 12   | U   |
| 591-78-6-----   | 2-Hexanone                 | 12   | U   |
| 127-18-4-----   | Tetrachloroethene          | 12   | U   |
| 79-34-5-----    | 1,1,2,2-Tetrachloroethane  | 12   | U   |
| 108-88-3-----   | Toluene                    | 12   | U   |
| 108-90-7-----   | Chlorobenzene              | 12   | U   |
| 100-41-4-----   | Ethylbenzene               | 12   | U   |
| 100-42-5-----   | Styrene                    | 12   | U   |
| 1330-20-7-----  | Xylene (total)             | 12   | U   |

1E  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SB-6 (16-18)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) Soil

Lab Sample ID: L62574-1

Sample wt/vol: 5 (g/mL) g

Lab File ID: CD081007

Level: (low/med) LOW

Date Received: 08/08/00

%Moisture: not dec. 17

Date Analyzed: 08/10/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 1

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.         | No Volatiles Found |    |            |   |
| 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

SAMPLE NO.

SB-6 (18-20)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) Soil

Lab Sample ID: L62574-4

Sample wt/vol: 5 (g/mL) g

Lab File ID: CD081005

Level: (low/med) LOW

Date Received: 08/08/00

%Moisture: not dec. 18

Date Analyzed: 08/10/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 1

Soil Extract Volume: (uL)

Soil Aliquot Volume: (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) ug/kg Q

CAS NO.

COMPOUND

|                 |                            |      |     |
|-----------------|----------------------------|------|-----|
| 74-87-3-----    | Chloromethane              | 12   | U   |
| 74-83-9-----    | Bromomethane               | 12   | U   |
| 75-01-4-----    | Vinyl Chloride             | 12   | U   |
| 75-00-3-----    | Chloroethane               | 12   | U   |
| 75-09-2-----    | Methylene Chloride         | 12 2 | JBW |
| 67-64-1-----    | Acetone                    | 12   | U   |
| 75-15-0-----    | Carbon Disulfide           | 12   | U   |
| 75-35-4-----    | 1,1-Dichloroethene         | 12   | U   |
| 75-34-3-----    | 1,1-Dichloroethane         | 12   | U   |
| 540-59-0-----   | 1,2-Dichloroethene (Total) | 12   | U   |
| 67-66-3-----    | Chloroform                 | 12   | U   |
| 107-06-2-----   | 1,2-Dichloroethane         | 12   | U   |
| 78-93-3-----    | 2-Butanone                 | 12   | U   |
| 71-55-6-----    | 1,1,1-Trichloroethane      | 12   | U   |
| 56-23-5-----    | Carbon Tetrachloride       | 12   | U   |
| 75-27-4-----    | Bromodichloromethane       | 12   | U   |
| 78-87-5-----    | 1,2-Dichloropropane        | 12   | U   |
| 10061-01-5----- | cis-1,3-Dichloropropene    | 12   | U   |
| 79-01-6-----    | Trichloroethene            | 12   | U   |
| 124-48-1-----   | Dibromochloromethane       | 12   | U   |
| 79-00-5-----    | 1,1,2-Trichloroethane      | 12   | U   |
| 71-43-2-----    | Benzene                    | 12   | U   |
| 10061-02-6----- | trans-1,3-Dichloropropene  | 12   | U   |
| 75-25-2-----    | Bromoform                  | 12   | U   |
| 108-10-1-----   | 4-Methyl-2-Pentanone       | 12   | U   |
| 591-78-6-----   | 2-Hexanone                 | 12   | U   |
| 127-18-4-----   | Tetrachloroethene          | 12   | U   |
| 79-34-5-----    | 1,1,2,2-Tetrachloroethane  | 12   | U   |
| 108-88-3-----   | Toluene                    | 12   | U   |
| 108-90-7-----   | Chlorobenzene              | 12   | U   |
| 100-41-4-----   | Ethylbenzene               | 12   | U   |
| 100-42-5-----   | Styrene                    | 12   | U   |
| 1330-20-7-----  | Xylene (total)             | 12   | U   |

1E  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SB-6 (18-20)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) Soil

Lab Sample ID: L62574-4

Sample wt/vol: 5 (g/mL) g

Lab File ID: CD081005

Level: (low/med) LOW

Date Received: 08/08/00

%Moisture: not dec. 18

Date Analyzed: 08/10/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 1

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. _____   | No Volatiles Found _____ |       |            |       |
| 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

SAMPLE NO.

SB-8 (6.5-8)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) Soil

Lab Sample ID: L62615-1

Sample wt/vol: 2 (g/mL) g

Lab File ID: CD081106

Level: (low/med) LOW

Date Received: 08/10/00

%Moisture: not dec. 75

Date Analyzed: 08/11/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 2.5

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   |
|------------|----------------------------|---|-----|
| 74-87-3    | Chloromethane              | 100   | U   |
| 74-83-9    | Bromomethane               | 100   | U   |
| 75-01-4    | Vinyl Chloride             | 100   | U   |
| 75-00-3    | Chloroethane               | 100   | U   |
| 75-09-2    | Methylene Chloride         | 100   | U   |
| 67-64-1    | Acetone                    | 580   |     |
| 75-15-0    | Carbon Disulfide           | 100   | U   |
| 75-35-4    | 1,1-Dichloroethene         | 100   | U   |
| 75-34-3    | 1,1-Dichloroethane         | 100   | U   |
| 540-59-0   | 1,2-Dichloroethene (Total) | 100   | U   |
| 67-66-3    | Chloroform                 | 100   | U   |
| 107-06-2   | 1,2-Dichloroethane         | 100   | U   |
| 78-93-3    | 2-Butanone                 | 250   |     |
| 71-55-6    | 1,1,1-Trichloroethane      | 100   | U   |
| 56-23-5    | Carbon Tetrachloride       | 100   | U   |
| 75-27-4    | Bromodichloromethane       | 100   | U   |
| 78-87-5    | 1,2-Dichloropropane        | 100   | U   |
| 10061-01-5 | cis-1,3-Dichloropropene    | 100   | U   |
| 79-01-6    | Trichloroethene            | 100   | U   |
| 124-48-1   | Dibromochloromethane       | 100   | U   |
| 79-00-5    | 1,1,2-Trichloroethane      | 100   | U   |
| 71-43-2    | Benzene                    | 7600 8500                                     | E D |
| 10061-02-6 | trans-1,3-Dichloropropene  | 100   | U   |
| 75-25-2    | Bromoform                  | 100   | U   |
| 108-10-1   | 4-Methyl-2-Pentanone       | 100   | U   |
| 591-78-6   | 2-Hexanone                 | 100   | U   |
| 127-18-4   | Tetrachloroethene          | 100   | U   |
| 79-34-5    | 1,1,2,2-Tetrachloroethane  | 100   | U   |
| 108-88-3   | Toluene                    | 300   |     |
| 108-90-7   | Chlorobenzene              | 100   | U   |
| 100-41-4   | Ethylbenzene               | 5600 9200                                     | E D |
| 100-42-5   | Styrene                    | 100   | U   |
| 1330-20-7  | Xylene (total)             | 9800 17000                                    | E D |

1E  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SB-8 (6.5-8)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) Soil

Lab Sample ID: L62615-1

Sample wt/vol: 2 (g/mL) g

Lab File ID: CD081106

Level: (low/med) LOW

Date Received: 08/10/00

%Moisture: not dec. 75

Date Analyzed: 08/11/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 2.5

Soil Extract Volume: (uL)

Soil Aliquot Volume: (uL)

Number TICS found: 10

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg

| CAS NUMBER      | COMPOUND NAME          | RT    | EST. CONC. | Q |
|-----------------|------------------------|-------|------------|---|
| 1.              | C3-Substituted Benzene | 27.89 | 1300       | J |
| 2.              | Benzene, trimethyl-    | 28.08 | 2000       | J |
| 3.              | Benzene, trimethyl-    | 29.08 | 4300       | J |
| 4.              | Benzene, trimethyl-    | 30.22 | 1800       | J |
| 5. 000496-11-7  | Indane                 | 30.83 | 22000      | J |
| 6. 000095-13-6  | Indene                 | 31.41 | 11000      | J |
| 7.              | Unknown aromatic       | 32.92 | 1300       | J |
| 8.              | Unknown aromatic       | 33.22 | 1600       | J |
| 9. 000091-20-3  | Naphthalene            | 36.57 | 14000      | J |
| 10. 000270-82-6 | Benzo[c]thiophene      | 36.99 | 1900       | J |
| 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

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

SB-8 (6.5-8) DL

Lab Code: Case No.: 2

SAS No.: SDG No.: L62574

Matrix: (soil/water) Soil

Lab Sample ID: L62615-1DL

Sample wt/vol: 4 (g/mL) g

Lab File ID: CD081506

Level: (low/med) MED

Date Received: 08/10/00

%Moisture: not dec. 75

Date Analyzed: 08/15/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 1

Soil Extract Volume: 10000 (uL)

Soil Aliquot Volume: 100 (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) ug/kg Q

CAS NO.

COMPOUND

|                 |                            |       |     |
|-----------------|----------------------------|-------|-----|
| 74-87-3-----    | Chloromethane              | 5000  | U W |
| 74-83-9-----    | Bromomethane               | 5000  | U   |
| 75-01-4-----    | Vinyl Chloride             | 5000  | U W |
| 75-00-3-----    | Chloroethane               | 5000  | U   |
| 75-09-2-----    | Methylene Chloride         | 1100  | JD  |
| 67-64-1-----    | Acetone                    | 5000  | U   |
| 75-15-0-----    | Carbon Disulfide           | 5000  | U   |
| 75-35-4-----    | 1,1-Dichloroethene         | 5000  | U   |
| 75-34-3-----    | 1,1-Dichloroethane         | 5000  | U   |
| 540-59-0-----   | 1,2-Dichloroethene (Total) | 5000  | U   |
| 67-66-3-----    | Chloroform                 | 5000  | U   |
| 107-06-2-----   | 1,2-Dichloroethane         | 5000  | U   |
| 78-93-3-----    | 2-Butanone                 | 2900  | JD  |
| 71-55-6-----    | 1,1,1-Trichloroethane      | 5000  | U   |
| 56-23-5-----    | Carbon Tetrachloride       | 5000  | U   |
| 75-27-4-----    | Bromodichloromethane       | 5000  | U   |
| 78-87-5-----    | 1,2-Dichloropropane        | 5000  | U   |
| 10061-01-5----- | cis-1,3-Dichloropropene    | 5000  | U   |
| 79-01-6-----    | Trichloroethene            | 5000  | U   |
| 124-48-1-----   | Dibromochloromethane       | 5000  | U   |
| 79-00-5-----    | 1,1,2-Trichloroethane      | 5000  | U   |
| 71-43-2-----    | Benzene                    | 8500  | D   |
| 10061-02-6----- | trans-1,3-Dichloropropene  | 5000  | U   |
| 75-25-2-----    | Bromoform                  | 5000  | U   |
| 108-10-1-----   | 4-Methyl-2-Pentanone       | 5000  | U   |
| 591-78-6-----   | 2-Hexanone                 | 5000  | U   |
| 127-18-4-----   | Tetrachloroethene          | 5000  | U   |
| 79-34-5-----    | 1,1,2,2-Tetrachloroethane  | 5000  | U   |
| 108-88-3-----   | Toluene                    | 5000  | U   |
| 108-90-7-----   | Chlorobenzene              | 5000  | U   |
| 100-41-4-----   | Ethylbenzene               | 9200  | D   |
| 100-42-5-----   | Styrene                    | 5000  | U   |
| 1330-20-7-----  | Xylene (total)             | 17000 | D   |

1E  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SB-8 (6.5-8) DL

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) Soil

Lab Sample ID: L62615-1 DL

Sample wt/vol: 4 (g/mL) g

Lab File ID: CD081506

Level: (low/med) MED

Date Received: 08/10/00

%Moisture: not dec. 75

Date Analyzed: 08/15/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 1

Soil Extract Volume: 10000 (uL)

Soil Aliquot Volume: 100 (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) ug/kg

Number TICS found: 10

| CAS NUMBER      | COMPOUND NAME                | RT    | EST. CONC. | Q |
|-----------------|------------------------------|-------|------------|---|
| 1.              | Benzene, trimethyl-          | 28.08 | 5200       | J |
| 2.              | Benzene, trimethyl-          | 29.08 | 11000      | J |
| 3.              | Benzene, trimethyl-          | 30.22 | 4600       | J |
| 4. 000496-11-7  | Indane                       | 30.83 | 67000      | J |
| 5. 000095-13-6  | Indene                       | 31.41 | 21000      | J |
| 6.              | Unknown Aromatic             | 32.93 | 4000       | J |
| 7.              | Unknown Aromatic             | 33.21 | 3900       | J |
| 8. 000874-35-1  | 1H-Indene, 2,3-dihydro-5-met | 34.47 | 4100       | J |
| 9. 000091-20-3  | Naphthalene                  | 36.57 | 71000      | J |
| 10. 000270-82-6 | Benzo[c]thiophene            | 37.00 | 5000       | J |
| 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

SAMPLE NO.

SB-10 (4-7)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) Soil

Lab Sample ID: L62615-2

Sample wt/vol: 5 (g/mL) g

Lab File ID: CD081114

Level: (low/med) LOW

Date Received: 08/10/00

%Moisture: not dec. 18

Date Analyzed: 08/11/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 1

Soil Extract Volume: (uL)

Soil Aliquot Volume: (uL)

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|                 |                            |    |   |
|-----------------|----------------------------|----|---|
| 74-87-3-----    | Chloromethane              | 12 | U |
| 74-83-9-----    | Bromomethane               | 12 | U |
| 75-01-4-----    | Vinyl Chloride             | 12 | U |
| 75-00-3-----    | Chloroethane               | 12 | U |
| 75-09-2-----    | Methylene Chloride         | 12 | U |
| 67-64-1-----    | Acetone                    | 21 |   |
| 75-15-0-----    | Carbon Disulfide           | 12 | U |
| 75-35-4-----    | 1,1-Dichloroethene         | 12 | U |
| 75-34-3-----    | 1,1-Dichloroethane         | 12 | U |
| 540-59-0-----   | 1,2-Dichloroethene (Total) | 12 | U |
| 67-66-3-----    | Chloroform                 | 12 | U |
| 107-06-2-----   | 1,2-Dichloroethane         | 12 | U |
| 78-93-3-----    | 2-Butanone                 | 12 | U |
| 71-55-6-----    | 1,1,1-Trichloroethane      | 12 | U |
| 56-23-5-----    | Carbon Tetrachloride       | 12 | U |
| 75-27-4-----    | Bromodichloromethane       | 12 | U |
| 78-87-5-----    | 1,2-Dichloropropane        | 12 | U |
| 10061-01-5----- | cis-1,3-Dichloropropene    | 12 | U |
| 79-01-6-----    | Trichloroethene            | 12 | U |
| 124-48-1-----   | Dibromochloromethane       | 12 | U |
| 79-00-5-----    | 1,1,2-Trichloroethane      | 12 | U |
| 71-43-2-----    | Benzene                    | 2  | J |
| 10061-02-6----- | trans-1,3-Dichloropropene  | 12 | U |
| 75-25-2-----    | Bromoform                  | 12 | U |
| 108-10-1-----   | 4-Methyl-2-Pentanone       | 12 | U |
| 591-78-6-----   | 2-Hexanone                 | 12 | U |
| 127-18-4-----   | Tetrachloroethene          | 12 | U |
| 79-34-5-----    | 1,1,2,2-Tetrachloroethane  | 12 | U |
| 108-88-3-----   | Toluene                    | 3  | J |
| 108-90-7-----   | Chlorobenzene              | 12 | U |
| 100-41-4-----   | Ethylbenzene               | 12 | U |
| 100-42-5-----   | Styrene                    | 1  | J |
| 1330-20-7-----  | Xylene (total)             | 4  | J |

1E  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

SB-10 (4-7)

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) Soil

Lab Sample ID: L62615-2

Sample wt/vol: 5 (g/mL) g

Lab File ID: CD081114

Level: (low/med) LOW

Date Received: 08/10/00

%Moisture: not dec. 18

Date Analyzed: 08/11/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 1

Soil Extract Volume: (uL)

Soil Aliquot Volume: (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) ug/kg

Number TICS found: 4

| CAS NUMBER     | COMPOUND NAME    | RT    | EST. CONC. | Q |
|----------------|------------------|-------|------------|---|
| 1. 000496-11-7 | Indane           | 30.81 | 8          | J |
| 2. 000095-13-6 | Indene           | 31.41 | 17         | J |
| 3. 000091-20-3 | Naphthalene      | 36.56 | 290        | J |
| 4.             | Unknown aromatic | 36.97 | 7          | 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.            |                  |       |            |   |



1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SB-13 (4-6)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) Soil

Lab Sample ID: L62615-3

Sample wt/vol: 5 (g/mL) g

Lab File ID: CD081108

Level: (low/med) LOW

Date Received: 08/10/00

%Moisture: not dec. 26

Date Analyzed: 08/11/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 1

Soil Extract Volume: (uL)

Soil Aliquot Volume: (uL)

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|                 |                            |      |     |
|-----------------|----------------------------|------|-----|
| 74-87-3-----    | Chloromethane              | 14   | U   |
| 74-83-9-----    | Bromomethane               | 14   | U   |
| 75-01-4-----    | Vinyl Chloride             | 14   | U   |
| 75-00-3-----    | Chloroethane               | 14   | U   |
| 75-09-2-----    | Methylene Chloride         | 14 2 | JBU |
| 67-64-1-----    | Acetone                    | 8    | J   |
| 75-15-0-----    | Carbon Disulfide           | 14   | U   |
| 75-35-4-----    | 1,1-Dichloroethene         | 14   | U   |
| 75-34-3-----    | 1,1-Dichloroethane         | 14   | U   |
| 540-59-0-----   | 1,2-Dichloroethene (Total) | 14   | U   |
| 67-66-3-----    | Chloroform                 | 14   | U   |
| 107-06-2-----   | 1,2-Dichloroethane         | 14   | U   |
| 78-93-3-----    | 2-Butanone                 | 14   | U   |
| 71-55-6-----    | 1,1,1-Trichloroethane      | 14   | U   |
| 56-23-5-----    | Carbon Tetrachloride       | 14   | U   |
| 75-27-4-----    | Bromodichloromethane       | 14   | U   |
| 78-87-5-----    | 1,2-Dichloropropane        | 14   | U   |
| 10061-01-5----- | cis-1,3-Dichloropropene    | 14   | U   |
| 79-01-6-----    | Trichloroethene            | 14   | U   |
| 124-48-1-----   | Dibromochloromethane       | 14   | U   |
| 79-00-5-----    | 1,1,2-Trichloroethane      | 14   | U   |
| 71-43-2-----    | Benzene                    | 14   | U   |
| 10061-02-6----- | trans-1,3-Dichloropropene  | 14   | U   |
| 75-25-2-----    | Bromoform                  | 14   | U   |
| 108-10-1-----   | 4-Methyl-2-Pentanone       | 14   | U   |
| 591-78-6-----   | 2-Hexanone                 | 14   | U   |
| 127-18-4-----   | Tetrachloroethene          | 14   | U   |
| 79-34-5-----    | 1,1,2,2-Tetrachloroethane  | 14   | U   |
| 108-88-3-----   | Toluene                    | 14   | U   |
| 108-90-7-----   | Chlorobenzene              | 14   | U   |
| 100-41-4-----   | Ethylbenzene               | 14   | U   |
| 100-42-5-----   | Styrene                    | 2    | J   |
| 1330-20-7-----  | Xylene (total)             | 14   | U   |

1E  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SB-13 (4-6)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) Soil

Lab Sample ID: L62615-3

Sample wt/vol: 5 (g/mL) g

Lab File ID: CD081108

Level: (low/med) LOW

Date Received: 08/10/00

%Moisture: not dec. 26

Date Analyzed: 08/11/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 1

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. _____   | No Volatiles Found _____ |       |            |       |
| 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

SAMPLE NO.

SB-14 (5-7)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) Soil

Lab Sample ID: L62615-5

Sample wt/vol: 5 (g/mL) g

Lab File ID: CD081110

Level: (low/med) LOW

Date Received: 08/10/00

%Moisture: not dec. 28

Date Analyzed: 08/11/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 1

Soil Extract Volume: (uL)

Soil Aliquot Volume: (uL)

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|                 |                            |    |   |
|-----------------|----------------------------|----|---|
| 74-87-3-----    | Chloromethane              | 14 | U |
| 74-83-9-----    | Bromomethane               | 14 | U |
| 75-01-4-----    | Vinyl Chloride             | 14 | U |
| 75-00-3-----    | Chloroethane               | 14 | U |
| 75-09-2-----    | Methylene Chloride         | 14 | U |
| 67-64-1-----    | Acetone                    | 67 |   |
| 75-15-0-----    | Carbon Disulfide           | 14 | U |
| 75-35-4-----    | 1,1-Dichloroethene         | 14 | U |
| 75-34-3-----    | 1,1-Dichloroethane         | 14 | U |
| 540-59-0-----   | 1,2-Dichloroethene (Total) | 14 | U |
| 67-66-3-----    | Chloroform                 | 14 | U |
| 107-06-2-----   | 1,2-Dichloroethane         | 14 | U |
| 78-93-3-----    | 2-Butanone                 | 26 |   |
| 71-55-6-----    | 1,1,1-Trichloroethane      | 14 | U |
| 56-23-5-----    | Carbon Tetrachloride       | 14 | U |
| 75-27-4-----    | Bromodichloromethane       | 14 | U |
| 78-87-5-----    | 1,2-Dichloropropane        | 14 | U |
| 10061-01-5----- | cis-1,3-Dichloropropene    | 14 | U |
| 79-01-6-----    | Trichloroethene            | 14 | U |
| 124-48-1-----   | Dibromochloromethane       | 14 | U |
| 79-00-5-----    | 1,1,2-Trichloroethane      | 14 | U |
| 71-43-2-----    | Benzene                    | 33 |   |
| 10061-02-6----- | trans-1,3-Dichloropropene  | 14 | U |
| 75-25-2-----    | Bromoform                  | 14 | U |
| 108-10-1-----   | 4-Methyl-2-Pentanone       | 14 | U |
| 591-78-6-----   | 2-Hexanone                 | 14 | U |
| 127-18-4-----   | Tetrachloroethene          | 14 | U |
| 79-34-5-----    | 1,1,2,2-Tetrachloroethane  | 14 | U |
| 108-88-3-----   | Toluene                    | 4  | U |
| 108-90-7-----   | Chlorobenzene              | 14 | U |
| 100-41-4-----   | Ethylbenzene               | 19 | U |
| 100-42-5-----   | Styrene                    | 14 | U |
| 1330-20-7-----  | Xylene (total)             | 23 | U |

1E  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SB-14 (5-7)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) Soil

Lab Sample ID: L62615-5

Sample wt/vol: 5 (g/mL) g

Lab File ID: CD081110

Level: (low/med) LOW

Date Received: 08/10/00

%Moisture: not dec. 28

Date Analyzed: 08/11/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 1

Soil Extract Volume: (uL)

Soil Aliquot Volume: (uL)

Number TICS found: 10

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg

| CAS NUMBER      | COMPOUND NAME                | RT    | EST. CONC. | Q     |
|-----------------|------------------------------|-------|------------|-------|
| 1. 000099-82-1  | 1-Methyl-4-(1-methylethyl)-c | 28.00 | 74         | J     |
| 2. _____        | Unknown                      | 28.51 | 73         | J     |
| 3. _____        | C3-Substituted Benzene       | 28.67 | 98         | J     |
| 4. _____        | Benzene, trimethyl-          | 30.21 | 160        | J     |
| 5. 000496-11-7  | Indane                       | 30.82 | 2900       | J     |
| 6. 000095-13-6  | Indene                       | 31.40 | 150        | J     |
| 7. 000767-58-8  | Indan, 1-methyl-             | 32.30 | 57         | J     |
| 8. _____        | Unknown Aromatic             | 32.93 | 87         | J     |
| 9. _____        | Unknown Aromatic             | 33.24 | 61         | J     |
| 10. 000824-22-6 | 1H-Indene, 2,3-dihydro-4-met | 34.45 | 150        | J     |
| 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

SAMPLE NO.

SB-14 (5-7) RE

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) Soil

Lab Sample ID: L62615-5RE

Sample wt/vol: 5 (g/mL) g

Lab File ID: CD081116

Level: (low/med) LOW

Date Received: 08/10/00

%Moisture: not dec. 28

Date Analyzed: 08/11/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 1

Soil Extract Volume: (uL)

Soil Aliquot Volume: (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) ug/kg Q

CAS NO.

