



O'BRIEN & GERE
ENGINEERS, INC.

February 19, 2002

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NYSDEC
Bureau of Radiation & Hazardous Site Management
Division of Solid & Hazardous Materials
625 Broadway
Albany, New York 12233-7252

RECEIVED
MAY 6 2002
NYSDEC
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Re: Additional Investigation/IRM
Activities – Former Allied Specialty
Chemical Site, Tonawanda, New York

File: 1163/26080 #2

Dear Mr. Terbush:

This letter serves to transmit and summarize data generated pursuant to the implementation of the Additional Investigation and Interim Remedial Measures Work Plan (Work Plan) dated May 2000 for the former Allied Specialty Chemical site (the Site) located in Tonawanda, New York. Work tasks completed under the Work Plan included the following: storm sewer evaluation, ground water monitoring well installations, and ground water, storm water, and sediment sampling. The various work tasks presented in the Work Plan were completed between November 2000 and December 2001.

SECTION 1 - INTRODUCTION

Site Setting

The Site is located within the western section of the Town of Tonawanda, Erie County, New York (Figure 1). The Site is located on the southwest corner of an industrial property (presently the Tonawanda Coke Corporation) in an extensively industrialized area on the east side of River Road, approximately 1.5 miles south of the South Grand Island Bridge. The Site is bordered to the south by a petroleum terminal, and to the west by River Road. The north and east boundaries abut the active Tonawanda Coke facility.

Local Hydrogeology

The Site is located within the Erie-Ontario lowlands physiographic province. The province was formerly a lake bottom during Late Wisconsin deglaciation, and is characterized by generally flat topography. The land elevation ranges between 590 and 600 feet above sea level, and slopes to the west towards River Road.

The surficial on-Site material is composed of alluvial/lacustrine sediments that are overlain by a thin veneer of fill material. This fill material, which consists of slag, tile, brick, sand and silt, averages approximately eight feet in thickness, except in the location of the former disposal pit (Figure 2) where there is approximately 10 feet of fill material. The water-bearing fill unit is not used as a



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source of potable water, and is not locally in hydraulic connection with the underlying bedrock aquifer.

The underlying bedrock aquifer has been shown to regionally flow towards the Niagara River (ERM 1992). The underlying glaciolacustrine clay has a low hydraulic conductivity, averaging 1×10^{-7} cm/sec. The clay deposit is underlain by the Camillus Shale formation. This formation consists of brownish gray non-fossiliferous mudstone with numerous gypsum and anhydrite inclusions and seams (ERM, 1992).

Based on the June 2001 ground water elevations, the local direction of ground water flow in the fill unit is predominantly to the west. This westerly flow direction is consistent with that reported during the May 1999 investigation (O'Brien & Gere Engineers, Inc., November 1999), and during the post-remediation quarterly ground water sampling conducted between September 1991 and June 1992 (ERM-Northeast, August 1992). A 48-inch storm sewer crosses the central portion of the Site and is oriented northwest to southeast. In the western portion of the Site a 36-inch storm sewer is generally oriented north-south perpendicular to ground water flow. The ground water elevations relative to the invert elevation of the 36-inch storm sewer were evaluated. Review of these relative elevations indicates that the invert elevation of the 36-inch storm sewer is below the water table. This suggests that ground water has the potential to discharge to the storm sewer.

Summary of Environmental Investigations

The primary focus of previous environmental investigations completed at the Site was a former disposal pit utilized for the disposal of spent catalysts generated during Research and Development (R & D) studies at the facility. Discharge into the pit occurred between 1958 and 1962, and R & D activities were discontinued in 1965. The pit area, located on the western side of the property, was estimated, by former plant personnel and aerial photographs, to be approximately 30 feet in diameter and 6 feet in depth. The wastes disposed of in this pit allegedly included off-specification batches of magnesium-chromate catalyst.

A series of investigations (under an Administrative Consent Order) were completed around the disposal pit to evaluate the lateral extent of the materials and to evaluate potential remedial options. During May 1991, the soils within the former disposal pit and the sediments within the adjacent catch basin were removed in accordance with an approved Cleanup Plan. The catch basin was cleaned with a high-pressure potable-water wash, and sediment was disposed of along with the excavated disposal pit soil. The resulting excavations were backfilled with clean soil and revegetated. A Cleanup Plan Implementation Report dated July 1991 was submitted to the New York State Department of Environmental Conservation (NYSDEC) documenting remedial activities, and was subsequently approved by the NYSDEC on August 5, 1991 (ERM, 1992).