COMPOUND

|                 |                            |     |     |
|-----------------|----------------------------|-----|-----|
| 74-87-3-----    | Chloromethane              | 14  | U   |
| 74-83-9-----    | Bromomethane               | 14  | U   |
| 75-01-4-----    | Vinyl Chloride             | 14  | U   |
| 75-00-3-----    | Chloroethane               | 14  | U   |
| 75-09-2-----    | Methylene Chloride         | 14  | U   |
| 67-64-1-----    | Acetone                    | 110 |     |
| 75-15-0-----    | Carbon Disulfide           | 14  | U   |
| 75-35-4-----    | 1,1-Dichloroethene         | 14  | U   |
| 75-34-3-----    | 1,1-Dichloroethane         | 14  | U   |
| 540-59-0-----   | 1,2-Dichloroethene (Total) | 14  | U   |
| 67-66-3-----    | Chloroform                 | 14  | U   |
| 107-06-2-----   | 1,2-Dichloroethane         | 14  | U   |
| 78-93-3-----    | 2-Butanone                 | 42  |     |
| 71-55-6-----    | 1,1,1-Trichloroethane      | 14  | U U |
| 56-23-5-----    | Carbon Tetrachloride       | 14  | U U |
| 75-27-4-----    | Bromodichloromethane       | 14  | U U |
| 78-87-5-----    | 1,2-Dichloropropane        | 14  | U U |
| 10061-01-5----- | cis-1,3-Dichloropropene    | 14  | U U |
| 79-01-6-----    | Trichloroethene            | 14  | U U |
| 124-48-1-----   | Dibromochloromethane       | 14  | U U |
| 79-00-5-----    | 1,1,2-Trichloroethane      | 14  | U U |
| 71-43-2-----    | Benzene                    | 17  | J   |
| 10061-02-6----- | trans-1,3-Dichloropropene  | 14  | U U |
| 75-25-2-----    | Bromoform                  | 14  | U U |
| 108-10-1-----   | 4-Methyl-2-Pentanone       | 14  | U U |
| 591-78-6-----   | 2-Hexanone                 | 14  | U U |
| 127-18-4-----   | Tetrachloroethene          | 14  | U U |
| 79-34-5-----    | 1,1,2,2-Tetrachloroethane  | 14  | U U |
| 108-88-3-----   | Toluene                    | 3   | J   |
| 108-90-7-----   | Chlorobenzene              | 14  | U U |
| 100-41-4-----   | Ethylbenzene               | 9   | J   |
| 100-42-5-----   | Styrene                    | 14  | U U |
| 1330-20-7-----  | Xylene (total)             | 14  |     |

1E  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

SB-14 (5-7) RE

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) Soil

Lab Sample ID: L62615-5 RE

Sample wt/vol: 5 (g/mL) g

Lab File ID: CD081116

Level: (low/med) LOW

Date Received: 08/10/00

%Moisture: not dec. 28

Date Analyzed: 08/11/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 1

Soil Extract Volume: (uL)

Soil Aliquot Volume: (uL)

Number TICS found: 10

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg

| CAS NUMBER      | COMPOUND NAME                | RT    | EST. CONC. | Q |
|-----------------|------------------------------|-------|------------|---|
| 1.              | Unknown Hydrocarbon          | 28.50 | 96         | J |
| 2.              | C3-Substituted Benzene       | 28.67 | 79         | J |
| 3.              | Benzene, trimethyl-          | 30.20 | 140        | J |
| 4. 000496-11-7  | Indane                       | 30.82 | 2600       | J |
| 5. 000095-13-6  | Indene                       | 31.40 | 99         | J |
| 6. 000767-58-8  | Indan, 1-methyl-             | 32.28 | 53         | J |
| 7.              | Unknown Aromatic             | 32.91 | 86         | J |
| 8.              | Unknown Aromatic             | 33.23 | 57         | J |
| 9. 000824-22-6  | 1H-Indene, 2,3-dihydro-4-met | 34.46 | 160        | J |
| 10. 000095-15-8 | Benzo[b]thiophene            | 36.98 | 54         | J |
| 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

SAMPLE NO.

SB-15 (4-6)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) Soil

Lab Sample ID: L62615-4

Sample wt/vol: 5 (g/mL) g

Lab File ID: CD081404

Level: (low/med) LOW

Date Received: 08/10/00

%Moisture: not dec. 29

Date Analyzed: 08/14/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 1

Soil Extract Volume: (uL)

Soil Aliquot Volume: (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) ug/kg Q

CAS NO.

COMPOUND

|                 |                            |      |     |
|-----------------|----------------------------|------|-----|
| 74-87-3-----    | Chloromethane              | 14   | U   |
| 74-83-9-----    | Bromomethane               | 14   | U   |
| 75-01-4-----    | Vinyl Chloride             | 14   | U   |
| 75-00-3-----    | Chloroethane               | 14   | U   |
| 75-09-2-----    | Methylene Chloride         | 14 2 | JBU |
| 67-64-1-----    | Acetone                    | 13   | J   |
| 75-15-0-----    | Carbon Disulfide           | 14   | U   |
| 75-35-4-----    | 1,1-Dichloroethene         | 14   | U   |
| 75-34-3-----    | 1,1-Dichloroethane         | 14   | U   |
| 540-59-0-----   | 1,2-Dichloroethene (Total) | 14   | U   |
| 67-66-3-----    | Chloroform                 | 14   | U   |
| 107-06-2-----   | 1,2-Dichloroethane         | 14   | U   |
| 78-93-3-----    | 2-Butanone                 | 14   | U   |
| 71-55-6-----    | 1,1,1-Trichloroethane      | 14   | U   |
| 56-23-5-----    | Carbon Tetrachloride       | 14   | U   |
| 75-27-4-----    | Bromodichloromethane       | 14   | U   |
| 78-87-5-----    | 1,2-Dichloropropane        | 14   | U   |
| 10061-01-5----- | cis-1,3-Dichloropropene    | 14   | U   |
| 79-01-6-----    | Trichloroethene            | 14   | U   |
| 124-48-1-----   | Dibromochloromethane       | 14   | U   |
| 79-00-5-----    | 1,1,2-Trichloroethane      | 14   | U   |
| 71-43-2-----    | Benzene                    | 14   | U   |
| 10061-02-6----- | trans-1,3-Dichloropropene  | 14   | U   |
| 75-25-2-----    | Bromoform                  | 14   | U   |
| 108-10-1-----   | 4-Methyl-2-Pentanone       | 14   | U   |
| 591-78-6-----   | 2-Hexanone                 | 14   | U   |
| 127-18-4-----   | Tetrachloroethene          | 14   | U   |
| 79-34-5-----    | 1,1,2,2-Tetrachloroethane  | 14   | U   |
| 108-88-3-----   | Toluene                    | 14   | U   |
| 108-90-7-----   | Chlorobenzene              | 14   | U   |
| 100-41-4-----   | Ethylbenzene               | 14   | U   |
| 100-42-5-----   | Styrene                    | 14   | U   |
| 1330-20-7-----  | Xylene (total)             | 14   | U   |

1E  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SB-15 (4-6)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) Soil

Lab Sample ID: L62615-4

Sample wt/vol: 5 (g/mL) g

Lab File ID: CD081404

Level: (low/med) LOW

Date Received: 08/10/00

%Moisture: not dec. 29

Date Analyzed: 08/14/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 1

Soil Extract Volume: (uL)

Soil Aliquot Volume: (uL)

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg

Number TICS found: 0

| CAS NUMBER | COMPOUND NAME      | RT | EST. CONC. | Q |
|------------|--------------------|----|------------|---|
| 1.         | No Volatiles Found |    |            |   |
| 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

SAMPLE NO.

DUP-2

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) Soil

Lab Sample ID: L62615-6

Sample wt/vol: 5 (g/mL) g

Lab File ID: CD081111

Level: (low/med) LOW

Date Received: 08/10/00

%Moisture: not dec. 21

Date Analyzed: 08/11/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 1

Soil Extract Volume: (uL)

Soil Aliquot Volume: (uL)

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|                 |                            |                 |   |
|-----------------|----------------------------|-----------------|---|
| 74-87-3-----    | Chloromethane              | 13              | U |
| 74-83-9-----    | Bromomethane               | 13              | U |
| 75-01-4-----    | Vinyl Chloride             | 13              | U |
| 75-00-3-----    | Chloroethane               | 13              | U |
| 75-09-2-----    | Methylene Chloride         | 13              | U |
| 67-64-1-----    | Acetone                    | 13 <del>9</del> | U |
| 75-15-0-----    | Carbon Disulfide           | 13              | U |
| 75-35-4-----    | 1,1-Dichloroethene         | 13              | U |
| 75-34-3-----    | 1,1-Dichloroethane         | 13              | U |
| 540-59-0-----   | 1,2-Dichloroethene (Total) | 13              | U |
| 67-66-3-----    | Chloroform                 | 13              | U |
| 107-06-2-----   | 1,2-Dichloroethane         | 13              | U |
| 78-93-3-----    | 2-Butanone                 | 13              | U |
| 71-55-6-----    | 1,1,1-Trichloroethane      | 13              | U |
| 56-23-5-----    | Carbon Tetrachloride       | 13              | U |
| 75-27-4-----    | Bromodichloromethane       | 13              | U |
| 78-87-5-----    | 1,2-Dichloropropane        | 13              | U |
| 10061-01-5----- | cis-1,3-Dichloropropene    | 13              | U |
| 79-01-6-----    | Trichloroethene            | 13              | U |
| 124-48-1-----   | Dibromochloromethane       | 13              | U |
| 79-00-5-----    | 1,1,2-Trichloroethane      | 13              | U |
| 71-43-2-----    | Benzene                    | 13              | U |
| 10061-02-6----- | trans-1,3-Dichloropropene  | 13              | U |
| 75-25-2-----    | Bromoform                  | 13              | U |
| 108-10-1-----   | 4-Methyl-2-Pentanone       | 13              | U |
| 591-78-6-----   | 2-Hexanone                 | 13              | U |
| 127-18-4-----   | Tetrachloroethene          | 13              | U |
| 79-34-5-----    | 1,1,2,2-Tetrachloroethane  | 13              | U |
| 108-88-3-----   | Toluene                    | 13              | U |
| 108-90-7-----   | Chlorobenzene              | 13              | U |
| 100-41-4-----   | Ethylbenzene               | 13              | U |
| 100-42-5-----   | Styrene                    | 13              | U |
| 1330-20-7-----  | Xylene (total)             | 13              | U |

1E  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

DUP-2

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) Soil

Lab Sample ID: L62615-6

Sample wt/vol: 5 (g/mL) g

Lab File ID: CD081111

Level: (low/med) LOW

Date Received: 08/10/00

%Moisture: not dec. 21

Date Analyzed: 08/11/00

GC Column: HP-624 ID: .25 (mm)

Dilution Factor: 1

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 Aromatic | 30.83 | 6          | 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.        |                  |       |            |   |

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

SB-6 (16-18)

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62574-1

Sample wt/vol: 30 (g/mL) g

Lab File ID: ED81206

Level: (low/med) LOW

Date Received: 08/08/00

% Moisture: 17 decanted: (Y/N) N

Date Extracted: 08/10/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/12/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg Q

|               |                              |      |   |
|---------------|------------------------------|------|---|
| 108-95-2----- | Phenol                       | 400  | U |
| 111-44-4----- | bis(2-Chloroethyl) ether     | 400  | U |
| 95-57-8-----  | 2-Chlorophenol               | 400  | U |
| 541-73-1----- | 1,3-Dichlorobenzene          | 400  | U |
| 106-46-7----- | 1,4-Dichlorobenzene          | 400  | U |
| 95-50-1-----  | 1,2-Dichlorobenzene          | 400  | U |
| 95-48-7-----  | 2-Methylphenol               | 400  | U |
| 108-60-1----- | 2,2'-oxybis(1-Chloropropane) | 400  | U |
| 106-44-5----- | 4-Methylphenol               | 400  | U |
| 621-64-7----- | N-Nitroso-di-n-propylamine   | 400  | U |
| 67-72-1-----  | Hexachloroethane             | 400  | U |
| 98-95-3-----  | Nitrobenzene                 | 400  | U |
| 78-59-1-----  | Isophorone                   | 400  | U |
| 88-75-5-----  | 2-Nitrophenol                | 400  | U |
| 105-67-9----- | 2,4-Dimethylphenol           | 400  | U |
| 111-91-1----- | bis(2-Chloroethoxy) methane  | 400  | U |
| 120-83-2----- | 2,4-Dichlorophenol           | 400  | U |
| 120-82-1----- | 1,2,4-Trichlorobenzene       | 400  | U |
| 91-20-3-----  | Naphthalene                  | 400  | U |
| 106-47-8----- | 4-Chloroaniline              | 400  | U |
| 87-68-3-----  | Hexachlorobutadiene          | 400  | U |
| 59-50-7-----  | 4-Chloro-3-methylphenol      | 400  | U |
| 91-57-6-----  | 2-Methylnaphthalene          | 400  | U |
| 77-47-4-----  | Hexachlorocyclopentadiene    | 400  | U |
| 88-06-2-----  | 2,4,6-Trichlorophenol        | 400  | U |
| 95-95-4-----  | 2,4,5-Trichlorophenol        | 1000 | U |
| 91-58-7-----  | 2-Chloronaphthalene          | 400  | U |
| 88-74-4-----  | 2-Nitroaniline               | 1000 | U |
| 131-11-3----- | Dimethylphthalate            | 400  | U |
| 208-96-8----- | Acenaphthylene               | 400  | U |
| 606-20-2----- | 2,6-Dinitrotoluene           | 400  | U |
| 99-09-2-----  | 3-Nitroaniline               | 1000 | U |
| 83-32-9-----  | Acenaphthene                 | 400  | U |

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SB-6 (16-18)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62574-1

Sample wt/vol: 30 (g/mL) g

Lab File ID: ED81206

Level: (low/med) LOW

Date Received: 08/08/00

% Moisture: 17 decanted: (Y/N) N

Date Extracted: 08/10/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/12/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|                |                            |      |   |
|----------------|----------------------------|------|---|
| 51-28-5-----   | 2,4-Dinitrophenol          | 1000 | U |
| 100-02-7-----  | 4-Nitrophenol              | 1000 | U |
| 132-64-9-----  | Dibenzofuran               | 400  | U |
| 121-14-2-----  | 2,4-Dinitrotoluene         | 400  | U |
| 84-66-2-----   | Diethylphthalate           | 400  | U |
| 86-73-7-----   | Fluorene                   | 400  | U |
| 7005-72-3----- | 4-Chlorophenyl-phenylether | 400  | U |
| 100-01-6-----  | 4-Nitroaniline             | 1000 | U |
| 534-52-1-----  | 4,6-Dinitro-2-methylphenol | 1000 | U |
| 86-30-6-----   | N-Nitrosodiphenylamine     | 400  | U |
| 101-55-3-----  | 4-Bromophenyl-phenylether  | 400  | U |
| 118-74-1-----  | Hexachlorobenzene          | 400  | U |
| 87-86-5-----   | Pentachlorophenol          | 1000 | U |
| 85-01-8-----   | Phenanthrene               | 400  | U |
| 120-12-7-----  | Anthracene                 | 400  | U |
| 86-74-8-----   | Carbazole                  | 400  | U |
| 84-74-2-----   | Di-n-butylphthalate        | 400  | U |
| 206-44-0-----  | Fluoranthene               | 400  | U |
| 129-00-0-----  | Pyrene                     | 400  | U |
| 85-68-7-----   | Butylbenzylphthalate       | 400  | U |
| 56-55-3-----   | Benzo(a)anthracene         | 400  | U |
| 91-94-1-----   | 3,3'-Dichlorobenzidine     | 400  | U |
| 218-01-9-----  | Chrysene                   | 400  | U |
| 117-81-7-----  | bis(2-Ethylhexyl)phthalate | 400  | U |
| 117-84-0-----  | Di-n-octylphthalate        | 400  | U |
| 205-99-2-----  | Benzo(b)fluoranthene       | 400  | U |
| 207-08-9-----  | Benzo(k)fluoranthene       | 400  | U |
| 50-32-8-----   | Benzo(a)pyrene             | 400  | U |
| 193-39-5-----  | Indeno(1,2,3-cd)pyrene     | 400  | U |
| 53-70-3-----   | Dibenzo(a,h)anthracene     | 400  | U |
| 191-24-2-----  | Benzo(g,h,i)perylene       | 400  | U |

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SB-6 (16-18)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62574-1

Sample wt/vol: 30 (g/mL) g

Lab File ID: ED81206

Level: (low/med) LOW

Date Received: 08/08/00

% Moisture: 17 decanted: (Y/N) N

Date Extracted: 08/10/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/12/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg

Number TICS found: 3

| CAS NUMBER     | COMPOUND NAME                | RT   | EST. CONC. | Q   |
|----------------|------------------------------|------|------------|-----|
| 1.             | Unknown                      | 3.71 | 1400       | JB  |
| 2. 000141-79-7 | 3-Penten-2-one, 4-methyl-    | 4.79 | 67000      | JBA |
| 3. 000123-42-2 | 2-Pentanone, 4-hydroxy-4-met | 5.59 | 3600       | JBA |
| 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.            |                              |      |            |     |

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SB-6 (18-20)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62574-4

Sample wt/vol: 30 (g/mL) g

Lab File ID: ED81209

Level: (low/med) LOW

Date Received: 08/08/00

% Moisture: 18 decanted: (Y/N) N

Date Extracted: 08/10/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/12/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|               |                              |      |   |
|---------------|------------------------------|------|---|
| 108-95-2----- | Phenol                       | 410  | U |
| 111-44-4----- | bis(2-Chloroethyl) ether     | 410  | U |
| 95-57-8-----  | 2-Chlorophenol               | 410  | U |
| 541-73-1----- | 1,3-Dichlorobenzene          | 410  | U |
| 106-46-7----- | 1,4-Dichlorobenzene          | 410  | U |
| 95-50-1-----  | 1,2-Dichlorobenzene          | 410  | U |
| 95-48-7-----  | 2-Methylphenol               | 410  | U |
| 108-60-1----- | 2,2'-oxybis(1-Chloropropane) | 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 |
| 120-82-1----- | 1,2,4-Trichlorobenzene       | 410  | U |
| 91-20-3-----  | Naphthalene                  | 410  | U |
| 106-47-8----- | 4-Chloroaniline              | 410  | U |
| 87-68-3-----  | Hexachlorobutadiene          | 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 |
| 91-58-7-----  | 2-Chloronaphthalene          | 410  | U |
| 88-74-4-----  | 2-Nitroaniline               | 1000 | U |
| 131-11-3----- | Dimethylphthalate            | 410  | U |
| 208-96-8----- | Acenaphthylene               | 410  | U |
| 606-20-2----- | 2,6-Dinitrotoluene           | 410  | U |
| 99-09-2-----  | 3-Nitroaniline               | 1000 | U |
| 83-32-9-----  | Acenaphthene                 | 410  | U |

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SB-6 (18-20)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62574-4

Sample wt/vol: 30 (g/mL) g

Lab File ID: ED81209

Level: (low/med) LOW

Date Received: 08/08/00

% Moisture: 18 decanted: (Y/N) N

Date Extracted: 08/10/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/12/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) 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     | 410  | U |
| 101-55-3-----  | 4-Bromophenyl-phenylether  | 410  | U |
| 118-74-1-----  | Hexachlorobenzene          | 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 |
| 56-55-3-----   | Benzo(a)anthracene         | 410  | U |
| 91-94-1-----   | 3,3'-Dichlorobenzidine     | 410  | U |
| 218-01-9-----  | Chrysene                   | 410  | U |
| 117-81-7-----  | bis(2-Ethylhexyl)phthalate | 410  | U |
| 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-----   | Dibenzo(a,h)anthracene     | 410  | U |
| 191-24-2-----  | Benzo(g,h,i)perylene       | 410  | U |

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SB-6 (18-20)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62574-4

Sample wt/vol: 30 (g/mL) g

Lab File ID: ED81209

Level: (low/med) LOW

Date Received: 08/08/00

% Moisture: 18 decanted: (Y/N) N

Date Extracted: 08/10/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/12/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg

Number TICS found: 5

| CAS NUMBER     | COMPOUND NAME                | RT    | EST. CONC. | Q   |
|----------------|------------------------------|-------|------------|-----|
| 1.             | Unknown                      | 3.72  | 1500       | JB  |
| 2. 000141-79-7 | 3-Penten-2-one, 4-methyl-    | 4.76  | 64000      | JBA |
| 3. 000123-42-2 | 2-Pentanone, 4-hydroxy-4-met | 5.58  | 3900       | JBA |
| 4.             | Unknown Hydrocarbon          | 14.50 | 520        | J   |
| 5.             | Unknown                      | 24.82 | 3300       | J   |
| 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.            |                              |       |            |     |



1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SB-8 (6.5-8)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-1

Sample wt/vol: 30 (g/mL) g

Lab File ID: ED82108

Level: (low/med) LOW

Date Received: 08/10/00

% Moisture: 75 decanted: (Y/N) N

Date Extracted: 08/14/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg Q

|               |                              |              |     |
|---------------|------------------------------|--------------|-----|
| 108-95-2----- | Phenol                       | 1300         | U   |
| 111-44-4----- | bis(2-Chloroethyl) ether     | 1300         | U   |
| 95-57-8-----  | 2-Chlorophenol               | 1300         | U   |
| 541-73-1----- | 1,3-Dichlorobenzene          | 1300         | U   |
| 106-46-7----- | 1,4-Dichlorobenzene          | 1300         | U   |
| 95-50-1-----  | 1,2-Dichlorobenzene          | 1300         | U   |
| 95-48-7-----  | 2-Methylphenol               | 1300         | U   |
| 108-60-1----- | 2,2'-oxybis(1-Chloropropane) | 1300         | U   |
| 106-44-5----- | 4-Methylphenol               | 180          | J   |
| 621-64-7----- | N-Nitroso-di-n-propylamine   | 1300         | U   |
| 67-72-1-----  | Hexachloroethane             | 1300         | U   |
| 98-95-3-----  | Nitrobenzene                 | 1300         | U   |
| 78-59-1-----  | Isophorone                   | 1300         | U   |
| 88-75-5-----  | 2-Nitrophenol                | 1300         | U   |
| 105-67-9----- | 2,4-Dimethylphenol           | 880          | J   |
| 111-91-1----- | bis(2-Chloroethoxy)methane   | 1300         | U   |
| 120-83-2----- | 2,4-Dichlorophenol           | 1300         | U   |
| 120-82-1----- | 1,2,4-Trichlorobenzene       | 1300         | U   |
| 91-20-3-----  | Naphthalene                  | 120000 62000 | E D |
| 106-47-8----- | 4-Chloroaniline              | 1300         | U   |
| 87-68-3-----  | Hexachlorobutadiene          | 1300         | U   |
| 59-50-7-----  | 4-Chloro-3-methylphenol      | 1300         | U   |
| 91-57-6-----  | 2-Methylnaphthalene          | 2500         |     |
| 77-47-4-----  | Hexachlorocyclopentadiene    | 1300         | U   |
| 88-06-2-----  | 2,4,6-Trichlorophenol        | 1300         | U   |
| 95-95-4-----  | 2,4,5-Trichlorophenol        | 3300         | U   |
| 91-58-7-----  | 2-Chloronaphthalene          | 1300         | U   |
| 88-74-4-----  | 2-Nitroaniline               | 3300         | U   |
| 131-11-3----- | Dimethylphthalate            | 1300         | U   |
| 208-96-8----- | Acenaphthylene               | 1200         | J   |
| 606-20-2----- | 2,6-Dinitrotoluene           | 1300         | U   |
| 99-09-2-----  | 3-Nitroaniline               | 3300         | U   |
| 83-32-9-----  | Acenaphthene                 | 26000 19000  | E D |

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

SB-8 (6.5-8)

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-1

Sample wt/vol: 30 (g/mL) g

Lab File ID: ED82108

Level: (low/med) LOW

Date Received: 08/10/00

% Moisture: 75 decanted: (Y/N) N

Date Extracted: 08/14/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|                |                            |       |            |
|----------------|----------------------------|-------|------------|
| 51-28-5-----   | 2,4-Dinitrophenol          | 3300  | U          |
| 100-02-7-----  | 4-Nitrophenol              | 3300  | U          |
| 132-64-9-----  | Dibenzofuran               | 10000 |            |
| 121-14-2-----  | 2,4-Dinitrotoluene         | 1300  | U          |
| 84-66-2-----   | Diethylphthalate           | 1300  | U          |
| 86-73-7-----   | Fluorene                   | 11000 | 9400 E DJ  |
| 7005-72-3----- | 4-Chlorophenyl-phenylether | 1300  | U          |
| 100-01-6-----  | 4-Nitroaniline             | 3300  | U          |
| 534-52-1-----  | 4,6-Dinitro-2-methylphenol | 3300  | U          |
| 86-30-6-----   | N-Nitrosodiphenylamine     | 1300  | U          |
| 101-55-3-----  | 4-Bromophenyl-phenylether  | 1300  | U          |
| 118-74-1-----  | Hexachlorobenzene          | 1300  | U          |
| 87-86-5-----   | Pentachlorophenol          | 3300  | U          |
| 85-01-8-----   | Phenanthrene               | 14000 | 12600 E DJ |
| 120-12-7-----  | Anthracene                 | 3300  |            |
| 86-74-8-----   | Carbazole                  | 7200  |            |
| 84-74-2-----   | Di-n-butylphthalate        | 1300  | U          |
| 206-44-0-----  | Fluoranthene               | 6700  |            |
| 129-00-0-----  | Pyrene                     | 8500  | J          |
| 85-68-7-----   | Butylbenzylphthalate       | 1300  | U W J      |
| 56-55-3-----   | Benzo(a)anthracene         | 3300  | J W J      |
| 91-94-1-----   | 3,3'-Dichlorobenzidine     | 1300  | U W J      |
| 218-01-9-----  | Chrysene                   | 2700  | J W J      |
| 117-81-7-----  | bis(2-Ethylhexyl)phthalate | 1300  | U W J      |
| 117-84-0-----  | Di-n-octylphthalate        | 1300  | U W J      |
| 205-99-2-----  | Benzo(b)fluoranthene       | 2900  | J W J      |
| 207-08-9-----  | Benzo(k)fluoranthene       | 940   | J W J      |
| 50-32-8-----   | Benzo(a)pyrene             | 2300  | J W J      |
| 193-39-5-----  | Indeno(1,2,3-cd)pyrene     | 1000  | J W J      |
| 53-70-3-----   | Dibenzo(a,h)anthracene     | 1300  | U W J      |
| 191-24-2-----  | Benzo(g,h,i)perylene       | 900   | J W J      |

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SB-8 (6.5-8)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-1

Sample wt/vol: 30 (g/mL) g

Lab File ID: ED82108

Level: (low/med) LOW

Date Received: 08/10/00

% Moisture: 75 decanted: (Y/N) N

Date Extracted: 08/14/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg

Number TICS found: 22

| CAS NUMBER      | COMPOUND NAME                | RT    | EST. CONC. | Q   |
|-----------------|------------------------------|-------|------------|-----|
| 1.              | Unknown                      | 3.46  | 4800       | JB  |
| 2. 000141-79-7  | 3-Penten-2-one, 4-methyl-    | 4.54  | 220000     | JBA |
| 3. 000123-42-2  | 2-Pentanone, 4-hydroxy-4-met | 5.37  | 13000      | JBA |
| 4.              | Unknown                      | 7.35  | 4200       | J   |
| 5. 000526-73-8  | Benzene, 1,2,3-trimethyl-    | 8.18  | 3100       | J   |
| 6.              | Unknown C9H12                | 8.62  | 2500       | J   |
| 7. 000496-11-7  | Indane                       | 8.86  | 27000      | J   |
| 8. 000095-13-6  | Indene                       | 8.99  | 3400       | J   |
| 9.              | Substituted Naphthalene      | 13.39 | 3800       | J   |
| 10.             | Substituted Naphthalene      | 13.70 | 2700       | J   |
| 11.             | Substituted Naphthalene      | 13.85 | 4900       | J   |
| 12.             | Unknown                      | 14.07 | 2300       | J   |
| 13.             | Unknown                      | 14.11 | 1900       | J   |
| 14. 000613-46-7 | 2-Naphthalenecarbonitrile    | 14.66 | 1800       | J   |
| 15. 007320-53-8 | Dibenzofuran, 4-methyl-      | 15.94 | 1700       | J   |
| 16. 000203-64-5 | 4H-Cyclopenta[def]phenanthre | 18.68 | 2000       | J   |
| 17.             | Unknown Hydrocarbon          | 19.95 | 4400       | J   |
| 18.             | Unknown                      | 20.24 | 1800       | J   |
| 19.             | Unknown Hydrocarbon          | 21.43 | 2300       | J   |
| 20. 000511-15-9 | 2-Phenanthrenol, 4b,5,6,7,8, | 21.46 | 7100       | J   |
| 21. 000511-15-9 | 2-Phenanthrenol, 4b,5,6,7,8, | 21.67 | 3100       | J   |
| 22.             | Unknown                      | 21.77 | 2300       | J   |
| 23.             | benzo(e) pyrene              | 24.85 | 1300       | J   |
| 24.             |                              |       |            |     |
| 25.             |                              |       |            |     |
| 26.             |                              |       |            |     |
| 27.             |                              |       |            |     |
| 28.             |                              |       |            |     |
| 29.             |                              |       |            |     |
| 30.             |                              |       |            |     |

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

SB-8 (6.5-8) DL

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-1DL

Sample wt/vol: 30 (g/mL) g

Lab File ID: ED82114

Level: (low/med) LOW

Date Received: 08/10/00

% Moisture: 75 decanted: (Y/N) N

Date Extracted: 08/14/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 20.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) ug/kg Q

|          |                              |       |    |
|----------|------------------------------|-------|----|
| 108-95-2 | Phenol                       | 27000 | U  |
| 111-44-4 | bis(2-Chloroethyl) ether     | 27000 | U  |
| 95-57-8  | 2-Chlorophenol               | 27000 | U  |
| 541-73-1 | 1,3-Dichlorobenzene          | 27000 | U  |
| 106-46-7 | 1,4-Dichlorobenzene          | 27000 | U  |
| 95-50-1  | 1,2-Dichlorobenzene          | 27000 | U  |
| 95-48-7  | 2-Methylphenol               | 27000 | U  |
| 108-60-1 | 2,2'-oxybis(1-Chloropropane) | 27000 | U  |
| 106-44-5 | 4-Methylphenol               | 27000 | U  |
| 621-64-7 | N-Nitroso-di-n-propylamine   | 27000 | U  |
| 67-72-1  | Hexachloroethane             | 27000 | U  |
| 98-95-3  | Nitrobenzene                 | 27000 | U  |
| 78-59-1  | Isophorone                   | 27000 | U  |
| 88-75-5  | 2-Nitrophenol                | 27000 | U  |
| 105-67-9 | 2,4-Dimethylphenol           | 27000 | U  |
| 111-91-1 | bis(2-Chloroethoxy) methane  | 27000 | U  |
| 120-83-2 | 2,4-Dichlorophenol           | 27000 | U  |
| 120-82-1 | 1,2,4-Trichlorobenzene       | 27000 | U  |
| 91-20-3  | Naphthalene                  | 62000 | D  |
| 106-47-8 | 4-Chloroaniline              | 27000 | U  |
| 87-68-3  | Hexachlorobutadiene          | 27000 | U  |
| 59-50-7  | 4-Chloro-3-methylphenol      | 27000 | U  |
| 91-57-6  | 2-Methylnaphthalene          | 27000 | U  |
| 77-47-4  | Hexachlorocyclopentadiene    | 27000 | U  |
| 88-06-2  | 2,4,6-Trichlorophenol        | 27000 | U  |
| 95-95-4  | 2,4,5-Trichlorophenol        | 67000 | U  |
| 91-58-7  | 2-Chloronaphthalene          | 27000 | U  |
| 88-74-4  | 2-Nitroaniline               | 67000 | U  |
| 131-11-3 | Dimethylphthalate            | 27000 | U  |
| 208-96-8 | Acenaphthylene               | 27000 | U  |
| 606-20-2 | 2,6-Dinitrotoluene           | 27000 | U  |
| 99-09-2  | 3-Nitroaniline               | 67000 | U  |
| 83-32-9  | Acenaphthene                 | 19000 | JD |

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SB-8 (6.5-8) DL

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-1DL

Sample wt/vol: 30 (g/mL) g

Lab File ID: ED82114

Level: (low/med) LOW

Date Received: 08/10/00

% Moisture: 75 decanted: (Y/N) N

Date Extracted: 08/14/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 20.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|                |                            |       |    |
|----------------|----------------------------|-------|----|
| 51-28-5-----   | 2,4-Dinitrophenol          | 67000 | U  |
| 100-02-7-----  | 4-Nitrophenol              | 67000 | U  |
| 132-64-9-----  | Dibenzofuran               | 9200  | JD |
| 121-14-2-----  | 2,4-Dinitrotoluene         | 27000 | U  |
| 84-66-2-----   | Diethylphthalate           | 27000 | U  |
| 86-73-7-----   | Fluorene                   | 9400  | JD |
| 7005-72-3----- | 4-Chlorophenyl-phenylether | 27000 | U  |
| 100-01-6-----  | 4-Nitroaniline             | 67000 | U  |
| 534-52-1-----  | 4,6-Dinitro-2-methylphenol | 67000 | U  |
| 86-30-6-----   | N-Nitrosodiphenylamine     | 27000 | U  |
| 101-55-3-----  | 4-Bromophenyl-phenylether  | 27000 | U  |
| 118-74-1-----  | Hexachlorobenzene          | 27000 | U  |
| 87-86-5-----   | Pentachlorophenol          | 67000 | U  |
| 85-01-8-----   | Phenanthrene               | 12000 | JD |
| 120-12-7-----  | Anthracene                 | 3100  | JD |
| 86-74-8-----   | Carbazole                  | 5800  | JD |
| 84-74-2-----   | Di-n-butylphthalate        | 27000 | U  |
| 206-44-0-----  | Fluoranthene               | 5100  | JD |
| 129-00-0-----  | Pyrene                     | 8200  | JD |
| 85-68-7-----   | Butylbenzylphthalate       | 27000 | U  |
| 56-55-3-----   | Benzo(a)anthracene         | 2700  | JD |
| 91-94-1-----   | 3,3'-Dichlorobenzidine     | 27000 | U  |
| 218-01-9-----  | Chrysene                   | 27000 | U  |
| 117-81-7-----  | bis(2-Ethylhexyl)phthalate | 27000 | U  |
| 117-84-0-----  | Di-n-octylphthalate        | 27000 | U  |
| 205-99-2-----  | Benzo(b)fluoranthene       | 2900  | JD |
| 207-08-9-----  | Benzo(k)fluoranthene       | 27000 | U  |
| 50-32-8-----   | Benzo(a)pyrene             | 27000 | U  |
| 193-39-5-----  | Indeno(1,2,3-cd)pyrene     | 27000 | U  |
| 53-70-3-----   | Dibenzo(a,h)anthracene     | 27000 | U  |
| 191-24-2-----  | Benzo(g,h,i)perylene       | 27000 | U  |

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SB-8 (6.5-8) DL

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-1 DL

Sample wt/vol: 30 (g/mL) g

Lab File ID: ED82114

Level: (low/med) LOW

Date Received: 08/10/00

% Moisture: 75 decanted: (Y/N) N

Date Extracted: 08/14/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 20.0

GPC Cleanup: (Y/N) Y pH: 7.0

Number TICS found: 12

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg

| CAS NUMBER     | COMPOUND NAME                | RT    | EST. CONC. | Q   |
|----------------|------------------------------|-------|------------|-----|
| 1.             | Unknown                      | 4.42  | 140000     | J   |
| 2. 000123-42-2 | 2-Pentanone, 4-hydroxy-4-met | 5.31  | 10000      | JBA |
| 3.             | Unknown C9H10                | 8.83  | 22000      | J   |
| 4.             | Unknown PAH                  | 11.13 | 12000      | J   |
| 5.             | Substituted Naphthalene      | 12.62 | 15000      | J   |
| 6.             | Unknown                      | 21.41 | 5700       | J   |
| 7.             | Unknown                      | 21.47 | 12000      | J   |
| 8.             | Unknown                      | 24.01 | 12000      | J   |
| 9.             | Unknown                      | 24.26 | 16000      | J   |
| 10.            | Unknown                      | 24.37 | 7500       | J   |
| 11.            | Unknown                      | 24.86 | 140000     | J   |
| 12.            | Unknown                      | 25.21 | 11000      | J   |
| 13.            |                              |       |            |     |
| 14.            |                              |       |            |     |
| 15.            |                              |       |            |     |
| 16.            |                              |       |            |     |
| 17.            |                              |       |            |     |
| 18.            |                              |       |            |     |
| 19.            |                              |       |            |     |
| 20.            |                              |       |            |     |
| 21.            |                              |       |            |     |
| 22.            |                              |       |            |     |
| 23.            |                              |       |            |     |
| 24.            |                              |       |            |     |
| 25.            |                              |       |            |     |
| 26.            |                              |       |            |     |
| 27.            |                              |       |            |     |
| 28.            |                              |       |            |     |
| 29.            |                              |       |            |     |
| 30.            |                              |       |            |     |

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SB-10 (4-7)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-2

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED82110

Level: (low/med) LOW

Date Received: 08/10/00

% Moisture: 18 decanted: (Y/N) N

Date Extracted: 08/14/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

|         |          |   |   |
|---------|----------|---|---|
| CAS NO. | COMPOUND | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) ug/kg | Q |
|---------|----------|---|---|

|               |                              |      |   |
|---------------|------------------------------|------|---|
| 108-95-2----- | Phenol                       | 400  | U |
| 111-44-4----- | bis(2-Chloroethyl) ether     | 400  | U |
| 95-57-8-----  | 2-Chlorophenol               | 400  | U |
| 541-73-1----- | 1,3-Dichlorobenzene          | 400  | U |
| 106-46-7----- | 1,4-Dichlorobenzene          | 400  | U |
| 95-50-1-----  | 1,2-Dichlorobenzene          | 400  | U |
| 95-48-7-----  | 2-Methylphenol               | 69   | J |
| 108-60-1----- | 2,2'-oxybis(1-Chloropropane) | 400  | U |
| 106-44-5----- | 4-Methylphenol               | 270  | J |
| 621-64-7----- | N-Nitroso-di-n-propylamine   | 400  | U |
| 67-72-1-----  | Hexachloroethane             | 400  | U |
| 98-95-3-----  | Nitrobenzene                 | 400  | U |
| 78-59-1-----  | Isophorone                   | 400  | U |
| 88-75-5-----  | 2-Nitrophenol                | 400  | U |
| 105-67-9----- | 2,4-Dimethylphenol           | 110  | J |
| 111-91-1----- | bis(2-Chloroethoxy) methane  | 400  | U |
| 120-83-2----- | 2,4-Dichlorophenol           | 400  | U |
| 120-82-1----- | 1,2,4-Trichlorobenzene       | 400  | U |
| 91-20-3-----  | Naphthalene                  | 3000 |   |
| 106-47-8----- | 4-Chloroaniline              | 400  | U |
| 87-68-3-----  | Hexachlorobutadiene          | 400  | U |
| 59-50-7-----  | 4-Chloro-3-methylphenol      | 400  | U |
| 91-57-6-----  | 2-Methylnaphthalene          | 1800 |   |
| 77-47-4-----  | Hexachlorocyclopentadiene    | 400  | U |
| 88-06-2-----  | 2,4,6-Trichlorophenol        | 400  | U |
| 95-95-4-----  | 2,4,5-Trichlorophenol        | 1000 | U |
| 91-58-7-----  | 2-Chloronaphthalene          | 400  | U |
| 88-74-4-----  | 2-Nitroaniline               | 1000 | U |
| 131-11-3----- | Dimethylphthalate            | 400  | U |
| 208-96-8----- | Acenaphthylene               | 4500 | E |
| 606-20-2----- | 2,6-Dinitrotoluene           | 400  | U |
| 99-09-2-----  | 3-Nitroaniline               | 1000 | U |
| 83-32-9-----  | Acenaphthene                 | 930  |   |

SAMPLE NO.

Contract: Blasland, B

SB-10 (4-7)

SDG No. : L62574

Lab Sample ID: L62615-2

Lab File ID: ED82110

Date Received: 08/10/00

Date Extracted:08/14/00

Date Analyzed: 08/21/00

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y      pH: 7.0

(ug/L or ug/Kg) ug/kg

Q

|                |                            |       |   |
|----------------|----------------------------|-------|---|
| 51-28-5-----   | 2,4-Dinitrophenol          | 1000  | U |
| 100-02-7-----  | 4-Nitrophenol              | 1000  | U |
| 132-64-9-----  | Dibenzofuran               | 4200  | E |
| 121-14-2-----  | 2,4-Dinitrotoluene         | 400   | U |
| 84-66-2-----   | Diethylphthalate           | 400   | U |
| 86-73-7-----   | Fluorene                   | 4800  | E |
| 7005-72-3----- | 4-Chlorophenyl-phenylether | 400   | U |
| 100-01-6-----  | 4-Nitroaniline             | 1000  | U |
| 534-52-1-----  | 4,6-Dinitro-2-methylphenol | 1000  | U |
| 86-30-6-----   | N-Nitrosodiphenylamine     | 400   | U |
| 101-55-3-----  | 4-Bromophenyl-phenylether  | 400   | U |
| 118-74-1-----  | Hexachlorobenzene          | 400   | U |
| 87-86-5-----   | Pentachlorophenol          | 1000  | U |
| 85-01-8-----   | Phenanthrene               | 33000 | E |
| 120-12-7-----  | Anthracene                 | 9100  | E |
| 86-74-8-----   | Carbazole                  | 1700  |   |
| 84-74-2-----   | Di-n-butylphthalate        | 400   | U |
| 206-44-0-----  | Fluoranthene               | 21000 | E |
| 129-00-0-----  | Pyrene                     | 42000 | E |
| 85-68-7-----   | Butylbenzylphthalate       | 400   | U |
| 56-55-3-----   | Benzo (a) anthracene       | 11000 | E |
| 91-94-1-----   | 3,3'-Dichlorobenzidine     | 400   | U |
| 218-01-9-----  | Chrysene                   | 9300  | E |
| 117-81-7-----  | bis(2-Ethylhexyl)phthalate | 400   | U |
| 117-84-0-----  | Di-n-octylphthalate        | 400   | U |
| 205-99-2-----  | Benzo (b) fluoranthene     | 11000 | E |
| 207-08-9-----  | Benzo (k) fluoranthene     | 3800  | E |
| 50-32-8-----   | Benzo (a) pyrene           | 9200  | E |
| 193-39-5-----  | Indeno (1,2,3-cd) pyrene   | 5300  | E |
| 53-70-3-----   | Dibenzo (a,h) anthracene   | 810   |   |
| 191-24-2-----  | Benzo (g,h,i) perylene     | 5100  | E |



1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SB-10 (4-7)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-2

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED82110

Level: (low/med) LOW

Date Received: 08/10/00

% Moisture: 18 decanted: (Y/N) N

Date Extracted: 08/14/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:

Number TICS found: 22

(ug/L or ug/Kg) ug/kg

| CAS NUMBER      | COMPOUND NAME                | RT    | EST. CONC. | Q   |
|-----------------|------------------------------|-------|------------|-----|
| 1.              | Unknown                      | 3.46  | 950        | J   |
| 2.              | Unknown                      | 3.59  | 750        | JB  |
| 3. 010574-37-5  | 2-Pentene, 2,3-dimethyl-     | 4.56  | 45000      | JBA |
| 4. 000123-42-2  | 2-Pentanone, 4-hydroxy-4-met | 5.38  | 2600       | JBA |
| 5.              | Unknown                      | 7.36  | 1100       | J   |
| 6.              | Substituted Naphthalene      | 12.63 | 1500       | J   |
| 7.              | Substituted Naphthalene      | 13.70 | 660        | J   |
| 8.              | Substituted Naphthalene      | 13.86 | 1100       | J   |
| 9.              | Substituted Naphthalene      | 13.90 | 740        | J   |
| 10.             | Substituted Naphthalene      | 14.08 | 660        | J   |
| 11. 007320-53-8 | Dibenzofuran, 4-methyl-      | 15.95 | 1100       | J   |
| 12.             | Unknown PAH                  | 16.11 | 2900       | J   |
| 13.             | Unknown PAH                  | 17.01 | 620        | J   |
| 14. 000132-65-0 | Dibenzothiophene             | 17.26 | 1500       | J   |
| 15.             | Unknown PAH                  | 18.51 | 1400       | J   |
| 16.             | Unknown PAH                  | 18.57 | 1900       | J   |
| 17.             | Unknown PAH                  | 18.68 | 1000       | J   |
| 18.             | Unknown PAH                  | 18.73 | 2800       | J   |
| 19.             | Unknown PAH                  | 18.78 | 770        | J   |
| 20.             | Unknown PAH                  | 19.13 | 1000       | J   |
| 21.             | Unknown PAH                  | 19.62 | 690        | J   |
| 22.             | Unknown PAH                  | 20.11 | 1700       | J   |
| 23.             | benzo(e)pyrene               | 24.89 | 4600       | J   |
| 24.             |                              |       |            |     |
| 25.             |                              |       |            |     |
| 26.             |                              |       |            |     |
| 27.             |                              |       |            |     |
| 28.             |                              |       |            |     |
| 29.             |                              |       |            |     |
| 30.             |                              |       |            |     |

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SB-10 (4-7) DL

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-2DL

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED82116

Level: (low/med) LOW

Date Received: 08/10/00

% Moisture: 18 decanted: (Y/N) N

Date Extracted: 08/14/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 20.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|               |                              |       |    |
|---------------|------------------------------|-------|----|
| 108-95-2----- | Phenol                       | 8100  | U  |
| 111-44-4----- | bis(2-Chloroethyl) ether     | 8100  | U  |
| 95-57-8-----  | 2-Chlorophenol               | 8100  | U  |
| 541-73-1----- | 1,3-Dichlorobenzene          | 8100  | U  |
| 106-46-7----- | 1,4-Dichlorobenzene          | 8100  | U  |
| 95-50-1-----  | 1,2-Dichlorobenzene          | 8100  | U  |
| 95-48-7-----  | 2-Methylphenol               | 8100  | U  |
| 108-60-1----- | 2,2'-oxybis(1-Chloropropane) | 8100  | U  |
| 106-44-5----- | 4-Methylphenol               | 8100  | U  |
| 621-64-7----- | N-Nitroso-di-n-propylamine   | 8100  | U  |
| 67-72-1-----  | Hexachloroethane             | 8100  | U  |
| 98-95-3-----  | Nitrobenzene                 | 8100  | U  |
| 78-59-1-----  | Isophorone                   | 8100  | U  |
| 88-75-5-----  | 2-Nitrophenol                | 8100  | U  |
| 105-67-9----- | 2,4-Dimethylphenol           | 8100  | U  |
| 111-91-1----- | bis(2-Chloroethoxy) methane  | 8100  | U  |
| 120-83-2----- | 2,4-Dichlorophenol           | 8100  | U  |
| 120-82-1----- | 1,2,4-Trichlorobenzene       | 8100  | U  |
| 91-20-3-----  | Naphthalene                  | 3200  | JD |
| 106-47-8----- | 4-Chloroaniline              | 8100  | U  |
| 87-68-3-----  | Hexachlorobutadiene          | 8100  | U  |
| 59-50-7-----  | 4-Chloro-3-methylphenol      | 8100  | U  |
| 91-57-6-----  | 2-Methylnaphthalene          | 1700  | JD |
| 77-47-4-----  | Hexachlorocyclopentadiene    | 8100  | U  |
| 88-06-2-----  | 2,4,6-Trichlorophenol        | 8100  | U  |
| 95-95-4-----  | 2,4,5-Trichlorophenol        | 20000 | U  |
| 91-58-7-----  | 2-Chloronaphthalene          | 8100  | U  |
| 88-74-4-----  | 2-Nitroaniline               | 20000 | U  |
| 131-11-3----- | Dimethylphthalate            | 8100  | U  |
| 208-96-8----- | Acenaphthylene               | 4500  | JD |
| 606-20-2----- | 2,6-Dinitrotoluene           | 8100  | U  |
| 99-09-2-----  | 3-Nitroaniline               | 20000 | U  |
| 83-32-9-----  | Acenaphthene                 | 1000  | JD |

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SB-10 (4-7) DL

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-2DL

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED82116

Level: (low/med) LOW

Date Received: 08/10/00

% Moisture: 18 decanted: (Y/N) N

Date Extracted: 08/14/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 20.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|                |                            |       |    |
|----------------|----------------------------|-------|----|
| 51-28-5-----   | 2,4-Dinitrophenol          | 20000 | U  |
| 100-02-7-----  | 4-Nitrophenol              | 20000 | U  |
| 132-64-9-----  | Dibenzofuran               | 3800  | JD |
| 121-14-2-----  | 2,4-Dinitrotoluene         | 8100  | U  |
| 84-66-2-----   | Diethylphthalate           | 8100  | U  |
| 86-73-7-----   | Fluorene                   | 4500  | JD |
| 7005-72-3----- | 4-Chlorophenyl-phenylether | 8100  | U  |
| 100-01-6-----  | 4-Nitroaniline             | 20000 | U  |
| 534-52-1-----  | 4,6-Dinitro-2-methylphenol | 20000 | U  |
| 86-30-6-----   | N-Nitrosodiphenylamine     | 8100  | U  |
| 101-55-3-----  | 4-Bromophenyl-phenylether  | 8100  | U  |
| 118-74-1-----  | Hexachlorobenzene          | 8100  | U  |
| 87-86-5-----   | Pentachlorophenol          | 20000 | U  |
| 85-01-8-----   | Phenanthrene               | 23000 | D  |
| 120-12-7-----  | Anthracene                 | 7400  | JD |
| 86-74-8-----   | Carbazole                  | 1800  | JD |
| 84-74-2-----   | Di-n-butylphthalate        | 8100  | U  |
| 206-44-0-----  | Fluoranthene               | 18000 | D  |
| 129-00-0-----  | Pyrene                     | 32000 | D  |
| 85-68-7-----   | Butylbenzylphthalate       | 8100  | U  |
| 56-55-3-----   | Benzo(a)anthracene         | 11000 | D  |
| 91-94-1-----   | 3,3'-Dichlorobenzidine     | 8100  | U  |
| 218-01-9-----  | Chrysene                   | 9800  | D  |
| 117-81-7-----  | bis(2-Ethylhexyl)phthalate | 8100  | U  |
| 117-84-0-----  | Di-n-octylphthalate        | 8100  | U  |
| 205-99-2-----  | Benzo(b)fluoranthene       | 14000 | D  |
| 207-08-9-----  | Benzo(k)fluoranthene       | 5800  | JD |
| 50-32-8-----   | Benzo(a)pyrene             | 12000 | D  |
| 193-39-5-----  | Indeno(1,2,3-cd)pyrene     | 6500  | JD |
| 53-70-3-----   | Dibenzo(a,h)anthracene     | 1800  | JD |
| 191-24-2-----  | Benzo(g,h,i)perylene       | 5800  | JD |