In May 1999, samples were collected from eight ground water monitoring wells, four storm water locations, four sediment locations and five potential source areas in accordance with a Work Plan approved by the NYSDEC on April 15, 1999. The samples were analyzed for volatile organic compounds (VOCs), semivolatile organic compounds (SVOCs), and metals in accordance with the NYSDEC 1995 Analytical Services Protocol. The investigation results were presented to Honeywell and NYSDEC in a letter report (O'Brien & Gere Engineers, Inc., November 1999). Potential source area water samples were collected from five floor drains and one catch basin. The data indicated that low concentrations of VOCs and SVOCs were detected in floor drains. In addition, low

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concentrations of VOCs were detected in Catch Basin CB-1. The concentrations detected do not indicate a source, but rather residual contamination from incidental use over time. The May 1999 sediment, ground water, and storm water data are presented in the sections below for comparison to the more recently collected data.

To address NYSDEC comments regarding the November 1999 letter report, an Additional Investigation and Interim Remedial Measures Work Plan (O'Brien & Gere Engineers, Inc., May 2000) was prepared on behalf of Honeywell. Field efforts were implemented between November 2000 and December 2001. The following provides a discussion of the Work Plan implementation and associated analytical results.

SECTION 2 – FIELD INVESTIGATIONS AND ANALYTICAL RESULTS

Storm Sewer Evaluation

Task 1 of the Work Plan described the cleaning of the 48-inch and 36-inch storm sewers using hydraulic flushing and inspection of the sewers using video camera. National Vacuum was subcontracted to provide the hydraulic flushing and video inspection services. Pursuant to a request by NYSDEC, removal of sediment in the areas of Inlets A and B and Outfalls #1 and #2 was conducted prior to the initiation of the hydraulic flushing. During the sediment removal in the Outfall area, silt fencing was constructed to minimize sediment transport off-site. Excavated sediments were staged on plastic sheeting. Subsequent to sediment removal, confirmation sampling of the remaining sediment was conducted.

After the initial sediment removal and sampling, the 36-inch and 48-inch sewer were hydraulically flushed to remove accumulated sediments. Following the flushing, the sewers were televised to evaluate potential sources of inflow to the sewers.

Approximately 200 ft. of the 36-inch storm sewer, south from Manhole E toward the inlet, was cleaned. This is about 40% of the length of the pipe. Upon initiation of the cleaning, extensive quantities of sediment and debris were encountered in the sewer, which made further hydraulic flushing impossible. It is estimated that approximately 15 cubic yards of sediment had been removed. Approximately 100 ft. (20%) of the 36-inch pipe was televised, beginning at the sewer inlet. A large amount of debris was observable beyond the portion of the pipe that was hydraulically flushed. The debris precluded television inspection of the remaining length of pipe. There were no laterals or other connections observed, and groundwater infiltration was not noted in the televised section of pipe.

The 48-inch sewer could not be hydraulically flushed. The outfall of the 48-inch sewer appears to be more than half full of debris and sediments. Approximately 400 ft. (50%) of the 48-inch storm sewer was televised, starting at the inlet. Three manholes and several lateral connections were seen in the video. As with the 36-inch sewer, groundwater infiltration was not noted in the televised section of this pipe.

The hydraulic flushing and television inspection of the storm sewers has been completed to the extent possible. The presence of large amounts of debris in the sewers made further investigatory work impractical, without substantial rehabilitation of the pipelines. Based on the work completed, no significant sources of groundwater infiltration into the sewers have been observed. The portions of the two sewers that were televised, 20% of the 36-inch pipe and 50% of the 48-inch pipe, are reasonably

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representative, and it can be concluded that these sewers do not represent a significant pathway for contaminant migration.

Sediment Sampling

In accordance with Task 1 – Storm Sewer Evaluation of the IRM Work Plan and the NYSDEC correspondence dated June 12, 2000 providing conditional approval of the IRM Work Plan, removal of sediments in the immediate vicinity of the inlet and outfall areas of the 36-inch and 48-inch diameter sewers was conducted. The excavations were conducted to remove potentially contaminated sediments from the inlet and outfall areas, as well as to provide a basin to allow sediment generated during the sewer cleaning to settle at the outfall area.

Sediment samples were collected from the inlet and outfall area of the 36-inch diameter sewer (Inlet A and Outfall #1) and 48-inch diameter sewer (Inlet B and Outfall #2) subsequent to the excavations during November 2000. The samples were analyzed in accordance with NYSDEC ASP procedures for volatile organic compounds (VOCs) using USEPA Method 8260B, semivolatile organic compounds (SVOCs) using USEPA Method 8270C, metals using USEPA Method 6010B, mercury using USEPA Method 7471, and cyanide using USEPA Method 9010B/9014. The samples from Inlets A and B were collected approximately 0.5 ft below the bottom of each sewer pipe. The samples from Outfalls #1 and #2 were collected approximately 1-ft below the bottom of each sewer pipe. A laboratory case narrative for the sediment analytical program is presented in Appendix A.

The VOC, SVOC, and metals (including cyanide) analytical data are provided on attached Tables 1, 2, and 3, respectively and summarized below. The May 1999 analytical data are also presented for comparative purposes.