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SB-10 (4-7) DL

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-2 DL

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED82116

Level: (low/med) LOW

Date Received: 08/10/00

% Moisture: 18 decanted: (Y/N) N

Date Extracted: 08/14/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 20.0

GPC Cleanup: (Y/N) Y pH: 7.0

Number TICS found: 22

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg

| CAS NUMBER     | COMPOUND NAME                | RT    | EST. CONC. | Q   |
|----------------|------------------------------|-------|------------|-----|
| 1.             | Unknown                      | 3.46  | 1600       | JB  |
| 2. 000141-79-7 | 3-Penten-2-one, 4-methyl     | 4.42  | 57000      | JBA |
| 3. 000123-42-2 | 2-Pentanone, 4-hydroxy-4-met | 5.31  | 4500       | JBA |
| 4.             | Unknown                      | 7.32  | 1800       | J   |
| 5.             | Unknown PAH                  | 18.48 | 2000       | J   |
| 6.             | Unknown PAH                  | 18.52 | 3200       | J   |
| 7.             | Unknown PAH                  | 18.67 | 3900       | J   |
| 8.             | Unknown PAH                  | 18.73 | 1800       | J   |
| 9.             | Unknown PAH                  | 19.10 | 2500       | J   |
| 10.            | Unknown PAH                  | 20.05 | 2500       | J   |
| 11.            | Unknown PAH                  | 21.00 | 1900       | J   |
| 12.            | Unknown PAH                  | 21.14 | 2100       | J   |
| 13.            | Unknown PAH                  | 22.29 | 1700       | J   |
| 14.            | Unknown PAH                  | 23.51 | 1700       | J   |
| 15.            | Unknown PAH                  | 23.89 | 2500       | J   |
| 16.            | Unknown                      | 23.97 | 4900       | J   |
| 17.            | Unknown                      | 24.04 | 2700       | J   |
| 18.            | Unknown PAH                  | 24.64 | 5000       | J   |
| 19.            | Unknown                      | 25.45 | 1800       | J   |
| 20.            | Unknown                      | 26.12 | 19000      | J   |
| 21.            | Unknown                      | 26.26 | 2900       | J   |
| 22.            | Unknown PAH                  | 26.58 | 1700       | J   |
| 23.            | benzo(e)pyrene               | 24.85 | 6000       | J   |
| 24.            |                              |       |            |     |
| 25.            |                              |       |            |     |
| 26.            |                              |       |            |     |
| 27.            |                              |       |            |     |
| 28.            |                              |       |            |     |
| 29.            |                              |       |            |     |
| 30.            |                              |       |            |     |

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SB-13 (4-6)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-3

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED82111

Level: (low/med) LOW

Date Received: 08/10/00

% Moisture: 26 decanted: (Y/N) N

Date Extracted: 08/14/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|               |                              |      |   |
|---------------|------------------------------|------|---|
| 108-95-2----- | Phenol                       | 450  | U |
| 111-44-4----- | bis(2-Chloroethyl) ether     | 450  | U |
| 95-57-8-----  | 2-Chlorophenol               | 450  | U |
| 541-73-1----- | 1,3-Dichlorobenzene          | 450  | U |
| 106-46-7----- | 1,4-Dichlorobenzene          | 450  | U |
| 95-50-1-----  | 1,2-Dichlorobenzene          | 450  | U |
| 95-48-7-----  | 2-Methylphenol               | 450  | U |
| 108-60-1----- | 2,2'-oxybis(1-Chloropropane) | 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 |
| 120-82-1----- | 1,2,4-Trichlorobenzene       | 450  | U |
| 91-20-3-----  | Naphthalene                  | 300  | J |
| 106-47-8----- | 4-Chloroaniline              | 450  | U |
| 87-68-3-----  | Hexachlorobutadiene          | 450  | U |
| 59-50-7-----  | 4-Chloro-3-methylphenol      | 450  | U |
| 91-57-6-----  | 2-Methylnaphthalene          | 300  | J |
| 77-47-4-----  | Hexachlorocyclopentadiene    | 450  | U |
| 88-06-2-----  | 2,4,6-Trichlorophenol        | 450  | U |
| 95-95-4-----  | 2,4,5-Trichlorophenol        | 1100 | U |
| 91-58-7-----  | 2-Chloronaphthalene          | 450  | U |
| 88-74-4-----  | 2-Nitroaniline               | 1100 | U |
| 131-11-3----- | Dimethylphthalate            | 450  | U |
| 208-96-8----- | Acenaphthylene               | 770  |   |
| 606-20-2----- | 2,6-Dinitrotoluene           | 450  | U |
| 99-09-2-----  | 3-Nitroaniline               | 1100 | U |
| 83-32-9-----  | Acenaphthene                 | 210  | J |

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SB-13 (4-6)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-3

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED82111

Level: (low/med) LOW

Date Received: 08/10/00

% Moisture: 26 decanted: (Y/N) N

Date Extracted: 08/14/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

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               | 620   |     |
| 121-14-2-----  | 2,4-Dinitrotoluene         | 450   | U   |
| 84-66-2-----   | Diethylphthalate           | 450   | U   |
| 86-73-7-----   | Fluorene                   | 850   |     |
| 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     | 450   | U   |
| 101-55-3-----  | 4-Bromophenyl-phenylether  | 450   | U   |
| 118-74-1-----  | Hexachlorobenzene          | 450   | U   |
| 87-86-5-----   | Pentachlorophenol          | 1100  | U   |
| 85-01-8-----   | Phenanthrene               | 7700  | E D |
| 120-12-7-----  | Anthracene                 | 2000  |     |
| 86-74-8-----   | Carbazole                  | 430   | J   |
| 84-74-2-----   | Di-n-butylphthalate        | 450   | U   |
| 206-44-0-----  | Fluoranthene               | 6800  | E D |
| 129-00-0-----  | Pyrene                     | 15000 | E D |
| 85-68-7-----   | Butylbenzylphthalate       | 450   | U   |
| 56-55-3-----   | Benzo(a)anthracene         | 4500  | E D |
| 91-94-1-----   | 3,3'-Dichlorobenzidine     | 450   | U   |
| 218-01-9-----  | Chrysene                   | 4000  | E D |
| 117-81-7-----  | bis(2-Ethylhexyl)phthalate | 450   | U   |
| 117-84-0-----  | Di-n-octylphthalate        | 450   | U   |
| 205-99-2-----  | Benzo(b)fluoranthene       | 3800  | E D |
| 207-08-9-----  | Benzo(k)fluoranthene       | 1400  |     |
| 50-32-8-----   | Benzo(a)pyrene             | 3100  |     |
| 193-39-5-----  | Indeno(1,2,3-cd)pyrene     | 1700  |     |
| 53-70-3-----   | Dibenzo(a,h)anthracene     | 510   |     |
| 191-24-2-----  | Benzo(g,h,i)perylene       | 1500  |     |

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SB-13 (4-6)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-3

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED82111

Level: (low/med) LOW

Date Received: 08/10/00

% Moisture: 26 decanted: (Y/N) N

Date Extracted: 08/14/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg

Number TICS found: 11

| CAS NUMBER     | COMPOUND NAME                | RT    | EST. CONC. | Q   |
|----------------|------------------------------|-------|------------|-----|
| 1.             | Unknown                      | 3.47  | 1300       | JB  |
| 2. 000141-79-7 | 3-Penten-2-one, 4-methyl-    | 4.59  | 65000      | JBA |
| 3. 000123-42-2 | 2-Pentanone, 4-hydroxy-4-met | 5.39  | 3400       | JBA |
| 4.             | Unknown                      | 7.37  | 1600       | J   |
| 5.             | Unknown PAH                  | 18.50 | 1100       | J   |
| 6.             | Unknown PAH                  | 18.55 | 1600       | J   |
| 7.             | Unknown PAH                  | 18.70 | 1500       | J   |
| 8.             | Substituted Naphthalene      | 19.11 | 980        | J   |
| 9.             | Unknown PAH                  | 19.61 | 860        | J   |
| 10.            | Unknown PAH                  | 20.08 | 1000       | J   |
| 11.            | Unknown PAH                  | 21.03 | 310        | J   |
| 12.            | Benzofuran                   | 24.87 | 1600       | J   |
| 13.            |                              |       |            |     |
| 14.            |                              |       |            |     |
| 15.            |                              |       |            |     |
| 16.            |                              |       |            |     |
| 17.            |                              |       |            |     |
| 18.            |                              |       |            |     |
| 19.            |                              |       |            |     |
| 20.            |                              |       |            |     |
| 21.            |                              |       |            |     |
| 22.            |                              |       |            |     |
| 23.            |                              |       |            |     |
| 24.            |                              |       |            |     |
| 25.            |                              |       |            |     |
| 26.            |                              |       |            |     |
| 27.            |                              |       |            |     |
| 28.            |                              |       |            |     |
| 29.            |                              |       |            |     |
| 30.            |                              |       |            |     |

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

SB-13 (4-6) DL

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-3DL

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED82117

Level: (low/med) LOW

Date Received: 08/10/00

% Moisture: 26 decanted: (Y/N) N

Date Extracted: 08/14/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 10.0

GPC Cleanup: (Y/N) Y pH: 7.0

| CAS NO. | COMPOUND | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) ug/kg | Q |
|---------|----------|---|---|
|---------|----------|---|---|

|               |                              |       |    |
|---------------|------------------------------|-------|----|
| 108-95-2----- | Phenol                       | 4500  | U  |
| 111-44-4----- | bis(2-Chloroethyl) ether     | 4500  | U  |
| 95-57-8-----  | 2-Chlorophenol               | 4500  | U  |
| 541-73-1----- | 1,3-Dichlorobenzene          | 4500  | U  |
| 106-46-7----- | 1,4-Dichlorobenzene          | 4500  | U  |
| 95-50-1-----  | 1,2-Dichlorobenzene          | 4500  | U  |
| 95-48-7-----  | 2-Methylphenol               | 4500  | U  |
| 108-60-1----- | 2,2'-oxybis(1-Chloropropane) | 4500  | U  |
| 106-44-5----- | 4-Methylphenol               | 4500  | U  |
| 621-64-7----- | N-Nitroso-di-n-propylamine   | 4500  | U  |
| 67-72-1-----  | Hexachloroethane             | 4500  | U  |
| 98-95-3-----  | Nitrobenzene                 | 4500  | U  |
| 78-59-1-----  | Isophorone                   | 4500  | U  |
| 88-75-5-----  | 2-Nitrophenol                | 4500  | U  |
| 105-67-9----- | 2,4-Dimethylphenol           | 4500  | U  |
| 111-91-1----- | bis(2-Chloroethoxy) methane  | 4500  | U  |
| 120-83-2----- | 2,4-Dichlorophenol           | 4500  | U  |
| 120-82-1----- | 1,2,4-Trichlorobenzene       | 4500  | U  |
| 91-20-3-----  | Naphthalene                  | 4500  | U  |
| 106-47-8----- | 4-Chloroaniline              | 4500  | U  |
| 87-68-3-----  | Hexachlorobutadiene          | 4500  | U  |
| 59-50-7-----  | 4-Chloro-3-methylphenol      | 4500  | U  |
| 91-57-6-----  | 2-Methylnaphthalene          | 4500  | U  |
| 77-47-4-----  | Hexachlorocyclopentadiene    | 4500  | U  |
| 88-06-2-----  | 2,4,6-Trichlorophenol        | 4500  | U  |
| 95-95-4-----  | 2,4,5-Trichlorophenol        | 11000 | U  |
| 91-58-7-----  | 2-Chloronaphthalene          | 4500  | U  |
| 88-74-4-----  | 2-Nitroaniline               | 11000 | U  |
| 131-11-3----- | Dimethylphthalate            | 4500  | U  |
| 208-96-8----- | Acenaphthylene               | 660   | JD |
| 606-20-2----- | 2,6-Dinitrotoluene           | 4500  | U  |
| 99-09-2-----  | 3-Nitroaniline               | 11000 | U  |
| 83-32-9-----  | Acenaphthene                 | 4500  | U  |

FORM I SV-I

*use original*



1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

SB-13 (4-6) DL

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-3DL

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED82117

Level: (low/med) LOW

Date Received: 08/10/00

% Moisture: 26 decanted: (Y/N) N

Date Extracted: 08/14/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 10.0

GPC Cleanup: (Y/N) Y pH: 7.0

| CAS NO.   | COMPOUND                   | CONCENTRATION UNITS: |         |
|-----------|----------------------------|----------------------|---------|
|           |                            | (ug/L or ug/Kg)      | ug/kg Q |
| 51-28-5   | 2,4-Dinitrophenol          | 11000                | U       |
| 100-02-7  | 4-Nitrophenol              | 11000                | U       |
| 132-64-9  | Dibenzofuran               | 520                  | JD      |
| 121-14-2  | 2,4-Dinitrotoluene         | 4500                 | U       |
| 84-66-2   | Diethylphthalate           | 4500                 | U       |
| 86-73-7   | Fluorene                   | 710                  | JD      |
| 7005-72-3 | 4-Chlorophenyl phenylether | 4500                 | U       |
| 100-01-6  | 4-Nitroaniline             | 11000                | U       |
| 534-52-1  | 4,6-Dinitro-2-methylphenol | 11000                | U       |
| 86-30-6   | N-Nitrosodiphenylamine     | 4500                 | U       |
| 101-55-3  | 4-Bromophenyl phenylether  | 4500                 | U       |
| 118-74-1  | Hexachlorobenzene          | 4500                 | U       |
| 87-86-5   | Pentachlorophenol          | 11000                | U       |
| 85-01-8   | Phenanthrene               | 6200                 | D       |
| 120-12-7  | Anthracene                 | 1700                 | JD      |
| 86-74-8   | Carbazole                  | 4500                 | U       |
| 84-74-2   | Di-n-butylphthalate        | 4500                 | U       |
| 206-44-0  | Fluoranthene               | 5600                 | D       |
| 129-00-0  | Pyrene                     | 10000                | D J     |
| 85-68-7   | Butylbenzylphthalate       | 4500                 | U W     |
| 56-55-3   | Benzo(a)anthracene         | 3600                 | JD J    |
| 91-94-1   | 3,3'-Dichlorobenzidine     | 4500                 | U W     |
| 218-01-9  | Chrysene                   | 3200                 | JD J    |
| 117-81-7  | bis(2-Ethylhexyl)phthalate | 4500                 | U W     |
| 117-84-0  | Di-n-octylphthalate        | 4500                 | U W     |
| 205-99-2  | Benzo(b)fluoranthene       | 3500                 | JD J    |
| 207-08-9  | Benzo(k)fluoranthene       | 1400                 | JD J    |
| 50-32-8   | Benzo(a)pyrene             | 3300                 | JD J    |
| 193-39-5  | Indeno(1,2,3-cd)pyrene     | 1600                 | JD J    |
| 53-70-3   | Dibenzo(a,h)anthracene     | 4500                 | U W     |
| 191-24-2  | Benzo(g,h,i)perylene       | 1600                 | JD J    |

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

SB-13 (4-6) DL

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-3 DL

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED82117

Level: (low/med) LOW

Date Received: 08/10/00

% Moisture: 26 decanted: (Y/N) N

Date Extracted: 08/14/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 10.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg

Number TICS found: 9

| CAS NUMBER     | COMPOUND NAME                | RT    | EST. CONC. | Q   |
|----------------|------------------------------|-------|------------|-----|
| 1.             | Unknown                      | 3.45  | 2200       | JB  |
| 2.             | Unknown                      | 4.43  | 76000      | J   |
| 3. 000123-42-2 | 2-Pentanone, 4-hydroxy 4-met | 5.31  | 5200       | JBA |
| 4.             | Unknown                      | 7.32  | 2300       | J   |
| 5.             | Unknown PAH                  | 18.53 | 1300       | J   |
| 6.             | Unknown PAH                  | 18.67 | 1300       | J   |
| 7.             | Unknown PAH                  | 19.10 | 1100       | J   |
| 8.             | Unknown PAH                  | 21.00 | 980        | J   |
| 9.             | Unknown PAH                  | 21.13 | 910        | J   |
| 10.            | benzo(e)pyrene               | 24.85 | 1600       | J   |
| 11.            |                              |       |            |     |
| 12.            |                              |       |            |     |
| 13.            |                              |       |            |     |
| 14.            |                              |       |            |     |
| 15.            |                              |       |            |     |
| 16.            |                              |       |            |     |
| 17.            |                              |       |            |     |
| 18.            |                              |       |            |     |
| 19.            |                              |       |            |     |
| 20.            |                              |       |            |     |
| 21.            |                              |       |            |     |
| 22.            |                              |       |            |     |
| 23.            |                              |       |            |     |
| 24.            |                              |       |            |     |
| 25.            |                              |       |            |     |
| 26.            |                              |       |            |     |
| 27.            |                              |       |            |     |
| 28.            |                              |       |            |     |
| 29.            |                              |       |            |     |
| 30.            |                              |       |            |     |

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SB-14 (5-7)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-5

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED82112

Level: (low/med) LOW

Date Received: 08/10/00

% Moisture: 28 decanted: (Y/N) N

Date Extracted: 08/14/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg Q

|               |                              |            |                |
|---------------|------------------------------|------------|----------------|
| 108-95-2----- | Phenol                       | 460        | U              |
| 111-44-4----- | bis(2-Chloroethyl) ether     | 460        | U              |
| 95-57-8-----  | 2-Chlorophenol               | 460        | U              |
| 541-73-1----- | 1,3-Dichlorobenzene          | 460        | U              |
| 106-46-7----- | 1,4-Dichlorobenzene          | 460        | U              |
| 95-50-1-----  | 1,2-Dichlorobenzene          | 460        | U              |
| 95-48-7-----  | 2-Methylphenol               | 460        | U              |
| 108-60-1----- | 2,2'-oxybis(1-Chloropropane) | 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              |
| 120-82-1----- | 1,2,4-Trichlorobenzene       | 460        | U              |
| 91-20-3-----  | Naphthalene                  | 380        | J              |
| 106-47-8----- | 4-Chloroaniline              | 460        | U              |
| 87-68-3-----  | Hexachlorobutadiene          | 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              |
| 91-58-7-----  | 2-Chloronaphthalene          | 460        | U              |
| 88-74-4-----  | 2-Nitroaniline               | 1200       | U              |
| 131-11-3----- | Dimethylphthalate            | 460        | U              |
| 208-96-8----- | Acenaphthylene               | 580        |                |
| 606-20-2----- | 2,6-Dinitrotoluene           | 460        | U              |
| 99-09-2-----  | 3-Nitroaniline               | 1200       | U              |
| 83-32-9-----  | Acenaphthene                 | 10000 8700 | <del>E</del> D |

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

SB-14 (5-7)

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-5

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED82112

Level: (low/med) LOW

Date Received: 08/10/00

% Moisture: 28 decanted: (Y/N) N

Date Extracted: 08/14/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg Q

|                |                            |             |   |   |
|----------------|----------------------------|-------------|---|---|
| 51-28-5-----   | 2,4-Dinitrophenol          | 1200        | U |   |
| 100-02-7-----  | 4-Nitrophenol              | 1200        | U |   |
| 132-64-9-----  | Dibenzofuran               | 6000 5800   | E | D |
| 121-14-2-----  | 2,4-Dinitrotoluene         | 460         | U |   |
| 84-66-2-----   | Diethylphthalate           | 460         | U |   |
| 86-73-7-----   | Fluorene                   | 8400 7000   | E | D |
| 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     | 460         | U |   |
| 101-55-3-----  | 4-Bromophenyl-phenylether  | 460         | U |   |
| 118-74-1-----  | Hexachlorobenzene          | 460         | U |   |
| 87-86-5-----   | Pentachlorophenol          | 1200        | U |   |
| 85-01-8-----   | Phenanthrene               | 34000 23000 | E | D |
| 120-12-7-----  | Anthracene                 | 7500 6700   | E | D |
| 86-74-8-----   | Carbazole                  | 2800        |   |   |
| 84-74-2-----   | Di-n-butylphthalate        | 460         | U |   |
| 206-44-0-----  | Fluoranthene               | 11000 10000 | E | D |
| 129-00-0-----  | Pyrene                     | 18000 15000 | E | D |
| 85-68-7-----   | Butylbenzylphthalate       | 460         | U |   |
| 56-55-3-----   | Benzo (a) anthracene       | 3500        | U |   |
| 91-94-1-----   | 3,3'-Dichlorobenzidine     | 460         | U |   |
| 218-01-9-----  | Chrysene                   | 2700        | U |   |
| 117-81-7-----  | bis(2-Ethylhexyl)phthalate | 460         | U |   |
| 117-84-0-----  | Di-n-octylphthalate        | 460         | U |   |
| 205-99-2-----  | Benzo (b) fluoranthene     | 2600        | U |   |
| 207-08-9-----  | Benzo (k) fluoranthene     | 1100        | U |   |
| 50-32-8-----   | Benzo (a) pyrene           | 2400        | U |   |
| 193-39-5-----  | Indeno (1,2,3-cd) pyrene   | 1400        | U |   |
| 53-70-3-----   | Dibenzo (a,h) anthracene   | 340         | U |   |
| 191-24-2-----  | Benzo (g,h,i) perylene     | 1200        | U |   |

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SB-14 (5-7)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-5

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED82112

Level: (low/med) LOW

Date Received: 08/10/00

% Moisture: 28 decanted: (Y/N) N

Date Extracted: 08/14/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:

(ug/L or ug/Kg) ug/kg

Number TICS found: 22

| CAS NUMBER     | COMPOUND NAME                | RT    | EST. CONC. | Q     |
|----------------|------------------------------|-------|------------|-------|
| =====          | =====                        | ===== | =====      | ===== |
| 1.             | Unknown                      | 3.46  | 1100       | JB    |
| 2. 010574-37-5 | 2-Pentene, 2,3-dimethyl-     | 4.57  | 57000      | JBA   |
| 3. 000123-42-2 | 2-Pentanone, 4-hydroxy-4-met | 5.39  | 3100       | JBA   |
| 4.             | Unknown                      | 7.36  | 1200       | J     |
| 5. 000496-11-7 | Indane                       | 8.84  | 690        | J     |
| 6.             | Unknown C10H9N               | 12.58 | 860        | J     |
| 7.             | Substituted Naphthalene      | 12.64 | 3400       | J     |
| 8. 000827-54-3 | Naphthalene, 2-ethenyl-      | 13.40 | 1300       | J     |
| 9.             | Substituted Naphthalene      | 13.71 | 1200       | J     |
| 10.            | Substituted Naphthalene      | 13.73 | 860        | J     |
| 11.            | Substituted Naphthalene      | 13.87 | 2700       | J     |
| 12.            | Substituted Naphthalene      | 13.91 | 1000       | J     |
| 13.            | Substituted Naphthalene      | 14.08 | 1200       | J     |
| 14.            | Substituted Naphthalene      | 14.76 | 780        | J     |
| 15.            | Substituted Naphthalene      | 15.18 | 710        | J     |
| 16.            | Unknown PAH                  | 15.96 | 1800       | J     |
| 17.            | Unknown                      | 19.53 | 780        | J     |
| 18.            | Unknown                      | 20.25 | 1600       | J     |
| 19.            | Unknown PAH                  | 21.18 | 860        | J     |
| 20.            | Unknown                      | 21.66 | 1300       | J     |
| 21.            | Unknown                      | 21.89 | 840        | J     |
| 22.            | Unknown                      | 22.56 | 4000       | J     |
| 23.            | benzo (e) pyrene             | 24.87 | 1100       | J     |
| 24.            |                              |       |            |       |
| 25.            |                              |       |            |       |
| 26.            |                              |       |            |       |
| 27.            |                              |       |            |       |
| 28.            |                              |       |            |       |
| 29.            |                              |       |            |       |
| 30.            |                              |       |            |       |

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

SB-14 (5-7) DL

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-5DL

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED82118

Level: (low/med) LOW

Date Received: 08/10/00

% Moisture: 28 decanted: (Y/N) N

Date Extracted: 08/14/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 10.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:

(ug/L or ug/Kg) ug/kg Q

| CAS NO.  | COMPOUND                     | CONCENTRATION UNITS: | Q  |
|----------|------------------------------|----------------------|----|
| 108-95-2 | Phenol                       | 4600                 | U  |
| 111-44-4 | bis(2-Chloroethyl) ether     | 4600                 | U  |
| 95-57-8  | 2-Chlorophenol               | 4600                 | U  |
| 541-73-1 | 1,3-Dichlorobenzene          | 4600                 | U  |
| 106-46-7 | 1,4-Dichlorobenzene          | 4600                 | U  |
| 95-50-1  | 1,2-Dichlorobenzene          | 4600                 | U  |
| 95-48-7  | 2-Methylphenol               | 4600                 | U  |
| 108-60-1 | 2,2'-oxybis(1-Chloropropane) | 4600                 | U  |
| 106-44-5 | 4-Methylphenol               | 4600                 | U  |
| 621-64-7 | N-Nitroso-di-n-propylamine   | 4600                 | U  |
| 67-72-1  | Hexachloroethane             | 4600                 | U  |
| 98-95-3  | Nitrobenzene                 | 4600                 | U  |
| 78-59-1  | Isophorone                   | 4600                 | U  |
| 88-75-5  | 2-Nitrophenol                | 4600                 | U  |
| 105-67-9 | 2,4-Dimethylphenol           | 4600                 | U  |
| 111-91-1 | bis(2-Chloroethoxy) methane  | 4600                 | U  |
| 120-83-2 | 2,4-Dichlorophenol           | 4600                 | U  |
| 120-82-1 | 1,2,4-Trichlorobenzene       | 4600                 | U  |
| 91-20-3  | Naphthalene                  | 4600                 | U  |
| 106-47-8 | 4-Chloroaniline              | 4600                 | U  |
| 87-68-3  | Hexachlorobutadiene          | 4600                 | U  |
| 59-50-7  | 4-Chloro-3-methylphenol      | 4600                 | U  |
| 91-57-6  | 2-Methylnaphthalene          | 4600                 | U  |
| 77-47-4  | Hexachlorocyclopentadiene    | 4600                 | U  |
| 88-06-2  | 2,4,6-Trichlorophenol        | 4600                 | U  |
| 95-95-4  | 2,4,5-Trichlorophenol        | 12000                | U  |
| 91-58-7  | 2-Chloronaphthalene          | 4600                 | U  |
| 88-74-4  | 2-Nitroaniline               | 12000                | U  |
| 131-11-3 | Dimethylphthalate            | 4600                 | U  |
| 208-96-8 | Acenaphthylene               | 560                  | JD |
| 606-20-2 | 2,6-Dinitrotoluene           | 4600                 | U  |
| 99-09-2  | 3-Nitroaniline               | 12000                | U  |
| 83-32-9  | Acenaphthene                 | 8700                 | D  |

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SB-14 (5-7) DL

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-5DL

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED82118

Level: (low/med) LOW

Date Received: 08/10/00

% Moisture: 28 decanted: (Y/N) N

Date Extracted: 08/14/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 10.0

GPC Cleanup: (Y/N) Y pH: 7.0

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg Q

|                |                            |       |    |
|----------------|----------------------------|-------|----|
| 51-28-5-----   | 2,4-Dinitrophenol          | 12000 | U  |
| 100-02-7-----  | 4-Nitrophenol              | 12000 | U  |
| 132-64-9-----  | Dibenzofuran               | 5800  | D  |
| 121-14-2-----  | 2,4-Dinitrotoluene         | 4600  | U  |
| 84-66-2-----   | Diethylphthalate           | 4600  | U  |
| 86-73-7-----   | Fluorene                   | 7000  | D  |
| 7005-72-3----- | 4-Chlorophenyl-phenylether | 4600  | U  |
| 100-01-6-----  | 4-Nitroaniline             | 12000 | U  |
| 534-52-1-----  | 4,6-Dinitro-2-methylphenol | 12000 | U  |
| 86-30-6-----   | N-Nitrosodiphenylamine     | 4600  | U  |
| 101-55-3-----  | 4-Bromophenyl-phenylether  | 4600  | U  |
| 118-74-1-----  | Hexachlorobenzene          | 4600  | U  |
| 87-86-5-----   | Pentachlorophenol          | 12000 | U  |
| 85-01-8-----   | Phenanthrene               | 23000 | D  |
| 120-12-7-----  | Anthracene                 | 6700  | D  |
| 86-74-8-----   | Carbazole                  | 3000  | JD |
| 84-74-2-----   | Di-n-butylphthalate        | 4600  | U  |
| 206-44-0-----  | Fluoranthene               | 10000 | D  |
| 129-00-0-----  | Pyrene                     | 15000 | D  |
| 85-68-7-----   | Butylbenzylphthalate       | 4600  | U  |
| 56-55-3-----   | Benzo(a)anthracene         | 3700  | JD |
| 91-94-1-----   | 3,3'-Dichlorobenzidine     | 4600  | U  |
| 218-01-9-----  | Chrysene                   | 3000  | JD |
| 117-81-7-----  | bis(2-Ethylhexyl)phthalate | 4600  | U  |
| 117-84-0-----  | Di-n-octylphthalate        | 4600  | U  |
| 205-99-2-----  | Benzo(b)fluoranthene       | 3600  | JD |
| 207-08-9-----  | Benzo(k)fluoranthene       | 1000  | JD |
| 50-32-8-----   | Benzo(a)pyrene             | 3100  | JD |
| 193-39-5-----  | Indeno(1,2,3-cd)pyrene     | 1500  | JD |
| 53-70-3-----   | Dibenzo(a,h)anthracene     | 4600  | U  |
| 191-24-2-----  | Benzo(g,h,i)perylene       | 1300  | JD |

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SB-14 (5-7) DL

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-5 DL

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: ED82118

Level: (low/med) LOW

Date Received: 08/10/00

% Moisture: 28 decanted: (Y/N) N

Date Extracted: 08/14/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 10.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:

(ug/L or ug/Kg) ug/kg

Number TICS found: 22

| CAS NUMBER     | COMPOUND NAME                | RT    | EST. CONC. | Q   |
|----------------|------------------------------|-------|------------|-----|
| 1.             | Unknown                      | 3.45  | 2000       | JB  |
| 2.             | Unknown                      | 4.43  | 72000      | J   |
| 3. 000123-42-2 | 2-Pentanone, 4-hydroxy-4-met | 5.31  | 5300       | JBA |
| 4.             | Unknown                      | 7.32  | 1900       | JBA |
| 5.             | Substituted Naphthalene      | 12.61 | 3800       | J   |
| 6.             | Substituted Naphthalene      | 13.84 | 2900       | J   |
| 7.             | Substituted Naphthalene      | 14.06 | 1700       | J   |
| 8. 000132-65-0 | Dibenzothiophene             | 17.23 | 2000       | J   |
| 9.             | Unknown PAH                  | 18.48 | 1800       | J   |
| 10.            | Unknown PAH                  | 18.53 | 2200       | J   |
| 11.            | Unknown PAH                  | 18.67 | 3500       | J   |
| 12.            | Unknown PAH                  | 18.73 | 1700       | J   |
| 13.            | Unknown                      | 19.00 | 1700       | J   |
| 14.            | Unknown                      | 19.21 | 4600       | J   |
| 15.            | Unknown                      | 19.45 | 22000      | J   |
| 16.            | Unknown                      | 19.67 | 4800       | J   |
| 17.            | Unknown                      | 20.05 | 4600       | J   |
| 18.            | Unknown                      | 20.18 | 4100       | J   |
| 19.            | Unknown PAH                  | 21.00 | 1800       | J   |
| 20.            | Unknown PAH                  | 21.13 | 2000       | J   |
| 21.            | Unknown                      | 22.42 | 2000       | J   |
| 22.            | Unknown                      | 24.63 | 1800       | J   |
| 23.            | benzo(e) pyrene              | 24.85 | 3100       | J   |
| 24.            |                              |       |            |     |
| 25.            |                              |       |            |     |
| 26.            |                              |       |            |     |
| 27.            |                              |       |            |     |
| 28.            |                              |       |            |     |
| 29.            |                              |       |            |     |
| 30.            |                              |       |            |     |



1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SB-15 (4-6)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-4

Sample wt/vol: 30.2 (g/mL) g

Lab File ID: ED82109

Level: (low/med) LOW

Date Received: 08/10/00

% Moisture: 29 decanted: (Y/N) N

Date Extracted: 08/14/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|               |                              |      |   |
|---------------|------------------------------|------|---|
| 108-95-2----- | Phenol                       | 470  | U |
| 111-44-4----- | bis(2-Chloroethyl) ether     | 470  | U |
| 95-57-8-----  | 2-Chlorophenol               | 470  | U |
| 541-73-1----- | 1,3-Dichlorobenzene          | 470  | U |
| 106-46-7----- | 1,4-Dichlorobenzene          | 470  | U |
| 95-50-1-----  | 1,2-Dichlorobenzene          | 470  | U |
| 95-48-7-----  | 2-Methylphenol               | 470  | U |
| 108-60-1----- | 2,2'-oxybis(1-Chloropropane) | 470  | U |
| 106-44-5----- | 4-Methylphenol               | 470  | U |
| 621-64-7----- | N-Nitroso-di-n-propylamine   | 470  | U |
| 67-72-1-----  | Hexachloroethane             | 470  | U |
| 98-95-3-----  | Nitrobenzene                 | 470  | U |
| 78-59-1-----  | Isophorone                   | 470  | U |
| 88-75-5-----  | 2-Nitrophenol                | 470  | U |
| 105-67-9----- | 2,4-Dimethylphenol           | 470  | U |
| 111-91-1----- | bis(2-Chloroethoxy) methane  | 470  | U |
| 120-83-2----- | 2,4-Dichlorophenol           | 470  | U |
| 120-82-1----- | 1,2,4-Trichlorobenzene       | 470  | U |
| 91-20-3-----  | Naphthalene                  | 2100 |   |
| 106-47-8----- | 4-Chloroaniline              | 470  | U |
| 87-68-3-----  | Hexachlorobutadiene          | 470  | U |
| 59-50-7-----  | 4-Chloro-3-methylphenol      | 470  | U |
| 91-57-6-----  | 2-Methylnaphthalene          | 160  | J |
| 77-47-4-----  | Hexachlorocyclopentadiene    | 470  | U |
| 88-06-2-----  | 2,4,6-Trichlorophenol        | 470  | U |
| 95-95-4-----  | 2,4,5-Trichlorophenol        | 1200 | U |
| 91-58-7-----  | 2-Chloronaphthalene          | 470  | U |
| 88-74-4-----  | 2-Nitroaniline               | 1200 | U |
| 131-11-3----- | Dimethylphthalate            | 470  | U |
| 208-96-8----- | Acenaphthylene               | 48   | J |
| 606-20-2----- | 2,6-Dinitrotoluene           | 470  | U |
| 99-09-2-----  | 3-Nitroaniline               | 1200 | U |
| 83-32-9-----  | Acenaphthene                 | 260  | J |

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SB-15 (4-6)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-4

Sample wt/vol: 30.2 (g/mL) g

Lab File ID: ED82109

Level: (low/med) LOW

Date Received: 08/10/00

% Moisture: 29 decanted: (Y/N) N

Date Extracted: 08/14/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CAS NO.

COMPOUND

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg

Q

|                |                            |      |   |
|----------------|----------------------------|------|---|
| 51-28-5-----   | 2,4-Dinitrophenol          | 1200 | U |
| 100-02-7-----  | 4-Nitrophenol              | 1200 | U |
| 132-64-9-----  | Dibenzofuran               | 84   | J |
| 121-14-2-----  | 2,4-Dinitrotoluene         | 470  | U |
| 84-66-2-----   | Diethylphthalate           | 470  | U |
| 86-73-7-----   | Fluorene                   | 150  | J |
| 7005-72-3----- | 4-Chlorophenyl-phenylether | 470  | U |
| 100-01-6-----  | 4-Nitroaniline             | 1200 | U |
| 534-52-1-----  | 4,6-Dinitro-2-methylphenol | 1200 | U |
| 86-30-6-----   | N-Nitrosodiphenylamine     | 470  | U |
| 101-55-3-----  | 4-Bromophenyl-phenylether  | 470  | U |
| 118-74-1-----  | Hexachlorobenzene          | 470  | U |
| 87-86-5-----   | Pentachlorophenol          | 1200 | U |
| 85-01-8-----   | Phenanthrene               | 370  | J |
| 120-12-7-----  | Anthracene                 | 84   | J |
| 86-74-8-----   | Carbazole                  | 230  | J |
| 84-74-2-----   | Di-n-butylphthalate        | 470  | U |
| 206-44-0-----  | Fluoranthene               | 670  |   |
| 129-00-0-----  | Pyrene                     | 1100 |   |
| 85-68-7-----   | Butylbenzylphthalate       | 470  | U |
| 56-55-3-----   | Benzo (a) anthracene       | 420  | J |
| 91-94-1-----   | 3,3'-Dichlorobenzidine     | 470  | U |
| 218-01-9-----  | Chrysene                   | 480  |   |
| 117-81-7-----  | bis(2-Ethylhexyl)phthalate | 470  | U |
| 117-84-0-----  | Di-n-octylphthalate        | 470  | U |
| 205-99-2-----  | Benzo (b) fluoranthene     | 710  |   |
| 207-08-9-----  | Benzo (k) fluoranthene     | 230  | J |
| 50-32-8-----   | Benzo (a) pyrene           | 500  |   |
| 193-39-5-----  | Indeno (1,2,3-cd) pyrene   | 290  | J |
| 53-70-3-----   | Dibenzo (a,h) anthracene   | 88   | J |
| 191-24-2-----  | Benzo (g,h,i) perylene     | 260  | J |

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SB-15 (4-6)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-4

Sample wt/vol: 30.2 (g/mL) g

Lab File ID: ED82109

Level: (low/med) LOW

Date Received: 08/10/00

% Moisture: 29 decanted: (Y/N) N

Date Extracted: 08/14/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

Number TICS found: 5

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg

| CAS NUMBER     | COMPOUND NAME             | RT    | EST. CONC. | Q   |
|----------------|---------------------------|-------|------------|-----|
| 1.             | Unknown                   | 3.47  | 1200       | JB  |
| 2. 000141-79-7 | 3-Penten-2-one, 4-methyl- | 4.57  | 55000      | JBA |
| 3.             | Unknown                   | 5.39  | 3100       | J   |
| 4.             | Unknown                   | 7.36  | 890        | J   |
| 5.             | Unknown                   | 9.07  | 650        | J   |
| 6.             | benzo(e) pyrene           | 24.85 | 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.            |                           |       |            |     |

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

SB-15 (4-6) R

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-4R E

Sample wt/vol: 30.2 (g/mL) g

Lab File ID: ED82115

Level: (low/med) LOW

Date Received: 08/10/00

% Moisture: 29 decanted: (Y/N) N

Date Extracted: 08/14/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:

(ug/L or ug/Kg) ug/kg Q

| CAS NO.  | COMPOUND                     | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) ug/kg | Q |
|----------|------------------------------|---|---|
| 108-95-2 | Phenol                       | 470   | U |
| 111-44-4 | bis(2-Chloroethyl) ether     | 470   | U |
| 95-57-8  | 2-Chlorophenol               | 470   | U |
| 541-73-1 | 1,3-Dichlorobenzene          | 470   | U |
| 106-46-7 | 1,4-Dichlorobenzene          | 470   | U |
| 95-50-1  | 1,2-Dichlorobenzene          | 470   | U |
| 95-48-7  | 2-Methylphenol               | 470   | U |
| 108-60-1 | 2,2'-oxybis(1-Chloropropane) | 470   | U |
| 106-44-5 | 4-Methylphenol               | 470   | U |
| 621-64-7 | N-Nitroso-di-n-propylamine   | 470   | U |
| 67-72-1  | Hexachloroethane             | 470   | U |
| 98-95-3  | Nitrobenzene                 | 470   | U |
| 78-59-1  | Isophorone                   | 470   | U |
| 88-75-5  | 2-Nitrophenol                | 470   | U |
| 105-67-9 | 2,4-Dimethylphenol           | 470   | U |
| 111-91-1 | bis(2-Chloroethoxy) methane  | 470   | U |
| 120-83-2 | 2,4-Dichlorophenol           | 470   | U |
| 120-82-1 | 1,2,4-Trichlorobenzene       | 470   | U |
| 91-20-3  | Naphthalene                  | 2100  |   |
| 106-47-8 | 4-Chloroaniline              | 470   | U |
| 87-68-3  | Hexachlorobutadiene          | 470   | U |
| 59-50-7  | 4-Chloro-3-methylphenol      | 470   | U |
| 91-57-6  | 2-Methylnaphthalene          | 150   | J |
| 77-47-4  | Hexachlorocyclopentadiene    | 470   | U |
| 88-06-2  | 2,4,6-Trichlorophenol        | 470   | U |
| 95-95-4  | 2,4,5-Trichlorophenol        | 1200  | U |
| 91-58-7  | 2-Chloronaphthalene          | 470   | U |
| 88-74-4  | 2-Nitroaniline               | 1200  | U |
| 131-11-3 | Dimethylphthalate            | 470   | U |
| 208-96-8 | Acenaphthylene               | 54  | J |
| 606-20-2 | 2,6-Dinitrotoluene           | 470   | U |
| 99-09-2  | 3-Nitroaniline               | 1200  | U |
| 83-32-9  | Acenaphthene                 | 260   | J |

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SB-15 (4-6) R

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-4R

Sample wt/vol: 30.2 (g/mL) g

Lab File ID: ED82115

Level: (low/med) LOW

Date Received: 08/10/00

% Moisture: 29 decanted: (Y/N) N

Date Extracted: 08/14/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg Q

|                |                            |      |   |
|----------------|----------------------------|------|---|
| 51-28-5-----   | 2,4-Dinitrophenol          | 1200 | U |
| 100-02-7-----  | 4-Nitrophenol              | 1200 | U |
| 132-64-9-----  | Dibenzofuran               | 78   | J |
| 121-14-2-----  | 2,4-Dinitrotoluene         | 470  | U |
| 84-66-2-----   | Diethylphthalate           | 470  | U |
| 86-73-7-----   | Fluorene                   | 150  | J |
| 7005-72-3----- | 4-Chlorophenyl-phenylether | 470  | U |
| 100-01-6-----  | 4-Nitroaniline             | 1200 | U |
| 534-52-1-----  | 4,6-Dinitro-2-methylphenol | 1200 | U |
| 86-30-6-----   | N-Nitrosodiphenylamine     | 470  | U |
| 101-55-3-----  | 4-Bromophenyl-phenylether  | 470  | U |
| 118-74-1-----  | Hexachlorobenzene          | 470  | U |
| 87-86-5-----   | Pentachlorophenol          | 1200 | U |
| 85-01-8-----   | Phenanthrene               | 370  | J |
| 120-12-7-----  | Anthracene                 | 83   | J |
| 86-74-8-----   | Carbazole                  | 210  | J |
| 84-74-2-----   | Di-n-butylphthalate        | 470  | U |
| 206-44-0-----  | Fluoranthene               | 620  |   |
| 129-00-0-----  | Pyrene                     | 1200 | J |
| 85-68-7-----   | Butylbenzylphthalate       | 470  | U |
| 56-55-3-----   | Benzo(a)anthracene         | 420  | J |
| 91-94-1-----   | 3,3'-Dichlorobenzidine     | 470  | U |
| 218-01-9-----  | Chrysene                   | 460  | J |
| 117-81-7-----  | bis(2-Ethylhexyl)phthalate | 470  | U |
| 117-84-0-----  | Di-n-octylphthalate        | 470  | U |
| 205-99-2-----  | Benzo(b)fluoranthene       | 790  | J |
| 207-08-9-----  | Benzo(k)fluoranthene       | 220  | J |
| 50-32-8-----   | Benzo(a)pyrene             | 640  | J |
| 193-39-5-----  | Indeno(1,2,3-cd)pyrene     | 400  | J |
| 53-70-3-----   | Dibenzo(a,h)anthracene     | 120  | J |
| 191-24-2-----  | Benzo(g,h,i)perylene       | 390  | J |

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

SB-15 (4-6) R

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-4 R

Sample wt/vol: 30.2 (g/mL) g

Lab File ID: ED82115

Level: (low/med) LOW

Date Received: 08/10/00

% Moisture: 29 decanted: (Y/N) N

Date Extracted: 08/14/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

Number TICS found: 4

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg

| CAS NUMBER     | COMPOUND NAME                | RT    | EST. CONC. | Q   |
|----------------|------------------------------|-------|------------|-----|
| 1.             | Unknown                      | 3.47  | 1200       | JB  |
| 2. 000141-79-7 | 3-Penten-2-one, 4-methyl     | 4.56  | 55000      | JBA |
| 3. 000123-42-2 | 2-Pentanone, 4-hydroxy-4-met | 5.38  | 3000       | JBA |
| 4.             | Unknown                      | 7.35  | 890        | J   |
| 5.             | benzole) pyrene              | 24.85 | 340        | J   |
| 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.            |                              |       |            |     |

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

DUP-2

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-6

Sample wt/vol: 30.2 (g/mL) g

Lab File ID: ED82113

Level: (low/med) LOW

Date Received: 08/10/00

% Moisture: 21 decanted: (Y/N) N

Date Extracted: 08/14/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|               |                              |      |   |
|---------------|------------------------------|------|---|
| 108-95-2----- | Phenol                       | 420  | U |
| 111-44-4----- | bis(2-Chloroethyl) ether     | 420  | U |
| 95-57-8-----  | 2-Chlorophenol               | 420  | U |
| 541-73-1----- | 1,3-Dichlorobenzene          | 420  | U |
| 106-46-7----- | 1,4-Dichlorobenzene          | 420  | U |
| 95-50-1-----  | 1,2-Dichlorobenzene          | 420  | U |
| 95-48-7-----  | 2-Methylphenol               | 420  | U |
| 108-60-1----- | 2,2'-oxybis(1-Chloropropane) | 420  | U |
| 106-44-5----- | 4-Methylphenol               | 67   | J |
| 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           | 50   | J |
| 111-91-1----- | bis(2-Chloroethoxy) methane  | 420  | U |
| 120-83-2----- | 2,4-Dichlorophenol           | 420  | U |
| 120-82-1----- | 1,2,4-Trichlorobenzene       | 420  | U |
| 91-20-3-----  | Naphthalene                  | 460  |   |
| 106-47-8----- | 4-Chloroaniline              | 420  | U |
| 87-68-3-----  | Hexachlorobutadiene          | 420  | U |
| 59-50-7-----  | 4-Chloro-3-methylphenol      | 420  | U |
| 91-57-6-----  | 2-Methylnaphthalene          | 500  |   |
| 77-47-4-----  | Hexachlorocyclopentadiene    | 420  | U |
| 88-06-2-----  | 2,4,6-Trichlorophenol        | 420  | U |
| 95-95-4-----  | 2,4,5-Trichlorophenol        | 1000 | U |
| 91-58-7-----  | 2-Chloronaphthalene          | 420  | U |
| 88-74-4-----  | 2-Nitroaniline               | 1000 | U |
| 131-11-3----- | Dimethylphthalate            | 420  | U |
| 208-96-8----- | Acenaphthylene               | 1200 |   |
| 606-20-2----- | 2,6-Dinitrotoluene           | 420  | U |
| 99-09-2-----  | 3-Nitroaniline               | 1000 | U |
| 83-32-9-----  | Acenaphthene                 | 340  | J |

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

DUP-2

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-6

Sample wt/vol: 30.2 (g/mL) g

Lab File ID: ED82113

Level: (low/med) LOW

Date Received: 08/10/00

% Moisture: 21 decanted: (Y/N) N

Date Extracted: 08/14/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CAS NO.

COMPOUND

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg

Q

|                |                            |       |   |
|----------------|----------------------------|-------|---|
| 51-28-5-----   | 2,4-Dinitrophenol          | 1000  | U |
| 100-02-7-----  | 4-Nitrophenol              | 1000  | U |
| 132-64-9-----  | Dibenzofuran               | 920   |   |
| 121-14-2-----  | 2,4-Dinitrotoluene         | 420   | U |
| 84-66-2-----   | Diethylphthalate           | 420   | U |
| 86-73-7-----   | Fluorene                   | 1500  |   |
| 7005-72-3----- | 4-Chlorophenyl-phenylether | 420   | U |
| 100-01-6-----  | 4-Nitroaniline             | 1000  | U |
| 534-52-1-----  | 4,6-Dinitro-2-methylphenol | 1000  | U |
| 86-30-6-----   | N-Nitrosodiphenylamine     | 420   | U |
| 101-55-3-----  | 4-Bromophenyl-phenylether  | 420   | U |
| 118-74-1-----  | Hexachlorobenzene          | 420   | U |
| 87-86-5-----   | Pentachlorophenol          | 1000  | U |
| 85-01-8-----   | Phenanthrene               | 12000 | E |
| 120-12-7-----  | Anthracene                 | 3200  |   |
| 86-74-8-----   | Carbazole                  | 770   |   |
| 84-74-2-----   | Di-n-butylphthalate        | 420   | U |
| 206-44-0-----  | Fluoranthene               | 10000 | E |
| 129-00-0-----  | Pyrene                     | 20000 | E |
| 85-68-7-----   | Butylbenzylphthalate       | 420   | U |
| 56-55-3-----   | Benzo(a)anthracene         | 5700  | E |
| 91-94-1-----   | 3,3'-Dichlorobenzidine     | 420   | U |
| 218-01-9-----  | Chrysene                   | 4900  | E |
| 117-81-7-----  | bis(2-Ethylhexyl)phthalate | 420   | U |
| 117-84-0-----  | Di-n-octylphthalate        | 420   | U |
| 205-99-2-----  | Benzo(b)fluoranthene       | 5000  | E |
| 207-08-9-----  | Benzo(k)fluoranthene       | 1600  | J |
| 50-32-8-----   | Benzo(a)pyrene             | 4200  | E |
| 193-39-5-----  | Indeno(1,2,3-cd)pyrene     | 2200  | J |
| 53-70-3-----   | Dibenzo(a,h)anthracene     | 720   | J |
| 191-24-2-----  | Benzo(g,h,i)perylene       | 2000  | J |



1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

DUP-2

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-6

Sample wt/vol: 30.2 (g/mL) g

Lab File ID: ED82113

Level: (low/med) LOW

Date Received: 08/10/00

% Moisture: 21 decanted: (Y/N) N

Date Extracted: 08/14/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/21/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

Number TICS found: 22

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg

| CAS NUMBER      | COMPOUND NAME                    | RT    | EST. CONC. | Q   |
|-----------------|----------------------------------|-------|------------|-----|
| 1.              | Unknown                          | 3.46  | 1300       | J   |
| 2.              | Unknown                          | 3.58  | 860        | JB  |
| 3. 000141-79-7  | 3-Penten-2-one, 4-methyl-        | 4.53  | 58000      | JBA |
| 4. 000123-42-2  | 2-Pentanone, 4-hydroxy-4-methyl- | 5.36  | 3600       | JBA |
| 5.              | Unknown                          | 7.36  | 1600       | J   |
| 6.              | Unknown PAH                      | 15.94 | 740        | J   |
| 7. 001730-37-6  | 9H-Fluorene, 1-methyl-           | 16.72 | 1000       | J   |
| 8.              | Unknown                          | 17.00 | 1200       | J   |
| 9. 000486-25-9  | 9H-Fluoren-9-one                 | 17.10 | 730        | J   |
| 10.             | Unknown                          | 17.22 | 1100       | J   |
| 11. 000132-65-0 | Dibenzothiophene                 | 17.25 | 1100       | J   |
| 12.             | Unknown PAH                      | 18.50 | 2300       | J   |
| 13.             | Unknown PAH                      | 18.56 | 2800       | J   |
| 14.             | Unknown PAH                      | 18.66 | 1700       | J   |
| 15.             | Unknown PAH                      | 18.71 | 3600       | J   |
| 16.             | Unknown PAH                      | 18.77 | 1500       | J   |
| 17.             | Unknown PAH                      | 19.12 | 1600       | J   |
| 18.             | Unknown PAH                      | 19.15 | 830        | J   |
| 19. 001576-67-6 | Phenanthrene, 3,6-dimethyl-      | 19.46 | 930        | J   |
| 20.             | Unknown PAH                      | 19.61 | 1600       | J   |
| 21.             | Unknown                          | 19.77 | 1400       | J   |
| 22.             | Unknown PAH                      | 20.09 | 2300       | J   |
| 23.             | Benzo(e)pyrene                   | 24.66 | 2100       | J   |
| 24.             |                                  |       |            |     |
| 25.             |                                  |       |            |     |
| 26.             |                                  |       |            |     |
| 27.             |                                  |       |            |     |
| 28.             |                                  |       |            |     |
| 29.             |                                  |       |            |     |
| 30.             |                                  |       |            |     |