VOCs

As summarized below, VOCs were detected at both inlet areas (Inlets A and B), and both outfall areas (Outfalls #1 and #2). Regarding the 36-inch sewer, the data indicate that total VOC concentrations were higher at Inlet A (170 µg/Kg) than Outfall #1 (146 µg/Kg). Conversely, the data associated with the 48-inch diameter sewer indicate that the total VOC concentrations at Inlet B (estimated at 2 µg/Kg) were less than at Outfall #2 (1,117.7 µg/Kg). The elevated VOCs detected in the sample from Outfall #2 are likely associated with an observed coal tar-like material in the vicinity of the outfall potentially resulting from operations on the adjacent coke plant site.

Table 2-1 VOC analytical data summary - sediment

| | Inlet A May 1999 | Inlet A November 2000 | Outfall #1 May 1999 | Outfall #1 November 2000 |
|------------------|---------------------|--------------------------|------------------------|-----------------------------|
| Acetone | 47 | 120 | 95 | 55 |
| Benzene | ND | ND | 27 | 27 |
| 2-Butanone | 20 J | 31 | ND | 22 |
| Carbon disulfide | ND | 8 | 2 J | 2 J |
| Ethylbenzene | ND | ND | 15 | 5 |
| Toluene | 400 | ND | 7 J | 4 J |

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| | Inlet A May 1999 | Inlet A November 2000 | Outfall #1 May 1999 | Outfall #1 November 2000 |
|--------------------|---------------------|--------------------------|------------------------|-----------------------------|
| Trichloroethene | ND | 4 J | 13 | 6 |
| Xylene | ND | ND | 31 | 11 |
| cis-1,2-DCE | ND | ND | 35 | 9 |
| trans-1,2-DCE | ND | ND | ND | 2 J |
| Vinyl chloride | ND | ND | ND | 1 J |
| Methylene chloride | 7 J | 7 J | 11 J | 2 J |
| Total VOCs | 474 27 | 170 50 | 236 141 | 146 91 |
| | Inlet B May 1999 | Inlet B November 2000 | Outfall #2 May 1999 | Outfall #2 November 2000 |
| Acetone | 24 | ND | 110 | 28 |
| Benzene | 2 J | ND | 780 | 170 |
| Carbon disulfide | ND | ND | ND | 0.7 J |
| Ethylbenzene | ND | ND | 950 D | 300 |
| Styrene | ND | ND | 900 | 6 |
| Tetrachloroethene | ND | ND | 55 | 1 J |
| Toluene | ND | ND | 900 D | 100 |
| trans-1,2-DCE | ND | ND | 6 J | 5 |
| Trichloroethene | ND | ND | 37 | 14 |
| Vinyl chloride | ND | ND | 14 J | 2 J |
| cis-1,2-DCE | ND | ND | 18 | 10 |
| Xylene | ND | ND | 4,800 D | 480 |
| Methylene chloride | ND | 2 J | ND | 1 J |
| Total VOCs | 26 2 | 2 2 | 8,570 8,460 | 1,117.7 1,089.7 |

Note: Concentrations in $\mu\text{g/L}$

ND – Not detected

J – Estimated concentration

D – Diluted concentration

Source: O'Brien & Gere Engineers, Inc.

SVOCs

SVOCs were detected in both the inlet and outfall samples. The majority of the SVOCs detected are polynuclear aromatic hydrocarbons (PAHs) typically derived from incomplete combustion and tar-

like substances. The PAHs detected in the samples include the following compounds: acenaphthene, acenaphthylene, anthracene, benzo(a)anthracene, benzo(b)fluoranthene, benzo(ghi)perylene, benzo(k)fluoranthene, benzo(a)pyrene, chrysene, dibenzo(a,h)anthracene, fluoranthene, fluorene, indeno(1,2,3-cd)pyrene, naphthalene, phenanthrene, and pyrene. Total PAH concentrations were higher at the outfalls than at the inlets, indicating potential impacts from the adjacent coke plant operations.

Table 2-2 SVOC analytical data summary - sediment

| | Inlet A May 1999 | Inlet A November 2000 | Outfall #1 May 1999 | Outfall #1 November 2000 |
|-------------|---------------------|--------------------------|------------------------|-----------------------------|
| Total PAHs | 64,230 | 41,000 | 204,550 | 1,601,100 |
| Other SVOCs | 1,150 | 4,900 | 4,400 | 17,500 |
| | Inlet B May 1999 | Inlet B November 2000 | Outfall #2 May 1999 | Outfall #2 November 2000 |
| Total PAHs | 55,040 | 75,070 | 3,057,000 | 2,809,900 |
| Other SVOCs | 720 | 2,390 | 172,000 | 52,400 |

Note: Concentrations in $\mu\text{g/L}$

Source: O'Brien & Gere Engineers, Inc.

Metals

Metals were detected in both inlet and outfall samples. In evaluating the metals analytical data, those metals that were greater than one half an order of magnitude higher in the outfall samples, as compared to the inlet samples, were considered site-related. Based on this evaluation, calcium was the only metal constituent that was detected at a concentration greater than one half an order of magnitude in the outfall sample (Outfall #2) compared to the inlet sample (Inlet B). In general, the metals concentrations are higher at the inlets than the outfalls, suggesting an off-site origin. Cyanide was not detected in the inlet or outfall samples.