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

DUP-2 DL

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-6DL

Sample wt/vol: 30.2 (g/mL) g

Lab File ID: ED82222

Level: (low/med) LOW

Date Received: 08/10/00

% Moisture: 21 decanted: (Y/N) N

Date Extracted: 08/14/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/22/00

Injection Volume: 2.0 (uL)

Dilution Factor: 10.0

GPC Cleanup: (Y/N) Y pH: 7.0

|         |          |   |   |
|---------|----------|---|---|
| CAS NO. | COMPOUND | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) ug/kg | Q |
|---------|----------|---|---|

|               |                              |       |    |
|---------------|------------------------------|-------|----|
| 108-95-2----- | Phenol                       | 4200  | U  |
| 111-44-4----- | bis(2-Chloroethyl) ether     | 4200  | U  |
| 95-57-8-----  | 2-Chlorophenol               | 4200  | U  |
| 541-73-1----- | 1,3-Dichlorobenzene          | 4200  | U  |
| 106-46-7----- | 1,4-Dichlorobenzene          | 4200  | U  |
| 95-50-1-----  | 1,2-Dichlorobenzene          | 4200  | U  |
| 95-48-7-----  | 2-Methylphenol               | 4200  | U  |
| 108-60-1----- | 2,2'-oxybis(1-Chloropropane) | 4200  | U  |
| 106-44-5----- | 4-Methylphenol               | 4200  | U  |
| 621-64-7----- | N-Nitroso-di-n-propylamine   | 4200  | U  |
| 67-72-1-----  | Hexachloroethane             | 4200  | U  |
| 98-95-3-----  | Nitrobenzene                 | 4200  | U  |
| 78-59-1-----  | Isophorone                   | 4200  | U  |
| 88-75-5-----  | 2-Nitrophenol                | 4200  | U  |
| 105-67-9----- | 2,4-Dimethylphenol           | 4200  | U  |
| 111-91-1----- | bis(2-Chloroethoxy) methane  | 4200  | U  |
| 120-83-2----- | 2,4-Dichlorophenol           | 4200  | U  |
| 120-82-1----- | 1,2,4-Trichlorobenzene       | 4200  | U  |
| 91-20-3-----  | Naphthalene                  | 510   | JD |
| 106-47-8----- | 4-Chloroaniline              | 4200  | U  |
| 87-68-3-----  | Hexachlorobutadiene          | 4200  | U  |
| 59-50-7-----  | 4-Chloro-3-methylphenol      | 4200  | U  |
| 91-57-6-----  | 2-Methylnaphthalene          | 520   | JD |
| 77-47-4-----  | Hexachlorocyclopentadiene    | 4200  | U  |
| 88-06-2-----  | 2,4,6-Trichlorophenol        | 4200  | U  |
| 95-95-4-----  | 2,4,5-Trichlorophenol        | 10000 | U  |
| 91-58-7-----  | 2-Chloronaphthalene          | 4200  | U  |
| 88-74-4-----  | 2-Nitroaniline               | 10000 | U  |
| 131-11-3----- | Dimethylphthalate            | 4200  | U  |
| 208-96-8----- | Acenaphthylene               | 1300  | JD |
| 606-20-2----- | 2,6-Dinitrotoluene           | 4200  | U  |
| 99-09-2-----  | 3-Nitroaniline               | 10000 | U  |
| 83-32-9-----  | Acenaphthene                 | 4200  | U  |

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

DUP-2 DL

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-6DL

Sample wt/vol: 30.2 (g/mL) g

Lab File ID: ED82222

Level: (low/med) LOW

Date Received: 08/10/00

% Moisture: 21 decanted: (Y/N) N

Date Extracted: 08/14/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/22/00

Injection Volume: 2.0 (uL)

Dilution Factor: 10.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|                |                            |       |    |
|----------------|----------------------------|-------|----|
| 51-28-5-----   | 2,4-Dinitrophenol          | 10000 | U  |
| 100-02-7-----  | 4-Nitrophenol              | 10000 | U  |
| 132-64-9-----  | Dibenzofuran               | 1100  | JD |
| 121-14-2-----  | 2,4-Dinitrotoluene         | 4200  | U  |
| 84-66-2-----   | Diethylphthalate           | 4200  | U  |
| 86-73-7-----   | Fluorene                   | 1800  | JD |
| 7005-72-3----- | 4-Chlorophenyl-phenylether | 4200  | U  |
| 100-01-6-----  | 4-Nitroaniline             | 10000 | U  |
| 534-52-1-----  | 4,6-Dinitro-2-methylphenol | 10000 | U  |
| 86-30-6-----   | N-Nitrosodiphenylamine     | 4200  | U  |
| 101-55-3-----  | 4-Bromophenyl-phenylether  | 4200  | U  |
| 118-74-1-----  | Hexachlorobenzene          | 4200  | U  |
| 87-86-5-----   | Pentachlorophenol          | 10000 | U  |
| 85-01-8-----   | Phenanthrene               | 10000 | D  |
| 120-12-7-----  | Anthracene                 | 3600  | JD |
| 86-74-8-----   | Carbazole                  | 920   | JD |
| 84-74-2-----   | Di-n-butylphthalate        | 4200  | U  |
| 206-44-0-----  | Fluoranthene               | 12000 | D  |
| 129-00-0-----  | Pyrene                     | 12000 | D  |
| 85-68-7-----   | Butylbenzylphthalate       | 4200  | U  |
| 56-55-3-----   | Benzo(a)anthracene         | 6400  | D  |
| 91-94-1-----   | 3,3'-Dichlorobenzidine     | 4200  | U  |
| 218-01-9-----  | Chrysene                   | 5500  | D  |
| 117-81-7-----  | bis(2-Ethylhexyl)phthalate | 4200  | U  |
| 117-84-0-----  | Di-n-octylphthalate        | 4200  | U  |
| 205-99-2-----  | Benzo(b)fluoranthene       | 7300  | D  |
| 207-08-9-----  | Benzo(k)fluoranthene       | 2600  | JD |
| 50-32-8-----   | Benzo(a)pyrene             | 5600  | D  |
| 193-39-5-----  | Indeno(1,2,3-cd)pyrene     | 1700  | JD |
| 53-70-3-----   | Dibenzo(a,h)anthracene     | 530   | JD |
| 191-24-2-----  | Benzo(g,h,i)perylene       | 1600  | JD |

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NO.

DUP-2 DL

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-6 DL

Sample wt/vol: 30.2 (g/mL) g

Lab File ID: ED82222

Level: (low/med) LOW

Date Received: 08/10/00

% Moisture: 21 decanted: (Y/N) N

Date Extracted: 08/14/00

Concentrated Extract Volume: 500 (uL)

Date Analyzed: 08/22/00

Injection Volume: 2.0 (uL)

Dilution Factor: 10.0

GPC Cleanup: (Y/N) Y pH: 7.0

Number TICS found: 21

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg

| CAS NUMBER     | COMPOUND NAME            | RT    | EST. CONC. | Q   |
|----------------|--------------------------|-------|------------|-----|
| 1.             | Unknown                  | 3.23  | 1600       | J   |
| 2.             | Unknown                  | 3.33  | 1400       | JB  |
| 3. 000141-79-7 | 3-Penten-2-one, 4-methyl | 4.19  | 46000      | JBA |
| 4.             | Unknown                  | 5.09  | 2900       | J   |
| 5.             | Unknown                  | 7.12  | 1700       | J   |
| 6.             | Unknown PAH              | 15.89 | 1000       | J   |
| 7.             | Unknown PAH              | 18.27 | 1900       | J   |
| 8.             | Unknown PAH              | 18.33 | 2600       | J   |
| 9.             | Unknown PAH              | 18.42 | 1200       | J   |
| 10.            | Unknown PAH              | 18.47 | 2900       | J   |
| 11.            | Unknown PAH              | 18.53 | 1900       | J   |
| 12.            | Unknown PAH              | 18.90 | 2000       | J   |
| 13.            | Unknown PAH              | 19.39 | 1400       | J   |
| 14.            | Unknown                  | 19.44 | 1200       | J   |
| 15.            | Unknown PAH              | 19.56 | 900        | J   |
| 16.            | Unknown PAH              | 19.85 | 1700       | J   |
| 17.            | Unknown PAH              | 20.80 | 1600       | J   |
| 18.            | Unknown PAH              | 20.93 | 1300       | J   |
| 19.            | Unknown PAH              | 22.09 | 980        | J   |
| 20.            | Unknown PAH              | 23.32 | 910        | J   |
| 21.            | Unknown PAH              | 24.48 | 2300       | J   |
| 22.            | benzo(e)pyrene           | 24.70 | 3000       | J   |
| 23.            |                          |       |            |     |
| 24.            |                          |       |            |     |
| 25.            |                          |       |            |     |
| 26.            |                          |       |            |     |
| 27.            |                          |       |            |     |
| 28.            |                          |       |            |     |
| 29.            |                          |       |            |     |
| 30.            |                          |       |            |     |

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

SB-6 (16-18)

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62574-1

Sample wt/vol: 30 (g/mL) g

Lab File ID: HP13A\0817C021

% Moisture: 17 decanted: (Y/N) N

Date Received: 08/08/00

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 08/10/00

Concentrated Extract Volume: 5000 (uL)

Date Analyzed: 08/18/00

Injection Volume: 1.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

Sulfur Cleanup: (Y/N) N

CAS NO.

COMPOUND

CONCENTRATION UNITS:

(ug/L or ug/Kg) ug/kg

Q

|                 |                     |     |   |
|-----------------|---------------------|-----|---|
| 319-84-6-----   | alpha-BHC           | 2.0 | U |
| 319-85-7-----   | beta-BHC            | 2.0 | U |
| 319-86-8-----   | delta-BHC           | 2.0 | U |
| 58-89-9-----    | gamma-BHC (Lindane) | 2.0 | U |
| 76-44-8-----    | Heptachlor          | 2.0 | U |
| 309-00-2-----   | Aldrin              | 2.0 | U |
| 1024-57-3-----  | Heptachlor epoxide  | 2.0 | U |
| 959-98-8-----   | Endosulfan I        | 2.0 | U |
| 60-57-1-----    | Dieldrin            | 4.0 | U |
| 72-55-9-----    | 4,4'-DDE            | 4.0 | U |
| 72-20-8-----    | Endrin              | 4.0 | U |
| 332-13-659----- | Endosulfan II       | 4.0 | U |
| 72-54-8-----    | 4,4'-DDD            | 4.0 | U |
| 103-10-78-----  | Endosulfan sulfate  | 4.0 | U |
| 50-29-3-----    | 4,4'-DDT            | 4.0 | U |
| 72-43-5-----    | Methoxychlor        | 20  | U |
| 53494-70-5----- | Endrin ketone       | 4.0 | U |
| 7421-36-3-----  | Endrin aldehyde     | 4.0 | U |
| 5103-71-9-----  | alpha-Chlordane     | 2.0 | U |
| 5103-7-42-----  | gamma-Chlordane     | 2.0 | U |
| 8001-35-2-----  | Toxaphene           | 200 | U |
| 12674-11-2----- | Aroclor-1016        | 40  | U |
| 11104-28-2----- | Aroclor-1221        | 80  | U |
| 11141-16-5----- | Aroclor-1232        | 40  | U |
| 53469-21-9----- | Aroclor-1242        | 40  | U |
| 12672-29-6----- | Aroclor-1248        | 40  | U |
| 11097-69-1----- | Aroclor-1254        | 40  | U |
| 11096-82-5----- | Aroclor-1260        | 40  | U |

FORM I PEST

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

SB-6 (18-20)

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62574-4

Sample wt/vol: 30 (g/mL) g

Lab File ID: HP13A\0817C024

% Moisture: 18 decanted: (Y/N) N

Date Received: 08/08/00

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 08/10/00

Concentrated Extract Volume: 5000 (uL)

Date Analyzed: 08/18/00

Injection Volume: 1.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

Sulfur Cleanup: (Y/N) N

| CAS NO. | COMPOUND | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) ug/kg | Q |
|---------|----------|---|---|
|---------|----------|---|---|

|                 |                     |     |   |
|-----------------|---------------------|-----|---|
| 319-84-6-----   | alpha-BHC           | 1.7 | U |
| 319-85-7-----   | beta-BHC            | 1.7 | U |
| 319-86-8-----   | delta-BHC           | 1.7 | U |
| 58-89-9-----    | gamma-BHC (Lindane) | 1.7 | U |
| 76-44-8-----    | Heptachlor          | 1.7 | U |
| 309-00-2-----   | Aldrin              | 1.7 | U |
| 1024-57-3-----  | Heptachlor epoxide  | 1.7 | U |
| 959-98-8-----   | Endosulfan I        | 1.7 | U |
| 60-57-1-----    | Dieldrin            | 3.3 | U |
| 72-55-9-----    | 4,4'-DDE            | 3.3 | U |
| 72-20-8-----    | Endrin              | 3.3 | U |
| 332-13-659----- | Endosulfan II       | 3.3 | U |
| 72-54-8-----    | 4,4'-DDD            | 3.3 | U |
| 103-10-78-----  | Endosulfan sulfate  | 3.3 | U |
| 50-29-3-----    | 4,4'-DDT            | 3.3 | U |
| 72-43-5-----    | Methoxychlor        | 17  | U |
| 53494-70-5----- | Endrin ketone       | 3.3 | U |
| 7421-36-3-----  | Endrin aldehyde     | 3.3 | U |
| 5103-71-9-----  | alpha-Chlordane     | 1.7 | U |
| 5103-7-42-----  | gamma-Chlordane     | 1.7 | U |
| 8001-35-2-----  | Toxaphene           | 170 | U |
| 12674-11-2----- | Aroclor-1016        | 33  | U |
| 11104-28-2----- | Aroclor-1221        | 67  | U |
| 11141-16-5----- | Aroclor-1232        | 33  | U |
| 53469-21-9----- | Aroclor-1242        | 33  | U |
| 12672-29-6----- | Aroclor-1248        | 33  | U |
| 11097-69-1----- | Aroclor-1254        | 33  | U |
| 11096-82-5----- | Aroclor-1260        | 33  | U |

FORM I PEST

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SB-8 (6.5-8)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-1

Sample wt/vol: 30.2 (g/mL) g

Lab File ID: HP13A\0817C026

% Moisture: 75 decanted: (Y/N) N

Date Received: 08/10/00

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 08/14/00

Concentrated Extract Volume: 5000 (uL)

Date Analyzed: 08/18/00

Injection Volume: 1.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg Q

|                 |                     |     |   |
|-----------------|---------------------|-----|---|
| 319-84-6-----   | alpha-BHC           | 1.6 | U |
| 319-85-7-----   | beta-BHC            | 1.6 | U |
| 319-86-8-----   | delta-BHC           | 1.6 | U |
| 58-89-9-----    | gamma-BHC (Lindane) | 1.6 | U |
| 76-44-8-----    | Heptachlor          | 1.6 | U |
| 309-00-2-----   | Aldrin              | 1.6 | U |
| 1024-57-3-----  | Heptachlor epoxide  | 1.6 | U |
| 959-98-8-----   | Endosulfan I        | 1.6 | U |
| 60-57-1-----    | Dieldrin            | 3.3 | U |
| 72-55-9-----    | 4,4'-DDE            | 3.3 | U |
| 72-20-8-----    | Endrin              | 3.3 | U |
| 332-13-659----- | Endosulfan II       | 3.3 | U |
| 72-54-8-----    | 4,4'-DDD            | 3.3 | U |
| 103-10-78-----  | Endosulfan sulfate  | 3.3 | U |
| 50-29-3-----    | 4,4'-DDT            | 3.3 | U |
| 72-43-5-----    | Methoxychlor        | 16  | U |
| 53494-70-5----- | Endrin ketone       | 3.3 | U |
| 7421-36-3-----  | Endrin aldehyde     | 3.3 | U |
| 5103-71-9-----  | alpha-Chlordane     | 1.6 | U |
| 5103-7-42-----  | gamma-Chlordane     | 1.6 | U |
| 8001-35-2-----  | Toxaphene           | 160 | U |
| 12674-11-2----- | Aroclor-1016        | 33  | U |
| 11104-28-2----- | Aroclor-1221        | 66  | U |
| 11141-16-5----- | Aroclor-1232        | 33  | U |
| 53469-21-9----- | Aroclor-1242        | 33  | U |
| 12672-29-6----- | Aroclor-1248        | 33  | U |
| 11097-69-1----- | Aroclor-1254        | 33  | U |
| 11096-82-5----- | Aroclor-1260        | 33  | U |

FORM I PEST

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

SB-10 (4-7)

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-2

Sample wt/vol: 30.3 (g/mL) g

Lab File ID: HP13A\0817C027

% Moisture: 18 decanted: (Y/N) N

Date Received: 08/10/00

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 08/14/00

Concentrated Extract Volume: 5000 (uL)

Date Analyzed: 08/18/00

Injection Volume: 1.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

Sulfur Cleanup: (Y/N) N

CAS NO.

COMPOUND

CONCENTRATION UNITS:

(ug/L or ug/Kg) ug/kg

Q

|                 |                     |         |   |
|-----------------|---------------------|---------|---|
| 319-84-6-----   | alpha-BHC           | 1.6     | U |
| 319-85-7-----   | beta-BHC            | 1.6     | U |
| 319-86-8-----   | delta-BHC           | 1.6     | U |
| 58-89-9-----    | gamma-BHC (Lindane) | 7.5     | U |
| 76-44-8-----    | Heptachlor          | 1.6     | U |
| 309-00-2-----   | Aldrin              | 1.6     | U |
| 1024-57-3-----  | Heptachlor epoxide  | 1.6     | U |
| 959-98-8-----   | Endosulfan I        | 1.6     | U |
| 60-57-1-----    | Dieldrin            | 3.3     | U |
| 72-55-9-----    | 4,4'-DDE            | 80 120  | U |
| 72-20-8-----    | Endrin              | 3.3     | U |
| 332-13-659----- | Endosulfan II       | 3.3     | U |
| 72-54-8-----    | 4,4'-DDD            | 3.3     | U |
| 103-10-78-----  | Endosulfan sulfate  | 3.3     | U |
| 50-29-3-----    | 4,4'-DDT            | 3.3     | U |
| 72-43-5-----    | Methoxychlor        | 16      | U |
| 53494-70-5----- | Endrin ketone       | 230 300 | U |
| 7421-36-3-----  | Endrin aldehyde     | 3.3     | U |
| 5103-71-9-----  | alpha-Chlordane     | 1.6     | U |
| 5103-7-42-----  | gamma-Chlordane     | 1.6     | U |
| 8001-35-2-----  | Toxaphene           | 160     | U |
| 12674-11-2----- | Aroclor-1016        | 33      | U |
| 11104-28-2----- | Aroclor-1221        | 66      | U |
| 11141-16-5----- | Aroclor-1232        | 33      | U |
| 53469-21-9----- | Aroclor-1242        | 33      | U |
| 12672-29-6----- | Aroclor-1248        | 33      | U |
| 11097-69-1----- | Aroclor-1254        | 33      | U |
| 11096-82-5----- | Aroclor-1260        | 33      | U |

FORM I PEST

Results may be biased high due to matrix-interference.



1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SB-10 (4-7) DL

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-2DL

Sample wt/vol: 30.3 (g/mL) g

Lab File ID: HP13A\0817C029

% Moisture: 18 decanted: (Y/N) N

Date Received: 08/10/00

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 08/14/00

Concentrated Extract Volume: 5000 (uL)

Date Analyzed: 08/18/00

Injection Volume: 1.0 (uL)

Dilution Factor: 10.0

GPC Cleanup: (Y/N) Y pH: 7.0

Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|                 |                     |       |        |
|-----------------|---------------------|-------|--------|
| 319-84-6-----   | alpha-BHC           | 16    | UD     |
| 319-85-7-----   | beta-BHC            | 16    | UD     |
| 319-86-8-----   | delta-BHC           | 16    | UD     |
| 58-89-9-----    | gamma-BHC (Lindane) | 16    | UD     |
| 76-44-8-----    | Heptachlor          | 16    | UD     |
| 309-00-2-----   | Aldrin              | 16    | UD     |
| 1024-57-3-----  | Heptachlor epoxide  | 16    | UD     |
| 959-98-8-----   | Endosulfan I        | 16    | UD     |
| 60-57-1-----    | Dieldrin            | 33    | UD     |
| 72-55-9-----    | 4,4'-DDE            | 80    | UD JND |
| 72-20-8-----    | Endrin              | 33    | UD     |
| 332-13-659----- | Endosulfan II       | 33    | UD     |
| 72-54-8-----    | 4,4'-DDD            | 33    | UD     |
| 103-10-78-----  | Endosulfan sulfate  | 33    | UD     |
| 50-29-3-----    | 4,4'-DDT            | 33    | UD     |
| 72-43-5-----    | Methoxychlor        | 160   | UD     |
| 53494-70-5----- | Endrin ketone       | 23033 | UD JND |
| 7421-36-3-----  | Endrin aldehyde     | 33    | UD     |
| 5103-71-9-----  | alpha-Chlordane     | 16    | UD     |
| 5103-7-42-----  | gamma-Chlordane     | 16    | UD     |
| 8001-35-2-----  | Toxaphene           | 1600  | UD     |
| 12674-11-2----- | Aroclor-1016        | 330   | UD     |
| 11104-28-2----- | Aroclor-1221        | 660   | UD     |
| 11141-16-5----- | Aroclor-1232        | 330   | UD     |
| 53469-21-9----- | Aroclor-1242        | 330   | UD     |
| 12672-29-6----- | Aroclor-1248        | 330   | UD     |
| 11097-69-1----- | Aroclor-1254        | 330   | UD     |
| 11096-82-5----- | Aroclor-1260        | 330   | UD     |

FORM I PEST

Results may be biased high due to matrix interference

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

SB-13 (4-6)

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-3

Sample wt/vol: 30.2 (g/mL) g

Lab File ID: HP13A\0817C030

% Moisture: 26 decanted: (Y/N) N

Date Received: 08/10/00

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 08/14/00

Concentrated Extract Volume: 5000 (uL)

Date Analyzed: 08/18/00

Injection Volume: 1.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|                 |                     |     |    |
|-----------------|---------------------|-----|----|
| 319-84-6-----   | alpha-BHC           | 1.6 | U  |
| 319-85-7-----   | beta-BHC            | 1.6 | U  |
| 319-86-8-----   | delta-BHC           | 1.6 | U  |
| 58-89-9-----    | gamma-BHC (Lindane) | 2.2 | JP |
| 76-44-8-----    | Heptachlor          | 1.6 | U  |
| 309-00-2-----   | Aldrin              | 1.6 | U  |
| 1024-57-3-----  | Heptachlor epoxide  | 1.6 | U  |
| 959-98-8-----   | Endosulfan I        | 1.6 | U  |
| 60-57-1-----    | Dieldrin            | 3.3 | U  |
| 72-55-9-----    | 4,4'-DDE            | 3.3 | U  |
| 72-20-8-----    | Endrin              | 3.3 | U  |
| 332-13-659----- | Endosulfan II       | 3.3 | U  |
| 72-54-8-----    | 4,4'-DDD            | 3.3 | U  |
| 103-10-78-----  | Endosulfan sulfate  | 3.3 | U  |
| 50-29-3-----    | 4,4'-DDT            | 3.3 | U  |
| 72-43-5-----    | Methoxychlor        | 16  | U  |
| 53494-70-5----- | Endrin ketone       | 3.3 | U  |
| 7421-36-3-----  | Endrin aldehyde     | 3.3 | U  |
| 5103-71-9-----  | alpha-Chlordane     | 1.6 | U  |
| 5103-7-42-----  | gamma-Chlordane     | 1.6 | U  |
| 8001-35-2-----  | Toxaphene           | 160 | U  |
| 12674-11-2----- | Aroclor-1016        | 33  | U  |
| 11104-28-2----- | Aroclor-1221        | 66  | U  |
| 11141-16-5----- | Aroclor-1232        | 33  | U  |
| 53469-21-9----- | Aroclor-1242        | 33  | U  |
| 12672-29-6----- | Aroclor-1248        | 33  | U  |
| 11097-69-1----- | Aroclor-1254        | 33  | U  |
| 11096-82-5----- | Aroclor-1260        | 33  | U  |

FORM I PEST

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SB-13 (4-6) DL

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-3DL

Sample wt/vol: 30.2 (g/mL) g

Lab File ID: HP13A\0817C031

% Moisture: 26 decanted: (Y/N) N

Date Received: 08/10/00

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 08/14/00

Concentrated Extract Volume: 5000 (uL)

Date Analyzed: 08/18/00

Injection Volume: 1.0 (uL)

Dilution Factor: 10.0 \*

GPC Cleanup: (Y/N) Y pH: 7.0

Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|                 |                     |      |    |
|-----------------|---------------------|------|----|
| 319-84-6-----   | alpha-BHC           | 16   | UD |
| 319-85-7-----   | beta-BHC            | 16   | UD |
| 319-86-8-----   | delta-BHC           | 16   | UD |
| 58-89-9-----    | gamma-BHC (Lindane) | 16   | UD |
| 76-44-8-----    | Heptachlor          | 16   | UD |
| 309-00-2-----   | Aldrin              | 16   | UD |
| 1024-57-3-----  | Heptachlor epoxide  | 16   | UD |
| 959-98-8-----   | Endosulfan I        | 16   | UD |
| 60-57-1-----    | Dieldrin            | 33   | UD |
| 72-55-9-----    | 4,4'-DDE            | 33   | UD |
| 72-20-8-----    | Endrin              | 33   | UD |
| 332-13-659----- | Endosulfan II       | 33   | UD |
| 72-54-8-----    | 4,4'-DDD            | 33   | UD |
| 103-10-78-----  | Endosulfan sulfate  | 33   | UD |
| 50-29-3-----    | 4,4'-DDT            | 33   | UD |
| 72-43-5-----    | Methoxychlor        | 160  | UD |
| 53494-70-5----- | Endrin ketone       | 33   | UD |
| 7421-36-3-----  | Endrin aldehyde     | 33   | UD |
| 5103-71-9-----  | alpha-Chlordane     | 16   | UD |
| 5103-7-42-----  | gamma-Chlordane     | 16   | UD |
| 8001-35-2-----  | Toxaphene           | 1600 | UD |
| 12674-11-2----- | Aroclor-1016        | 330  | UD |
| 11104-28-2----- | Aroclor-1221        | 660  | UD |
| 11141-16-5----- | Aroclor-1232        | 330  | UD |
| 53469-21-9----- | Aroclor-1242        | 330  | UD |
| 12672-29-6----- | Aroclor-1248        | 330  | UD |
| 11097-69-1----- | Aroclor-1254        | 330  | UD |
| 11096-82-5----- | Aroclor-1260        | 330  | UD |

FORM I PEST

\* Diluted due to matrix interference.

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

SB-14 (5-7)

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-5

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: HP13A\0817C041

% Moisture: 28 decanted: (Y/N) N

Date Received: 08/10/00

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 08/14/00

Concentrated Extract Volume: 5000 (uL)

Date Analyzed: 08/18/00

Injection Volume: 1.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|                 |                     |     |    |
|-----------------|---------------------|-----|----|
| 319-84-6-----   | alpha-BHC           | 1.7 | U  |
| 319-85-7-----   | beta-BHC            | 1.7 | U  |
| 319-86-8-----   | delta-BHC           | 9.4 |    |
| 58-89-9-----    | gamma-BHC (Lindane) | 1.7 | U  |
| 76-44-8-----    | Heptachlor          | 1.7 | U  |
| 309-00-2-----   | Aldrin              | 1.7 | U  |
| 1024-57-3-----  | Heptachlor epoxide  | 1.7 | U  |
| 959-98-8-----   | Endosulfan I        | 1.7 | U  |
| 60-57-1-----    | Dieldrin            | 3.3 | U  |
| 72-55-9-----    | 4,4'-DDE            | 5.4 |    |
| 72-20-8-----    | Endrin              | 3.3 | U  |
| 332-13-659----- | Endosulfan II       | 3.3 | U  |
| 72-54-8-----    | 4,4'-DDD            | 3.3 | U  |
| 103-10-78-----  | Endosulfan sulfate  | 3.3 | U  |
| 50-29-3-----    | 4,4'-DDT            | 3.3 | U  |
| 72-43-5-----    | Methoxychlor        | 15  | JP |
| 53494-70-5----- | Endrin ketone       | 3.3 | U  |
| 7421-36-3-----  | Endrin aldehyde     | 3.3 | U  |
| 5103-71-9-----  | alpha-Chlordane     | 1.7 | U  |
| 5103-7-42-----  | gamma-Chlordane     | 1.7 | U  |
| 8001-35-2-----  | Toxaphene           | 170 | U  |
| 12674-11-2----- | Aroclor-1016        | 33  | U  |
| 11104-28-2----- | Aroclor-1221        | 66  | U  |
| 11141-16-5----- | Aroclor-1232        | 33  | U  |
| 53469-21-9----- | Aroclor-1242        | 33  | U  |
| 12672-29-6----- | Aroclor-1248        | 33  | U  |
| 11097-69-1----- | Aroclor-1254        | 33  | U  |
| 11096-82-5----- | Aroclor-1260        | 33  | U  |

FORM I PEST

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

SB-15 (4-6)

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-4

Sample wt/vol: 30.1 (g/mL) g

Lab File ID: HP13A\0817C032

% Moisture: 29 decanted: (Y/N) N

Date Received: 08/10/00

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 08/14/00

Concentrated Extract Volume: 5000 (uL)

Date Analyzed: 08/18/00

Injection Volume: 1.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

Sulfur Cleanup: (Y/N) N

CAS NO.

COMPOUND

CONCENTRATION UNITS:

(ug/L or ug/Kg) ug/kg Q

|                 |                     |     |   |
|-----------------|---------------------|-----|---|
| 319-84-6-----   | alpha-BHC           | 1.7 | U |
| 319-85-7-----   | beta-BHC            | 1.7 | U |
| 319-86-8-----   | delta-BHC           | 1.7 | U |
| 58-89-9-----    | gamma-BHC (Lindane) | 1.7 | U |
| 76-44-8-----    | Heptachlor          | 1.7 | U |
| 309-00-2-----   | Aldrin              | 1.7 | U |
| 1024-57-3-----  | Heptachlor epoxide  | 1.7 | U |
| 959-98-8-----   | Endosulfan I        | 1.7 | U |
| 60-57-1-----    | Dieldrin            | 3.3 | U |
| 72-55-9-----    | 4,4'-DDE            | 3.3 | U |
| 72-20-8-----    | Endrin              | 3.3 | U |
| 332-13-659----- | Endosulfan II       | 3.3 | U |
| 72-54-8-----    | 4,4'-DDD            | 3.3 | U |
| 103-10-78-----  | Endosulfan sulfate  | 3.3 | U |
| 50-29-3-----    | 4,4'-DDT            | 3.3 | U |
| 72-43-5-----    | Methoxychlor        | 17  | U |
| 53494-70-5----- | Endrin ketone       | 3.3 | U |
| 7421-36-3-----  | Endrin aldehyde     | 3.3 | U |
| 5103-71-9-----  | alpha-Chlordane     | 1.7 | U |
| 5103-7-42-----  | gamma-Chlordane     | 1.7 | U |
| 8001-35-2-----  | Toxaphene           | 170 | U |
| 12674-11-2----- | Aroclor-1016        | 33  | U |
| 11104-28-2----- | Aroclor-1221        | 66  | U |
| 11141-16-5----- | Aroclor-1232        | 33  | U |
| 53469-21-9----- | Aroclor-1242        | 33  | U |
| 12672-29-6----- | Aroclor-1248        | 33  | U |
| 11097-69-1----- | Aroclor-1254        | 33  | U |
| 11096-82-5----- | Aroclor-1260        | 33  | U |

FORM I PEST

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

DUP-2

Lab Code:

Case No.: 2

SAS No.:

SDG No.: L62574

Matrix: (soil/water) SOIL

Lab Sample ID: L62615-6

Sample wt/vol: 30 (g/mL) g

Lab File ID: HP13A\0817C043

% Moisture: 21 decanted: (Y/N) N

Date Received: 08/10/00

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 08/14/00

Concentrated Extract Volume: 5000 (uL)

Date Analyzed: 08/18/00

Injection Volume: 1.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|                 |                     |     |   |
|-----------------|---------------------|-----|---|
| 319-84-6-----   | alpha-BHC           | 1.7 | U |
| 319-85-7-----   | beta-BHC            | 1.7 | U |
| 319-86-8-----   | delta-BHC           | 1.7 | U |
| 58-89-9-----    | gamma-BHC (Lindane) | 1.7 | U |
| 76-44-8-----    | Heptachlor          | 1.7 | U |
| 309-00-2-----   | Aldrin              | 1.7 | U |
| 1024-57-3-----  | Heptachlor epoxide  | 1.7 | U |
| 959-98-8-----   | Endosulfan I        | 1.7 | U |
| 60-57-1-----    | Dieldrin            | 3.3 | U |
| 72-55-9-----    | 4,4'-DDE            | 3.3 | U |
| 72-20-8-----    | Endrin              | 3.3 | U |
| 332-13-659----- | Endosulfan II       | 3.3 | U |
| 72-54-8-----    | 4,4'-DDD            | 3.3 | U |
| 103-10-78-----  | Endosulfan sulfate  | 3.3 | U |
| 50-29-3-----    | 4,4'-DDT            | 3.3 | U |
| 72-43-5-----    | Methoxychlor        | 27  | U |
| 53494-70-5----- | Endrin ketone       | 3.3 | U |
| 7421-36-3-----  | Endrin aldehyde     | 3.3 | U |
| 5103-71-9-----  | alpha-Chlordane     | 1.7 | U |
| 5103-7-42-----  | gamma-Chlordane     | 1.7 | U |
| 8001-35-2-----  | Toxaphene           | 170 | U |
| 12674-11-2----- | Aroclor-1016        | 33  | U |
| 11104-28-2----- | Aroclor-1221        | 67  | U |
| 11141-16-5----- | Aroclor-1232        | 33  | U |
| 53469-21-9----- | Aroclor-1242        | 33  | U |
| 12672-29-6----- | Aroclor-1248        | 33  | U |
| 11097-69-1----- | Aroclor-1254        | 33  | U |
| 11096-82-5----- | Aroclor-1260        | 33  | U |

FORM I PEST

Results may be biased high due to matrix-interference!

## NYSDEC ASP

1  
INORGANIC ANALYSIS DATA SHEET

NYSDEC SAMPLE NO.

SB-6(16-18)

Lab Name: Galson Laboratories

Contract: BBL

Lab Code: 11626

Case No.:

SAS No.:

SDG No.: L62574

Matrix (soil/water): Soil

Lab Sample ID: L62574-1

Level (low/med): LOW

Date Received: 08/08/00

% Solids: 82.9

Concentration Units (ug/L or mg/kg dry weight): mg/kg

| CAS No.   | Analyte   | Concentration | C | Q | M  |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum  | 5120          | — |   | P  |
| 7440-36-0 | Antimony  | 1.2           | U | N | P  |
| 7440-38-2 | Arsenic   | 4.0           |   |   | P  |
| 7440-39-3 | Barium    | 24.0          | B |   | P  |
| 7440-41-7 | Beryllium | 0.40          | B |   | P  |
| 7440-43-9 | Cadmium   | 0.48          | U |   | P  |
| 7440-70-2 | Calcium   | 51400         |   | * | P  |
| 7440-47-3 | Chromium  | 7.6           |   |   | P  |
| 7440-48-4 | Cobalt    | 6.5           | B |   | P  |
| 7440-50-8 | Copper    | 21.4          |   | * | P  |
| 7439-89-6 | Iron      | 13500         |   |   | P  |
| 7439-92-1 | Lead      | 5.3           |   | * | P  |
| 7439-95-4 | Magnesium | 20000         |   | * | P  |
| 7439-96-5 | Manganese | 488           |   | N | P  |
| 7439-97-6 | Mercury   | 0.060         | U | * | AV |
| 7440-02-0 | Nickel    | 12.0          |   |   | P  |
| 7440-09-7 | Potassium | 936           | B |   | P  |
| 7782-49-2 | Selenium  | 0.97          | U |   | P  |
| 7440-22-4 | Silver    | 0.48          | U |   | P  |
| 7440-23-5 | Sodium    | 110           | B |   | P  |
| 7440-28-0 | Thallium  | 1.4           | U |   | P  |
| 7440-62-2 | Vanadium  | 10.0          | B |   | P  |
| 7440-66-6 | Zinc      | 25.7          |   | * | P  |
| 57-12-5   | Cyanide   | 1.2           |   |   | C  |

Color Before: brown

Clarity Before:

Texture: med

Color After: light yell

Clarity After: clear

Artifacts: small

Comments:

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## NYSDEC ASP

1  
INORGANIC ANALYSIS DATA SHEET

NYSDEC SAMPLE NO.

SB-6(18-20)

Lab Name: Galson Laboratories

Contract: BBL

Lab Code: 11626

Case No.:

SAS No.:

SDG No.: L62574

Matrix (soil/water): Soil

Lab Sample ID: L62574-4

Level (low/med): LOW

Date Received: 08/08/00

% Solids: 82.5

Concentration Units (ug/L or mg/kg dry weight): mg/kg

| CAS No.   | Analyte   | Concentration | C | Q | M  |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum  | 3380          |   |   | P  |
| 7440-36-0 | Antimony  | 1.2           | U | N | P  |
| 7440-38-2 | Arsenic   | 1.6           | B |   | P  |
| 7440-39-3 | Barium    | 16.3          | B |   | P  |
| 7440-41-7 | Beryllium | 0.29          | B |   | P  |
| 7440-43-9 | Cadmium   | 0.48          | U |   | P  |
| 7440-70-2 | Calcium   | 52700         |   | * | P  |
| 7440-47-3 | Chromium  | 5.2           |   |   | P  |
| 7440-48-4 | Cobalt    | 4.9           | B |   | P  |
| 7440-50-8 | Copper    | 16.5          |   | * | P  |
| 7439-89-6 | Iron      | 10000         |   |   | P  |
| 7439-92-1 | Lead      | 4.0           |   | * | P  |
| 7439-95-4 | Magnesium | 21500         |   | * | P  |
| 7439-96-5 | Manganese | 377           |   | N | P  |
| 7439-97-6 | Mercury   | 0.16          |   | * | AV |
| 7440-02-0 | Nickel    | 8.6           | B |   | P  |
| 7440-09-7 | Potassium | 709           | B |   | P  |
| 7782-49-2 | Selenium  | 0.97          | U |   | P  |
| 7440-22-4 | Silver    | 0.48          | U |   | P  |
| 7440-23-5 | Sodium    | 129           | B |   | P  |
| 7440-28-0 | Thallium  | 1.5           | U |   | P  |
| 7440-62-2 | Vanadium  | 7.8           | B |   | P  |
| 7440-66-6 | Zinc      | 19.8          |   | * | P  |
| 57-12-5   | Cyanide   | 2.7           |   |   | C  |

Color Before: brown

Clarity Before:

Texture: fine

Color After: light yell

Clarity After: clear

Artifacts: none

Comments:

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## NYSDEC ASP

1  
INORGANIC ANALYSIS DATA SHEET

NYSDEC SAMPLE NO.

SB-8(6.5-8)

Lab Name: Galson Laboratories

Contract: BBL

Lab Code: 11626

Case No.:

SAS No.:

SDG No.: L62574

Matrix (soil/water): Soil

Lab Sample ID: L62615-1

Level (low/med): LOW

Date Received: 08/10/00

% Solids: 25.0

Concentration Units (ug/L or mg/kg dry weight): mg/kg

| CAS No.   | Analyte   | Concentration | C | Q | M  |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum  | 1210          | - |   | P  |
| 7440-36-0 | Antimony  | 2.0           | U | N | P  |
| 7440-38-2 | Arsenic   | 1.6           | U |   | P  |
| 7440-39-3 | Barium    | 140           |   |   | P  |
| 7440-41-7 | Beryllium | 0.40          | U |   | P  |
| 7440-43-9 | Cadmium   | 0.80          | U |   | P  |
| 7440-70-2 | Calcium   | 21600         |   | * | P  |
| 7440-47-3 | Chromium  | 2.9           | B |   | P  |
| 7440-48-4 | Cobalt    | 0.83          | B |   | P  |
| 7440-50-8 | Copper    | 14.1          |   | * | P  |
| 7439-89-6 | Iron      | 18700         |   | * | P  |
| 7439-92-1 | Lead      | 4.5           |   | * | P  |
| 7439-95-4 | Magnesium | 917           | B | * | P  |
| 7439-96-5 | Manganese | 573           |   | N | P  |
| 7439-97-6 | Mercury   | 0.20          | B | * | AV |
| 7440-02-0 | Nickel    | 2.3           | B |   | P  |
| 7440-09-7 | Potassium | 146           | U |   | P  |
| 7782-49-2 | Selenium  | 2.6           |   |   | P  |
| 7440-22-4 | Silver    | 0.80          | U |   | P  |
| 7440-23-5 | Sodium    | 166           | U |   | P  |
| 7440-28-0 | Thallium  | 2.4           | U |   | P  |
| 7440-62-2 | Vanadium  | 2.4           | B |   | P  |
| 7440-66-6 | Zinc      | 10.6          |   | * | P  |
| 57-12-5   | Cyanide   | 1.9           | U |   | C  |

Color Before: dark brown

Clarity Before:

Texture: med

Color After: yellow

Clarity After: clear

Artifacts: rocks

Comments:

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## NYSDEC ASP

1  
INORGANIC ANALYSIS DATA SHEET

NYSDEC SAMPLE NO.

SB-10(4-7)

Lab Name: Galson Laboratories

Contract: BBL

Lab Code: 11626

Case No.:

SAS No.:

SDG No.: L62574

Matrix (soil/water): Soil

Lab Sample ID: L62615-2

Level (low/med): LOW

Date Received: 08/10/00

% Solids: 82.2

Concentration Units (ug/L or mg/kg dry weight): mg/kg

| CAS No.   | Analyte   | Concentration | C | Q | M  |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum  | 2500          |   |   | P  |
| 7440-36-0 | Antimony  | 0.61          | U | N | P  |
| 7440-38-2 | Arsenic   | 2.5           |   |   | P  |
| 7440-39-3 | Barium    | 12.1          | B |   | P  |
| 7440-41-7 | Beryllium | 0.22          | B |   | P  |
| 7440-43-9 | Cadmium   | 0.83          |   |   | P  |
| 7440-70-2 | Calcium   | 12000         |   | * | P  |
| 7440-47-3 | Chromium  | 5.1           |   |   | P  |
| 7440-48-4 | Cobalt    | 3.0           | B |   | P  |
| 7440-50-8 | Copper    | 26.2          |   | * | P  |
| 7439-89-6 | Iron      | 7990          |   |   | P  |
| 7439-92-1 | Lead      | 17.0          |   | * | P  |
| 7439-95-4 | Magnesium | 4540          |   | * | P  |
| 7439-96-5 | Manganese | 103           |   | N | P  |
| 7439-97-6 | Mercury   | 0.13          |   | * | AV |
| 7440-02-0 | Nickel    | 11.7          |   |   | P  |
| 7440-09-7 | Potassium | 487           | B |   | P  |
| 7782-49-2 | Selenium  | 0.49          | U |   | P  |
| 7440-22-4 | Silver    | 0.24          | U |   | P  |
| 7440-23-5 | Sodium    | 268           | B |   | P  |
| 7440-28-0 | Thallium  | 0.73          | U |   | P  |
| 7440-62-2 | Vanadium  | 4.8           | B |   | P  |
| 7440-66-6 | Zinc      | 2210          |   | * | P  |
| 57-12-5   | Cyanide   | 3.6           |   |   | -C |

Color Before: brown

Clarity Before:

Texture: med

Color After: yellow

Clarity After: clear

Artifacts: small

Comments:

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## NYSDEC ASP

1  
INORGANIC ANALYSIS DATA SHEET

NYSDEC SAMPLE NO.

SB-13(4-6)

Lab Name: Galson Laboratories

Contract: BBL

Lab Code: 11626

Case No.:

SAS No.:

SDG No.: L62574

Matrix (soil/water): Soil

Lab Sample ID: L62615-3

Level (low/med): LOW

Date Received: 8/10/00

% Solids: 74.5

Concentration Units (ug/L or mg/kg dry weight): mg/kg

| CAS No.   | Analyte   | Concentration        | C | Q   | M  |
|-----------|-----------|----------------------|---|-----|----|
| 7429-90-5 | Aluminum  | 3150                 |   |     | P  |
| 7440-36-0 | Antimony  | 5.5                  | B | N J | P  |
| 7440-38-2 | Arsenic   | 6.9                  |   |     | P  |
| 7440-39-3 | Barium    | 34.6                 |   |     | P  |
| 7440-41-7 | Beryllium | 0.28                 | B |     | P  |
| 7440-43-9 | Cadmium   | 8.5                  |   |     | P  |
| 7440-70-2 | Calcium   | 9440                 |   | *   | P  |
| 7440-47-3 | Chromium  | 17.2                 |   |     | P  |
| 7440-48-4 | Cobalt    | 3.7                  | B |     | P  |
| 7440-50-8 | Copper    | 74.3                 |   | *   | P  |
| 7439-89-6 | Iron      | 21000                |   |     | P  |
| 7439-92-1 | Lead      | 374                  |   | *   | P  |
| 7439-95-4 | Magnesium | 4060                 |   | *   | P  |
| 7439-96-5 | Manganese | 172                  |   | N J | P  |
| 7439-97-6 | Mercury   | 0.34                 |   | *   | AV |
| 7440-02-0 | Nickel    | 10.4                 |   |     | P  |
| 7440-09-7 | Potassium | 552                  | B |     | P  |
| 7782-49-2 | Selenium  | 1.2                  |   |     | P  |
| 7440-22-4 | Silver    | 0.27                 | U |     | P  |
| 7440-23-5 | Sodium    | 859                  |   |     | P  |
| 7440-28-0 | Thallium  | 0.81                 | U |     | P  |
| 7440-62-2 | Vanadium  | 8.2                  |   |     | P  |
| 7440-66-6 | Zinc      | 9040 <del>90.4</del> |   | * J | P  |
| 57-12-5   | Cyanide   | 0.86                 |   |     | C  |

Color Before: brown

Clarity Before:

Texture: med

Color After: yellow

Clarity After: clear

Artifacts: rocks

Comments:

FORM I - IN

10/95

## NYSDEC ASP

1  
INORGANIC ANALYSIS DATA SHEET

NYSDEC SAMPLE NO.

SB-14(5-7)

Lab Name: Galson Laboratories

Contract: BBL

Lab Code: 11626

Case No.:

SAS No.:

SDG No.: L62574

Matrix (soil/water): Soil

Lab Sample ID: L62615-5

Level (low/med): LOW

Date Received: 08/10/00

% Solids: 72.0

Concentration Units (ug/L or mg/kg dry weight): mg/kg

| CAS No.   | Analyte   | Concentration | C | Q    | M  |
|-----------|-----------|---------------|---|------|----|
| 7429-90-5 | Aluminum  | 3580          |   |      | P  |
| 7440-36-0 | Antimony  | 0.69          | U | N WJ | P  |
| 7440-38-2 | Arsenic   | 1.5           |   |      | P  |
| 7440-39-3 | Barium    | 38.8          |   |      | P  |
| 7440-41-7 | Beryllium | 0.24          | B |      | P  |
| 7440-43-9 | Cadmium   | 1.3           |   |      | P  |
| 7440-70-2 | Calcium   | 5760          |   | *    | P  |
| 7440-47-3 | Chromium  | 4.5           |   |      | P  |
| 7440-48-4 | Cobalt    | 3.3           | B |      | P  |
| 7440-50-8 | Copper    | 15.3          |   | *    | P  |
| 7439-89-6 | Iron      | 7550          |   | *    | P  |
| 7439-92-1 | Lead      | 168           |   | *    | P  |
| 7439-95-4 | Magnesium | 2710          |   | *    | P  |
| 7439-96-5 | Manganese | 106           |   | N J  | P  |
| 7439-97-6 | Mercury   | 1.5           |   | *    | AV |
| 7440-02-0 | Nickel    | 9.0           |   |      | P  |
| 7440-09-7 | Potassium | 522           | B |      | P  |
| 7782-49-2 | Selenium  | 0.56          | U |      | P  |
| 7440-22-4 | Silver    | 0.28          | U |      | P  |
| 7440-23-5 | Sodium    | 193           | B |      | P  |
| 7440-28-0 | Thallium  | 0.83          | U |      | P  |
| 7440-62-2 | Vanadium  | 8.5           |   |      | P  |
| 7440-66-6 | Zinc      | 1360          |   | * J  | P  |
| 57-12-5   | Cyanide   | 0.68          | U |      | -C |

Color Before: brown

Clarity Before:

Texture: med

Color After: yellow

Clarity After: clear

Artifacts: rocks

Comments:

FORM I - IN

10/95

## NYSDEC ASP

1  
INORGANIC ANALYSIS DATA SHEET

NYSDEC SAMPLE NO.

SB-15(4-6)

Lab Name: Galson Laboratories

Contract: BBL

Lab Code: 11626

Case No.:

SAS No.:

SDG No.: L62574

Matrix (soil/water): Soil

Lab Sample ID: L62615-4

Level (low/med): LOW

Date Received: 08/10/00

% Solids: 70.9

Concentration Units (ug/L or mg/kg dry weight): mg/kg

| CAS No.   | Analyte   | Concentration | C | Q | M  |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum  | 3960          | — |   | P  |
| 7440-36-0 | Antimony  | 0.71          | U | N | P  |
| 7440-38-2 | Arsenic   | 2.2           |   |   | P  |
| 7440-39-3 | Barium    | 22.6          | B |   | P  |
| 7440-41-7 | Beryllium | 0.31          | B |   | P  |
| 7440-43-9 | Cadmium   | 0.28          | U |   | P  |
| 7440-70-2 | Calcium   | 5060          |   | * | P  |
| 7440-47-3 | Chromium  | 5.1           |   |   | P  |
| 7440-48-4 | Cobalt    | 3.2           | B |   | P  |
| 7440-50-8 | Copper    | 21.4          |   | * | P  |
| 7439-89-6 | Iron      | 8410          |   |   | P  |
| 7439-92-1 | Lead      | 28.8          |   | * | P  |
| 7439-95-4 | Magnesium | 3090          |   | * | P  |
| 7439-96-5 | Manganese | 122           |   | N | P  |
| 7439-97-6 | Mercury   | 0.41          |   | * | AV |
| 7440-02-0 | Nickel    | 8.6           |   |   | P  |
| 7440-09-7 | Potassium | 542           | B |   | P  |
| 7782-49-2 | Selenium  | 0.56          | U |   | P  |
| 7440-22-4 | Silver    | 0.28          | U |   | P  |
| 7440-23-5 | Sodium    | 159           | B |   | P  |
| 7440-28-0 | Thallium  | 0.85          | U |   | P  |
| 7440-62-2 | Vanadium  | 7.0           | B |   | P  |
| 7440-66-6 | Zinc      | 1900          |   | * | P  |
| 57-12-5   | Cyanide   | 1.1           |   |   | -C |

Color Before: brown

Clarity Before:

Texture: med

Color After: yellow

Clarity After: clear

Artifacts: rocks

Comments:

FORM I - IN

10/95

## NYSDEC ASP

1  
INORGANIC ANALYSIS DATA SHEET

NYSDEC SAMPLE NO.