While the recently collected data provides additional indications of sediment impacts in the vicinity of the inlets and outfalls of the sewers, storm water results from the May 1999 investigation and December 2001 investigation (discussed later in this report) indicate that contaminants in the sediments are not having a significant impact on the quality of storm water discharge. The results show that water quality data for storm water at the inlets and outfalls contained very low concentrations of VOCs, SVOCs, and metals that are orders of magnitude less than the sediment concentrations.

Monitoring Well Installations

Four additional monitoring wells (MW-9, MW-10, MW-11, and MW-12) were installed at the Site during May 2001 to evaluate whether ground water is infiltrating into the 36-inch sewer and to evaluate if contaminants are migrating off-site via ground water. The locations of these monitoring wells, as well as the eight existing monitoring wells, are shown on Figure 2. Each additional monitoring well was constructed using 5-ft long, 2-inch diameter, schedule 40 PVC well screens with 0.010-inch slots flush-threaded to PVC riser casing. Each monitoring well was installed to screen the

fill materials underlying the Site. Ground water elevations for the existing and newly installed monitoring wells are presented in Attachment 1.

The Test Boring Logs contained in Appendix B provide descriptions of the soils encountered at the four monitoring well boring locations. As indicated on the logs, fill thickness encountered in the boring locations ranged from approximately 1.5 ft at MW-11 to 9.5 ft at MW-9 and MW-10.

It should be noted that two attempts were made to install MW-11 closer to the western property line as originally proposed. The initial attempt was advanced to a depth of 14 ft below grade. Dry fill materials were encountered from 0.5 ft to 1.5 ft below grade. Beneath the fill, a dense silt was encountered to 14 ft below grade which did not contain noticeable water. Based on the background geologic information for the area, this silt extends to the top of bedrock and isolates the water-bearing fill materials from the underlying bedrock. The second attempt was advanced to a depth of 10 ft below grade approximately 25 ft to the east of the initial borehole. Similar to the initial attempt, water bearing fill materials were not encountered and the borehole was abandoned. Monitoring well MW-11 was ultimately installed at the location as shown on Figure 2.

Similar to the attempted installations of MW-11, the initial attempt to install MW-12 near the western property line as originally proposed was abandoned. The initial attempt was advanced to a depth of 10 ft below grade with similar soil encountered as at the initial attempt for MW-11. MW-12 was ultimately installed at the location shown on Figure 2. The multiple installation attempts of monitoring wells MW-11 and MW-12 along the western property boundary indicated that the fill thins significantly to the west. As such, the potential for ground water flow off-site to the west in the fill materials is minimal.

Ground Water Sampling and Analytical Data

To evaluate current ground water quality conditions at the Site, ground water samples were collected from monitoring wells MW-1 through MW-11 during June 2001. Monitoring well MW-12 was dry at the time of sampling. The analytical results are presented on attached Tables 4, 5, and 6 and are summarized on the tables below. The samples were collected using low-flow ground water purging and sampling techniques described in the Work Plan. Specific sampling details are presented on the ground water sampling field logs contained in Appendix C. Prior to sample collection, depths to ground water were measured. The depth to water measurements were converted to ground water elevations. Figure 2 illustrates ground water flow based on the June 5, 2001 ground water elevations. As indicated on Figure 2, ground water in the fill materials flows to the west-southwest, which is consistent with historic data.

Ground water samples were analyzed for volatile organic compounds (VOCs) and semivolatile organic compounds (SVOCs) by USEPA method 8260B and 8270B, respectively. In addition, each sample was analyzed for target analyte list (TAL) metals by USEPA series 6000-7000 and cyanide by USEPA method 335. The laboratory analyses were completed by O'Brien & Gere Laboratories, Inc. Quality Assurance/Quality Control (QA/QC) samples analyzed included matrix spike, matrix spike duplicate, blind duplicate, and a trip blank. Category B laboratory deliverables were provided by the laboratory. A laboratory case narrative for the ground water analytical program is presented in Appendix A.

VOCs

The ground water VOC data are presented on Table 4 and summarized below. The May 1999 analytical data is also presented for comparative purposes. The data indicate that VOCs were not detected above ground water standards at existing wells MW-2, MW-4, and MW-5 or newly installed well MW-11. VOCs that were detected above NYS Class GA ground water standards during June 2001 are summarized below and results are compared to the previous ground water sampling event conducted during May 1999.