DUP-2

Lab Name: Galson Laboratories

Contract: BBL

Lab Code: 11626

Case No.:

SAS No.:

SDG No.: L62574

Matrix (soil/water): Soil

Lab Sample ID: L62615-6

Level (low/med): LOW

Date Received: 8/10/00

% 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  | 3400          |   |     | P  |
| 7440-36-0 | Antimony  | 2.9           | B | N J | P  |
| 7440-38-2 | Arsenic   | 4.0           |   |     | P  |
| 7440-39-3 | Barium    | 26.9          |   |     | P  |
| 7440-41-7 | Beryllium | 0.29          | B |     | P  |
| 7440-43-9 | Cadmium   | 10.6          |   |     | P  |
| 7440-70-2 | Calcium   | 11900         |   | *   | P  |
| 7440-47-3 | Chromium  | 14.3          |   |     | P  |
| 7440-48-4 | Cobalt    | 3.1           | B |     | P  |
| 7440-50-8 | Copper    | 77.9          |   | *   | P  |
| 7439-89-6 | Iron      | 20200         |   | *   | P  |
| 7439-92-1 | Lead      | 219           |   | *   | P  |
| 7439-95-4 | Magnesium | 6280          |   | *   | P  |
| 7439-96-5 | Manganese | 159           |   | N J | P  |
| 7439-97-6 | Mercury   | 0.23          |   | *   | AV |
| 7440-02-0 | Nickel    | 8.7           |   |     | P  |
| 7440-09-7 | Potassium | 564           | B |     | P  |
| 7782-49-2 | Selenium  | 1.4           |   |     | P  |
| 7440-22-4 | Silver    | 0.25          | U |     | P  |
| 7440-23-5 | Sodium    | 793           |   |     | P  |
| 7440-28-0 | Thallium  | 0.76          | U |     | P  |
| 7440-62-2 | Vanadium  | 8.0           |   |     | P  |
| 7440-66-6 | Zinc      | 8360          |   | * J | P  |
| 57-12-5   | Cyanide   | 0.62          | U |     | C  |

Color Before: brown

Clarity Before:

Texture: med

Color After: yellow

Clarity After: clear

Artifacts: rocks

Comments:

FORM I - IN

10/95

BLASLAND, BOUCK & LEE, INC.

LABORATORY DATA REVIEW REPORT

Project: NMPC - Cedar Street

Analytical Laboratory: Galson Laboratories

Laboratory Report Identification Number: L66535

Date of Laboratory Report: January 8, 2001

Date of Review: February 17, 2001

Reviewer: Laurie Indick

Number of Samples: 6

Sample Matrix: soil

Date of Collection: 12/07/00

Sample Analysis: PAH

Quality Control Checks

|  |            |           |                       |
|--|------------|-----------|-----------------------|
| 1. Field Chain-of-Custody complete   | <u>yes</u> | no        | not applicable        |
| 2. Proper methods for analysis used  | <u>yes</u> | no        | not applicable        |
| 3. All documentation supplied  | <u>yes</u> | no        | not applicable        |
| 4. Samples analyzed within specified holding times                             | <u>yes</u> | no        | not applicable        |
| 5. The minimum number of field and laboratory QC samples analyzed              | <u>yes</u> | no        | not applicable        |
| 6. Laboratory accuracy maintained within established ranges for the following: |            |           |                       |
| - %RSD, initial calibration  | <u>yes</u> | no        | not applicable        |
| - %D, continuing calibration   | <u>yes</u> | no        | not applicable        |
| - %Recovery, matrix spike  | <u>yes</u> | no        | not applicable        |
| - %Recovery, blank spike   | yes        | <u>no</u> | not applicable        |
| - %Recovery, surrogate   | yes        | <u>no</u> | not applicable        |
| - %Recovery, control sample  | yes        | no        | <u>not applicable</u> |

7. Laboratory precision maintained within established ranges for the following:
- |                     |            |    |                |
|---------------------|------------|----|----------------|
| - RPD, matrix spike | <u>yes</u> | no | not applicable |
| - RPD, duplicates   | <u>yes</u> | no | not applicable |
8. Target analyte concentrations below detection limit in all blank samples
- |  |            |    |                |
|--|------------|----|----------------|
|  | <u>yes</u> | no | not applicable |
|--|------------|----|----------------|

Notes: \_\_\_\_\_

Recovery for one surrogate was outside control limits in samples TP-3-8-6 and TP-4-10-6. Since recoveries for the remaining surrogates were within control limits, no data have been qualified based on the deviations.

The response for one or more internal standards were below established limits in all samples. Data for all compounds quantitated under the noncompliant standards have either been qualified as estimated or rejected, depending on the severity of the deviation.

The recovery for acenaphthylene was above control limits in the matrix spike blank. No action has been taken as a result of this deficiency since acceptable recoveries were demonstrated in the matrix spike/matrix spike duplicate.

Other than for the deviations noted in this review, all data quality parameters were within method-specified limits and the data is acceptable for use as reported by the laboratory.



Reviewed and Approved:

  
\_\_\_\_\_  
Quality Assurance Manager

\_\_\_\_\_  
Project Manager

## SAMPLE COMPLIANCE REPORT

NMPC - Cedar Street

| Sample Delivery Group | Sampling Date | ASP Protocol | Sample ID | Matrix | Compliance <sup>1</sup> |     |     |    | Noncompliance                    |
|-----------------------|---------------|--------------|-----------|--------|-------------------------|-----|-----|----|----------------------------------|
|                       |               |              |           |        | VOA                     | BNA | PCB | Cn |                                  |
| L62197                | 12/07/00      | 1995         | TP-2-11-3 | soil   | --                      | no  | --  | -- | BN - int std.                    |
| L62197                | 12/07/00      | 1995         | TP-3-8-6  | soil   | --                      | no  | --  | -- | BN - surr <sup>2</sup> , int std |
| L62197                | 12/07/00      | 1995         | TP-4-3-3  | soil   | --                      | no  | --  | -- | BN - int std.                    |
| L62197                | 12/07/00      | 1995         | TP-4-10-6 | soil   | --                      | no  | --  | -- | BN - surr <sup>2</sup> , int std |
| L62197                | 12/07/00      | 1995         | TP-5-15-5 | soil   | --                      | no  | --  | -- | BN - int std.                    |
| L62197                | 12/07/00      | 1995         | DUP-1     | soil   | --                      | no  | --  | -- | BN - int std.                    |
|                       |               |              |           |        |                         |     |     |    |                                  |
|                       |               |              |           |        |                         |     |     |    |                                  |
|                       |               |              |           |        |                         |     |     |    |                                  |
|                       |               |              |           |        |                         |     |     |    |                                  |
|                       |               |              |           |        |                         |     |     |    |                                  |
|                       |               |              |           |        |                         |     |     |    |                                  |
|                       |               |              |           |        |                         |     |     |    |                                  |
|                       |               |              |           |        |                         |     |     |    |                                  |
|                       |               |              |           |        |                         |     |     |    |                                  |
|                       |               |              |           |        |                         |     |     |    |                                  |

- 1 Samples which are compliant with no added validation qualifiers are listed as "yes". Samples which are non-compliant or which have added qualifiers are listed as "no". A "no" designation does not necessarily indicate that the data have been rejected or are otherwise unusable.
- 2 The noncompliance resulted in no qualification of data.
- 3 Although the deviation resulted in the qualification of data, the laboratory was method compliant.

SAMPLE NO.

DUP-1

Contract: Blasland, B

SDG No.: L66535

Lab Sample ID: L66535-8

Lab File ID: ED121315

Date Received: 12/07/00

Date Extracted:12/12/00

Date Analyzed: 12/13/00

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: 7.0

(ug/L or ug/Kg) ug/kg

Q

2424

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

DUP-1 DL

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L66535

Matrix: (soil/water) SOIL

Lab Sample ID: L66535-8DL

Sample wt/vol: 31.25 (g/mL) g

Lab File ID: ED121311

Level: (low/med) LOW

Date Received: 12/07/00

% Moisture: 27 decanted: (Y/N) N

Date Extracted: 12/12/00

Concentrated Extract Volume: 1000 (uL) Date Analyzed: 12/13/00

Injection Volume: 2.0 (uL)

Dilution Factor: 10.0

GPC Cleanup: (Y/N) N pH: 7.0

CONCENTRATION UNITS:

(ug/L or ug/Kg) ug/kg

Q

| CAS NO.  | COMPOUND               | CONCENTRATION UNITS: | Q   |
|----------|------------------------|----------------------|-----|
| 91-20-3  | Naphthalene            | 4400                 | U D |
| 91-57-6  | 2-Methylnaphthalene    | 4400                 | U D |
| 208-96-8 | Acenaphthylene         | 4400                 | U D |
| 83-32-9  | Acenaphthene           | 4400                 | U D |
| 86-73-7  | Fluorene               | 4400                 | U D |
| 85-01-8  | Phenanthrene           | 1600                 | JD  |
| 120-12-7 | Anthracene             | 440                  | JD  |
| 206-44-0 | Fluoranthene           | 4000                 | JD  |
| 129-00-0 | Pyrene                 | 9700                 | D   |
| 56-55-3  | Benzo(a)anthracene     | 5800                 | DJ  |
| 218-01-9 | Chrysene               | 5300                 | DJ  |
| 205-99-2 | Benzo(b)fluoranthene   | 7800                 | DJ  |
| 207-08-9 | Benzo(k)fluoranthene   | 3100                 | JD  |
| 50-32-8  | Benzo(a)pyrene         | 5900                 | DJ  |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 4700                 | DJ  |
| 53-70-3  | Dibenzo(a,h)anthracene | 930                  | JD  |
| 191-24-2 | Benzo(g,h,i)perylene   | 6500                 | DJ  |

SAMPLE NO.

Contract: Blasland, B

TP-2-11-3

SDG No. : L66535

Lab Sample ID: L66535-1

Lab File ID: ED121317

Date Received: 12/07/00

Date Extracted:12/12/00

Date Analyzed: 12/13/00

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: 7.0

CONCENTRATION UNITS:

COMPOUND

(ug/L or ug/Kg) ug/kg Q

|                                       |           |    |
|---------------------------------------|-----------|----|
| 91-20-3-----Naphthalene               | 110       | J  |
| 91-57-6-----2-Methylnaphthalene       | 430       | U  |
| 208-96-8-----Acenaphthylene           | 320       | J  |
| 83-32-9-----Acenaphthene              | 430       | U  |
| 86-73-7-----Fluorene                  | 89        | J  |
| 85-01-8-----Phenanthrene              | 3000      |    |
| 120-12-7-----Anthracene               | 580       |    |
| 206-44-0-----Fluoranthene             | 8100      | EJ |
| 129-00-0-----Pyrene                   | 9900 7500 | E  |
| 56-55-3-----Benzo (a) anthracene      | 4200 5000 | EJ |
| 218-01-9-----Chrysene                 | 3600 4200 | EJ |
| 205-99-2-----Benzo (b) fluoranthene   | 5200 7000 | EJ |
| 207-08-9-----Benzo (k) fluoranthene   | 1400      | J  |
| 50-32-8-----Benzo (a) pyrene          | 5800 5100 | EJ |
| 193-39-5-----Indeno (1,2,3-cd) pyrene | 2700      | J  |
| 53-70-3-----Dibenzo (a,h) anthracene  | 380       | J  |
| 191-24-2-----Benzo (g,h,i) perylene   | 3000      | J  |

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

TP-2-11-3 DL

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L66535

Matrix: (soil/water) SOIL

Lab Sample ID: L66535-1DL

Sample wt/vol: 31.3 (g/mL) g

Lab File ID: ED121313

Level: (low/med) LOW

Date Received: 12/07/00

% Moisture: 26 decanted: (Y/N) N

Date Extracted: 12/12/00

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 12/13/00

Injection Volume: 2.0 (uL)

Dilution Factor: 10.0

GPC Cleanup: (Y/N) N pH: 7.0

CONCENTRATION UNITS:

(ug/L or ug/Kg) ug/kg Q

| CAS NO.  | COMPOUND               | CONCENTRATION UNITS:<br>(ug/L or ug/Kg) ug/kg | Q   |
|----------|------------------------|---|-----|
| 91-20-3  | Naphthalene            | 4300  | U D |
| 91-57-6  | 2-Methylnaphthalene    | 4300  | U D |
| 208-96-8 | Acenaphthylene         | 4300  | U D |
| 83-32-9  | Acenaphthene           | 4300  | U D |
| 86-73-7  | Fluorene               | 4300  | U D |
| 85-01-8  | Phenanthrene           | 1600  | JD  |
| 120-12-7 | Anthracene             | 450   | JD  |
| 206-44-0 | Fluoranthene           | 3900  | JD  |
| 129-00-0 | Pyrene                 | 7900  | D   |
| 56-55-3  | Benzo(a)anthracene     | 5000  | DJ  |
| 218-01-9 | Chrysene               | 4200  | JD  |
| 205-99-2 | Benzo(b)fluoranthene   | 7000  | DJ  |
| 207-08-9 | Benzo(k)fluoranthene   | 4200  | JD  |
| 50-32-8  | Benzo(a)pyrene         | 5100  | DJ  |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 3600  | JD  |
| 53-70-3  | Dibenzo(a,h)anthracene | 760   | JD  |
| 191-24-2 | Benzo(g,h,i)perylene   | 4700  | DJ  |

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

TP-3-8-6

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L66535

Matrix: (soil/water) SOIL

Lab Sample ID: L66535-2

Sample wt/vol: 31.3 (g/mL) g

Lab File ID: ED121314

Level: (low/med) LOW

Date Received: 12/07/00

% Moisture: 38 decanted: (Y/N) N

Date Extracted: 12/12/00

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 12/13/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: 7.0

CONCENTRATION UNITS:

(ug/L or ug/Kg) ug/kg Q

| CAS NO.       | COMPOUND                 |      |   |
|---------------|--------------------------|------|---|
| 91-20-3-----  | Naphthalene              | 520  | U |
| 91-57-6-----  | 2-Methylnaphthalene      | 520  | U |
| 208-96-8----- | Acenaphthylene           | 66   | J |
| 83-32-9-----  | Acenaphthene             | 190  | J |
| 86-73-7-----  | Fluorene                 | 220  | J |
| 85-01-8-----  | Phenanthrene             | 380  | J |
| 120-12-7----- | Anthracene               | 260  | J |
| 206-44-0----- | Fluoranthene             | 1000 |   |
| 129-00-0----- | Pyrene                   | 1400 |   |
| 56-55-3-----  | Benzo (a) anthracene     | 740  | J |
| 218-01-9----- | Chrysene                 | 660  | J |
| 205-99-2----- | Benzo (b) fluoranthene   | 860  | J |
| 207-08-9----- | Benzo (k) fluoranthene   | 310  | J |
| 50-32-8-----  | Benzo (a) pyrene         | 830  | J |
| 193-39-5----- | Indeno (1,2,3-cd) pyrene | 500  | J |
| 53-70-3-----  | Dibenzo (a,h) anthracene | 100  | J |
| 191-24-2----- | Benzo (g,h,i) perylene   | 400  | J |

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

TP-3-8-6 R

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L66535

Matrix: (soil/water) SOIL

Lab Sample ID: L66535-2R

Sample wt/vol: 31.3 (g/mL) g

Lab File ID: ED121318

Level: (low/med) LOW

Date Received: 12/07/00

% Moisture: 38 decanted: (Y/N) N

Date Extracted: 12/12/00

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 12/13/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: 7.0

| CAS NO.  | COMPOUND                 | CONCENTRATION UNITS: |         |
|----------|--------------------------|----------------------|---------|
|          |                          | (ug/L or ug/Kg)      | ug/kg Q |
| 91-20-3  | Naphthalene              | 520                  | U       |
| 91-57-6  | 2-Methylnaphthalene      | 520                  | U       |
| 208-96-8 | Acenaphthylene           | 79                   | J       |
| 83-32-9  | Acenaphthene             | 180                  | J       |
| 86-73-7  | Fluorene                 | 230                  | J       |
| 85-01-8  | Phenanthrene             | 390                  | J       |
| 120-12-7 | Anthracene               | 280                  | J       |
| 206-44-0 | Fluoranthene             | 1400                 |         |
| 129-00-0 | Pyrene                   | 1500                 |         |
| 56-55-3  | Benzo (a) anthracene     | 720                  |         |
| 218-01-9 | Chrysene                 | 510                  | J       |
| 205-99-2 | Benzo (b) fluoranthene   | 760                  |         |
| 207-08-9 | Benzo (k) fluoranthene   | 230                  | J       |
| 50-32-8  | Benzo (a) pyrene         | 600                  |         |
| 193-39-5 | Indeno (1,2,3-cd) pyrene | 300                  | J       |
| 53-70-3  | Dibenzo (a,h) anthracene | 100                  | J       |
| 191-24-2 | Benzo (g,h,i) perylene   | 400                  | J       |



1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

TP-4-10-6

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L66535

Matrix: (soil/water) SOIL

Lab Sample ID: L66535-4

Sample wt/vol: 31.42 (g/mL) g

Lab File ID: ED121306

Level: (low/med) LOW

Date Received: 12/07/00

% Moisture: 73 decanted: (Y/N) N

Date Extracted: 12/12/00

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 12/13/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: 7.0

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) ug/kg

Q

|               |                        |      |    |
|---------------|------------------------|------|----|
| 91-20-3-----  | Naphthalene            | 1200 | U  |
| 91-57-6-----  | 2-Methylnaphthalene    | 1200 | U  |
| 208-96-8----- | Acenaphthylene         | 1200 | U  |
| 83-32-9-----  | Acenaphthene           | 150  | J  |
| 86-73-7-----  | Fluorene               | 1200 | U  |
| 85-01-8-----  | Phenanthrene           | 1200 | U  |
| 120-12-7----- | Anthracene             | 1200 | U  |
| 206-44-0----- | Fluoranthene           | 1200 | U  |
| 129-00-0----- | Pyrene                 | 1200 | U  |
| 56-55-3-----  | Benzo(a)anthracene     | 1200 | UJ |
| 218-01-9----- | Chrysene               | 1200 | UJ |
| 205-99-2----- | Benzo(b)fluoranthene   | 1200 | UJ |
| 207-08-9----- | Benzo(k)fluoranthene   | 1200 | UJ |
| 50-32-8-----  | Benzo(a)pyrene         | 1200 | UJ |
| 193-39-5----- | Indeno(1,2,3-cd)pyrene | 1200 | UJ |
| 53-70-3-----  | Dibenzo(a,h)anthracene | 1200 | UJ |
| 191-24-2----- | Benzo(g,h,i)perylene   | 1200 | UJ |

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

TP-4-3-3

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L66535

Matrix: (soil/water) SOIL

Lab Sample ID: L66535-3

Sample wt/vol: 30.92 (g/mL) g

Lab File ID: ED121316

Level: (low/med) LOW

Date Received: 12/07/00

% Moisture: 29 decanted: (Y/N) N

Date Extracted: 12/12/00

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 12/13/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: 7.0

CAS NO.

COMPOUND

CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg

Q

|               |                          |      |   |
|---------------|--------------------------|------|---|
| 91-20-3-----  | Naphthalene              | 46   | J |
| 91-57-6-----  | 2-Methylnaphthalene      | 460  | U |
| 208-96-8----- | Acenaphthylene           | 130  | J |
| 83-32-9-----  | Acenaphthene             | 52   | J |
| 86-73-7-----  | Fluorene                 | 71   | J |
| 85-01-8-----  | Phenanthrene             | 1200 |   |
| 120-12-7----- | Anthracene               | 260  | J |
| 206-44-0----- | Fluoranthene             | 1900 |   |
| 129-00-0----- | Pyrene                   | 2500 |   |
| 56-55-3-----  | Benzo (a) anthracene     | 1200 | J |
| 218-01-9----- | Chrysene                 | 1200 | J |
| 205-99-2----- | Benzo (b) fluoranthene   | 1500 | J |
| 207-08-9----- | Benzo (k) fluoranthene   | 460  | J |
| 50-32-8-----  | Benzo (a) pyrene         | 1200 | J |
| 193-39-5----- | Indeno (1,2,3-cd) pyrene | 710  | J |
| 53-70-3-----  | Dibenzo (a,h) anthracene | 170  | J |
| 191-24-2----- | Benzo (g,h,i) perylene   | 850  | J |

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

TP-4-3-3 R

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L66535

Matrix: (soil/water) SOIL

Lab Sample ID: L66535-3R

Sample wt/vol: 30.92 (g/mL) g

Lab File ID: ED121320

Level: (low/med) LOW

Date Received: 12/07/00

% Moisture: 29 decanted: (Y/N) N

Date Extracted: 12/12/00

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 12/13/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: 7.0

CONCENTRATION UNITS:

(ug/L or ug/Kg) ug/kg

Q

| CAS NO.  | COMPOUND               | CONCENTRATION UNITS: | Q |
|----------|------------------------|----------------------|---|
| 91-20-3  | Naphthalene            | 460                  | U |
| 91-57-6  | 2-Methylnaphthalene    | 460                  | U |
| 208-96-8 | Acenaphthylene         | 130                  | J |
| 83-32-9  | Acenaphthene           | 47                   | J |
| 86-73-7  | Fluorene               | 68                   | J |
| 85-01-8  | Phenanthrene           | 1200                 |   |
| 120-12-7 | Anthracene             | 270                  | J |
| 206-44-0 | Fluoranthene           | 1700                 |   |
| 129-00-0 | Pyrene                 | 3500                 |   |
| 56-55-3  | Benzo(a)anthracene     | 1300                 |   |
| 218-01-9 | Chrysene               | 1300                 |   |
| 205-99-2 | Benzo(b)fluoranthene   | 1400                 |   |
| 207-08-9 | Benzo(k)fluoranthene   | 400                  | J |
| 50-32-8  | Benzo(a)pyrene         | 1100                 |   |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 820                  |   |
| 53-70-3  | Dibenzo(a,h)anthracene | 160                  | J |
| 191-24-2 | Benzo(g,h,i)perylene   | 1200                 |   |

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

TP-5-15-5

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L66535

Matrix: (soil/water) SOIL

Lab Sample ID: L66535-7

Sample wt/vol: 31.27 (g/mL) g

Lab File ID: ED121305

Level: (low/med) LOW

Date Received: 12/07/00

% Moisture: 32 decanted: (Y/N) N

Date Extracted: 12/12/00

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 12/13/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: 7.0

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) ug/kg Q

|               |                          |     |    |
|---------------|--------------------------|-----|----|
| 91-20-3-----  | Naphthalene              | 470 | U  |
| 91-57-6-----  | 2-Methylnaphthalene      | 470 | U  |
| 208-96-8----- | Acenaphthylene           | 470 | U  |
| 83-32-9-----  | Acenaphthene             | 470 | U  |
| 86-73-7-----  | Fluorene                 | 470 | U  |
| 85-01-8-----  | Phenanthrene             | 470 | U  |
| 120-12-7----- | Anthracene               | 470 | U  |
| 206-44-0----- | Fluoranthene             | 470 | U  |
| 129-00-0----- | Pyrene                   | 470 | U  |
| 56-55-3-----  | Benzo (a) anthracene     | 470 | U  |
| 218-01-9----- | Chrysene                 | 470 | U  |
| 205-99-2----- | Benzo (b) fluoranthene   | 470 | UJ |
| 207-08-9----- | Benzo (k) fluoranthene   | 470 | UJ |
| 50-32-8-----  | Benzo (a) pyrene         | 470 | UJ |
| 193-39-5----- | Indeno (1,2,3-cd) pyrene | 470 | UJ |
| 53-70-3-----  | Dibenzo (a,h) anthracene | 470 | UJ |
| 191-24-2----- | Benzo (g,h,i) perylene   | 470 | UJ |

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

TP-5-15-5 R

Lab Name: GALSON LABORATORIES

Contract: Blasland, B

Lab Code:

Case No.: 1

SAS No.:

SDG No.: L66535

Matrix: (soil/water) SOIL

Lab Sample ID: L66535-7R

Sample wt/vol: 31.27 (g/mL) g

Lab File ID: ED121309

Level: (low/med) LOW

Date Received: 12/07/00

% Moisture: 32 decanted: (Y/N) N

Date Extracted: 12/12/00

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 12/13/00

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: 7.0

CONCENTRATION UNITS:

(ug/L or ug/Kg) ug/kg

Q

| CAS NO.  | COMPOUND               | CONCENTRATION UNITS: | Q |
|----------|------------------------|----------------------|---|
| 91-20-3  | Naphthalene            | 470                  | U |
| 91-57-6  | 2-Methylnaphthalene    | 470                  | U |
| 208-96-8 | Acenaphthylene         | 470                  | U |
| 83-32-9  | Acenaphthene           | 470                  | U |
| 86-73-7  | Fluorene               | 470                  | U |
| 85-01-8  | Phenanthrene           | 470                  | U |
| 120-12-7 | Anthracene             | 470                  | U |
| 206-44-0 | Fluoranthene           | 470                  | U |
| 129-00-0 | Pyrene                 | 470                  | U |
| 56-55-3  | Benzo(a)anthracene     | 470                  | U |
| 218-01-9 | Chrysene               | 470                  | U |
| 205-99-2 | Benzo(b)fluoranthene   | 470                  | U |
| 207-08-9 | Benzo(k)fluoranthene   | 470                  | U |
| 50-32-8  | Benzo(a)pyrene         | 470                  | U |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 470                  | U |
| 53-70-3  | Dibenzo(a,h)anthracene | 470                  | U |
| 191-24-2 | Benzo(g,h,i)perylene   | 470                  | U |

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