Table 2-4 VOC analytical data summary – ground water

| | VOCs | NYS Class GA Std | May 1999 | June 2001 |
|-------|------------------------|------------------|----------|-----------|
| MW-1 | Benzene | 1 | 67 D | 46 J |
| | Ethylbenzene | 5 | 28 | 14 J |
| | Toluene | 5 | 29 | 17 J |
| | Xylene | 5 | 200 D | 86 |
| MW-3 | Trichloroethene | 5 | 0.5 UD | 6 |
| | cis-1,2-dichloroethene | 5 | 0.5 UD | 20 |
| MW-6 | Benzene | 1 | 81 D | 79 |
| | Ethylbenzene | 5 | 20 | 17 J |
| | Toluene | 5 | 10 | ND |
| | Xylene | 5 | 54 | 49 J |
| | Methylene chloride | 5 | ND | 14 J |
| MW-7 | Benzene | 1 | 15 | 13 |
| | Xylene | 5 | 6 | 9 |
| MW-8 | 1,1-Dichloroethane | 5 | 9 | 7 |
| MW-9 | Benzene | 1 | NI | 5 |
| | Ethylbenzene | 5 | NI | 7 |
| | Vinyl chloride | 2 | NI | 5 |
| | cis-1,2-dichloroethene | 5 | NI | 8 |
| MW-10 | Benzene | 1 | NI | 16 |

Note: Concentrations in µg/L

NI – Not installed

ND – Not detected

J – Estimated concentration

D – Diluted concentration

Source: O'Brien & Gere Engineers, Inc.

SVOCs

The analytical results for the SVOCs analysis are presented on Table 5 and summarized below. The May 1999 analytical data is also presented for comparative purposes. The data indicate that SVOCs were not detected above Class GA ground water standards at MW-2, MW-5, MW-6, MW-8, or MW-9 during the June 2001 monitoring. The data indicate that SVOCs were detected above NYS Class GA ground water standards at MW-1, MW-3, MW-4, MW-7, MW-10, and MW-11 during the June 2001 monitoring.

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Table 2-5 SVOC analytical data summary – ground water

| | | MW-1 | | MW-3 | | MW-4 | | MW-7 | | MW-10 | | MW-11 | |
|--------------------|---|-----------------------|----------|-----------|----------|-----------|----------|-----------|----------|-----------|-----------|-----------|-----------|
| | | NYS Class GA Standard | May 1999 | June 2001 | June 2001 | June 2001 | June 2001 |
| 2,4-Dimethylphenol | 1 | 17 | 8 | ND | ND | ND | ND | 3 J | ND | ND | ND | ND | ND |
| 2-Methylphenol | 1 | 9 | 4 J | ND | ND | ND | ND | 2 J | ND | ND | ND | ND | ND |
| 4-Methylphenol | 1 | 13 | 7 J | ND | ND | ND | ND | 3 J | 2 J | ND | ND | ND | ND |
| Phenol | 1 | 4 J | 4 J | ND | ND | ND | ND | 2 J | 2 J | 1 J | ND | ND | ND |
| BEHP | 5 | ND | ND | ND | 14 | ND | 10 | 3 J | ND | ND | ND | ND | 12 |

Note: Concentrations in µg/L

ND - not detected

J – Estimated concentration

BEHP – Bis(2-ethylhexyl)phthalate

Source: O'Brien & Gere Engineers, Inc.

Metals

The analytical results of the metals analyses are presented on Table 6. The data indicated that sodium, manganese, and iron were detected in most wells at concentrations that exceed NYS Class GA standards. These elements are likely naturally occurring and are not likely Site related.

Cyanide was detected in most of the wells at concentrations that exceed Class GA standards. Antimony, arsenic, chromium, nickel, and lead were detected with less frequency. Elevated concentrations of these elements were detected in previous sampling events. It is likely that the presence of these metals at elevated levels is Site related. The wells where these elements were detected above standards, and the ranges in concentrations, are summarized below.

Table 2-6 Metals analytical data summary – ground water

| Metal | Well Detected Above Standard | NYS Class GA Std | Range in Conc. (mg/L) |
|-----------|---|------------------|---------------------------------------|
| Antimony | MW-2, MW-5, MW-7, MW-10 | 0.003 | 0.0032 J at MW-10 to 0.0081 J at MW-7 |
| Arsenic | MW-7 | 0.025 | 0.044 |
| Chromium | MW-1, MW-6, and MW-7 | 0.05 | 0.0661 at MW-6 to 0.22 at MW-7 |
| Cyanide | MW-1 through MW-9 | 0.2 | 0.36 at MW-3 to 2.1 at MW-2 |
| Iron | MW-1, MW-2, MW-4, MW-5, MW-6, MW-7, MW-8, MW-9, and MW-10 | 0.3 | 0.549 at MW-5 to 885 at MW-10 |
| Lead | MW-7 | 0.025 | 0.143 |
| Manganese | MW-1 through MW-7, MW-9 and MW-10 | 0.3 | 0.363 at MW-3 to 36.6 at MW-10 |

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| Metal | Well Detected Above Standard | NYS GA Std | Class | Range in Conc. (mg/L) |
|--------|--|---------------|-------|------------------------------|
| Nickel | MW-7 | 0.1 | 0.259 | |
| Sodium | MW-1, MW-2, MW-4, MW-5, MW-6, MW-8, MW-9, MW-10, and MW-11 | 20 | | 23.1 at MW-10 to 131 at MW-8 |

Source: O'Brien & Gere Engineers, Inc.

Storm Water Sampling and Analytical Data

Storm water samples were collected at Inlets A and B and Outfalls #1 and #2 on December 11, 2001. The storm water samples were analyzed for volatile organic compounds (VOCs) and semivolatile organic compounds (SVOCs) by ASP Method 95-1 and 95-2, respectively. In addition, each sample was analyzed for target analyte list metals (TAL) by USEPA series 6000-7000 and cyanide by USEPA method 335. The laboratory analyses were completed by O'Brien & Gere Laboratories, Inc. Quality Assurance/Quality Control (QA/QC) samples analyzed included matrix spike, matrix spike duplicate, blind duplicate, and a trip blank. Category B laboratory deliverables were provided by the laboratory. A laboratory case narrative for the storm water analytical program is presented in Appendix A.

VOCs

The storm water VOC data are presented on Table 7 and summarized below. The May 1999 analytical data are also presented for comparative purposes. The data indicate that other than methylene chloride, VOCs were not detected in the Inlet samples. Methylene chloride was detected in the QC trip blank and its detection in the Inlet samples is likely not site-related. VOCs detected in the Outfall samples are similar in both compounds and concentrations as compared to the May 1999 data.

Table 2-7 VOC analytical data summary – storm water

| | Inlet A May 1999 | Inlet A December 2001 | Outfall #1 May 1999 | Outfall #1 December 2001 |
|--------------------|---------------------|--------------------------|------------------------|-----------------------------|
| Acetone | 2 J | ND | 4 J | 3 J |
| Benzene | ND | ND | 6 | 6 J |
| Carbon disulfide | ND | ND | 0.2 J | 0.8 J |
| Ethylbenzene | ND | ND | 0.4 J | 0.6 J |
| Toluene | 0.1 J | ND | 0.9 | 2 J |
| Trichloroethene | ND | ND | 0.9 | ND |
| Xylene | ND | ND | 2 | 2 J |
| cis-1,2-DCE | ND | ND | 8 | 0.8 J |
| Methylene chloride | ND | 1 J | ND | 1 J |

| | Inlet B May 1999 | Inlet B December 2001 | Outfall #2 May 1999 | Outfall #2 December 2001 |
|-----------------------|---------------------|--------------------------|------------------------|-----------------------------|
| Acetone | 4 J | ND | 6 J | 3 J |
| Bromodichloromethane | ND | ND | 0.2 J | ND |
| Chloroform | ND | ND | 0.3 J | ND |
| Tetrachloroethene | ND | ND | 0.4 J | ND |
| 1,1,1-Trichloroethane | ND | ND | 0.1 J | ND |
| Trichloroethene | ND | ND | 0.2 J | ND |
| Vinyl chloride | ND | ND | 0.5 J | 0.5 J |
| cis-1,2-DCE | ND | ND | 1 | 0.7 J |
| Methylene chloride | ND | 1 J | ND | 1 J |

Note: Concentrations in µg/L

ND – Not detected

J – Estimated concentration

Source: O'Brien & Gere Engineers, Inc.

SVOCs

The storm water SVOC data are presented on Table 8. The data indicate that SVOCs were not detected in the Inlet samples. Naphthalene (57 µg/L), and estimated concentrations of 2-methylnaphthalene (6 µg/L), fluorene (2 µg/L), and phenanthrene (2 µg/L) were detected in the Outfall #1 sample. Naphthalene was detected in the Outfall #2 sample at a concentration of 22 µg/L. No other SVOCs were detected in the Outfall #2 sample.

Metals

The storm water metals data are presented on Table 9. To distinguish whether metals detected in the storm water Outfall samples are site-related or due to natural variations, those metal constituents that were detected at concentrations equal to or greater than one-half an order of magnitude, or detected in the Outfalls but not the Inlets, were identified. The following metal constituents were identified as fitting these criteria: iron, silver, cyanide for Outfall #1, and antimony and lead for Outfall #2. It should be noted that silver was detected in the Outfall #1 sample at an estimated concentration at the detection limit (1 µg/L).

SECTION 3 – CONCLUSIONS

Storm Sewer Evaluation

The hydraulic flushing and video inspection indicated that the 36-inch and 48-inch storm sewers contain a large amount of debris, making further investigatory work impractical without substantial rehabilitation of the pipelines. Based on the work completed, no significant sources of groundwater infiltration into the sewers were observed. The portions of the two sewers that were televised (20% of the 36-inch pipe and 50% of the 48-inch pipe) are reasonably representative. It can be concluded

that these sewers do not represent a significant pathway for contaminant migration. While the invert elevations of the storm sewer system are below the water table and the storm sewers may be receiving minimal amounts of ground water via infiltration, the concentrations are relatively low as indicated by the storm water sample data. In addition, the storm water sample data indicate that impacted sediments in the Inlet and Outfall areas have minimal impact on storm water quality.

Storm Sewer Sediment

In summary, the sediment data associated with the 36-inch storm sewer indicated higher concentrations of total VOCs at Inlet A than were detected at Outfall #1. Conversely, at the 48-inch storm sewer, higher concentrations of total VOCs were detected in Outfall #2 than were detected in Inlet B. SVOCs were detected in both the upstream and downstream sediment samples. The highest concentrations of SVOCs were detected Outfall #2. The elevated BTEX and SVOC concentrations at Outfall# 2 are likely attributed to the observed coal tar-like substance in that area and could potentially be related to the adjacent coke plant.

A comparison between VOC, SVOC, and metals analytical results for sediment samples collected during May 1999 as part of a voluntary site investigation (O'Brien & Gere Engineers, Inc., letter report dated November 9, 1999), and the recent sediment sampling (November 2000), is provided on Table 4. The data shown on Table 4 is summarized as follows:

- Total VOC concentrations were lower during November 2000 compared to May 1999.
- Total PAH concentrations were lower during November 2000 compared to May 1999 at Inlet A and Outfall #2, and higher at Inlet B and Outfall #1.
- The total concentration of other SVOCs was generally higher during November 2000 compared to May 1999, with the exception of Outfall #2.
- Metals concentrations that were greater than one half an order of magnitude in the outfall samples compared to the inlet samples were considered to be site-related. During the May 1999 sampling, concentrations of calcium at Outfall #2, mercury at Outfalls #1 and #2, and cyanide at Outfalls #1 and #2 were considered to be site-related. During the November 2000 sampling, the concentrations of mercury was lower at Inlet A, Outfalls #1 and #2, and higher at Inlet B compared to the May 1999 data, and cyanide was not detected.

Ground Water

VOC constituents detected in ground water above NYS Class GA ground water standards consist mainly of benzene, ethylbenzene, toluene, and xylene. One or more of these compounds were detected above standards in wells MW-1, MW-6, MW-7, MW-9, and MW-10. Chlorinated VOCs were detected in ground water with less frequency. 1,1-dichloroethane (detected in MW-8), trichloroethene (detected in MW-3), and cis-1,2-dichloroethene (detected in MW-3 and MW-9) were the only chlorinated VOCs detected in ground water at concentrations above standards. Low concentrations of SVOCs were detected above standards at monitoring wells MW-1, MW-3, MW-4,

Mr. Bruce R. Terbush, P.E.
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MW-7, MW-10, and MW-11. Consistent with previous data, chromium, cyanide, and lead were detected in most wells at concentrations that exceed NYS Class GA standards.

The VOC, SVOC, and metals concentrations detected in ground water are not indicative of a continuing source but rather residual contamination.

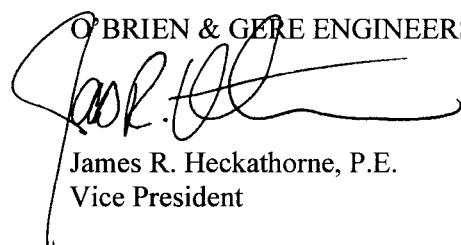
Storm Water

Based on the analytical data, concentrations of VOCs, SVOCs, or metals in the residual sediments in the sewer pipes and the Inlet and Outfall areas do not have a significant impact on storm water quality. Likewise, potential infiltration of ground water carrying concentrations of VOCs, SVOCs, and metals does not have a significant impact on storm water quality. The storm water sample data indicates that small concentrations of VOCs may be carried off-site via storm water but at concentrations that do not exceed New York State Ambient Water Quality Standards (NYSDEC, 1998). The analytical data indicates that migration of SVOCs and metals via storm water are not significant.

If you have any questions regarding the information presented herein, please contact me at your earliest convenience at (315) 437-6100.

Very truly yours,

O'BRIEN & GERE ENGINEERS, INC.



James R. Heckathorne, P.E.
Vice President

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cc: Maria Kaouris – Honeywell
David Flynn, Esq. – Phillips, Lytle, Hitchcock, Blaine & Huber, LLP
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Deborah Y. Wright – O'Brien & Gere Engineers, Inc.



O'BRIEN & GERE
ENGINEERS, INC.
Honeywell
Tonawanda, NY

Volatile Organic Compound Data

Storm Sediment Samples

| Compound | Sample ID | Outfall #1 | Outfall #1 RE | Outfall #1 | Outfall #1 DCP | Outfall #2 | Outfall #2 DL | Outfall #2 |
|-----------------------------|-----------|------------|---------------|------------|----------------|------------|---------------|------------|
| | | | | | | | | |
| | | | | | | | | |
| 1,1,1-Trichloroethane | 11 UD | 11 UD | 11 UD | 5 UD | 5 UD | 14 UD | 290 UD | 3 UD |
| 1,1,2,2-Tetrachloroethane | 11 UD | 11 UD | 11 UD | 5 UD | 5 UD | 14 UD | 290 UD | 3 UD |
| 1,1,2-Trichloroethane | 11 UD | 11 UD | 11 UD | 5 UD | 5 UD | 14 UD | 290 UD | 3 UD |
| 1,1-Dichloroethane | 11 UD | 11 UD | 11 UD | 5 UD | 5 UD | 14 UD | 290 UD | 3 UD |
| 1,1-Dichloroethene | 11 UD | 11 UD | 11 UD | 5 UD | 5 UD | 14 UD | 290 UD | 3 UD |
| 1,2-Dichloroethane | 11 UD | 11 UD | 11 UD | 5 UD | 5 UD | 14 UD | 290 UD | 3 UD |
| 1,2-Dichloropropane | 11 UD | 11 UD | 11 UD | 5 UD | 5 UD | 14 UD | 290 UD | 3 UD |
| 2-Butanone (MEK) | 43 UD | 43 UD | 22 | 21 J | 57 UD | 1100 UD | 13 UD | |
| 2-Hexanone | 22 UD | 22 UD | 9 UD | 10 UD | 29 UD | 570 UD | 6 UD | |
| 4-Methyl-2-pentanone (MIBK) | 22 UD | 22 UD | 9 UD | 10 UD | 29 UD | 570 UD | 6 UD | |
| Acetone | 95 | 66 | 55 | 51 | 110 | 1100 UD | 28 | |
| Benzene | 27 | 16 | 27 | 22 | 780 | 300 D | 170 | |
| Bromodichloromethane | 11 UD | 11 UD | 5 UD | 5 UD | 14 UD | 290 UD | 3 UD | |
| Bromoform | 11 UD | 11 UD | 5 UD | 5 UD | 14 UD | 290 UD | 3 UD | |
| Bromomethane | 22 UD | 22 UD | 9 UD | 10 UD | 29 UD | 570 UD | 6 UD | |
| Carbon disulfide | 2 J | 11 UD | 2 J | 2 J | 14 UD | 290 UD | 0.7 J | |
| Carbon tetrachloride | 11 UD | 11 UD | 5 UD | 5 UD | 14 UD | 290 UD | 3 UD | |
| Chlorobenzene | 11 UD | 11 UD | 5 UD | 5 UD | 14 UD | 290 UD | 3 UD | |
| Chloroethane | 22 UD | 22 UD | 9 UD | 10 UD | 29 UD | 570 UD | 6 UD | |
| Chloroform | 11 UD | 11 UD | 5 UD | 5 UD | 14 UD | 290 UD | 3 UD | |
| Chloromethane | 22 UD | 22 UD | 9 UD | 10 UD | 29 UD | 570 UD | 6 UD | |
| Dibromochloromethane | 11 UD | 11 UD | 5 UD | 5 UD | 14 UD | 290 UD | 3 UD | |
| Ethene, 1,2-dichloro-(E)- | 11 UD | 11 UD | 2 J | 2 J | 6 J | 290 UD | 5 | |
| Ethybenzene | 15 | 7 J | 5 | 5 J | 2400 E | 950 D | 300 | |
| Methylene chloride | 22 UD | 11 J | 2 J | 2 J | 29 UD | 570 UD | 1 J | |
| Styrene | 11 UD | 11 UD | 5 UD | 5 UD | 900 | 51 JD | 6 | |
| Tetrachloroethene | 11 UD | 11 UD | 5 UD | 5 UD | 55 | 290 UD | 1 J | |
| Toluene | 7 J | 6 J | 4 J | 3 J | 2200 E | 900 D | 100 | |
| Trichloroethene | 13 | 5 J | 6 | 5 J | 37 | 290 UD | 14 | |
| Vinyl chloride | 22 UD | 22 UD | 1 J | 1 J | 14 J | 570 UD | 2 J | |
| Xylene (total) | 17 | 31 | 11 | 9 | 1100 E | 4800 D | 480 | |
| cis-1,2-Dichloroethylene | 35 | 22 | 9 | 8 | 18 | 290 UD | 10 | |
| cis-1,3-Dichloropropene | 11 UD | 11 UD | 5 UD | 5 UD | 14 UD | 290 UD | 3 UD | |
| trans-1,3-Dichloropropene | 11 UD | 11 UD | 5 UD | 5 UD | 14 UD | 290 UD | 3 UD | |

NOTES: U - Not detected, J - Estimated value, E - outside linear range, D - from diluted analysis, --- - Not analyzed



Table 1
Honeywell
Tonawanda, NY
Volatile Organic Compound Data
Storm Sediment Samples

| Compound | Sample ID | Storm Inlet A | | Storm Inlet A RE | | Storm Inlet B | | Storm Inlet B RE | | Storm Inlet B | |
|-----------------------------|-----------|---------------|----------|------------------|----------|---------------|----------|------------------|----------|---------------|----------|
| | | Sample Date | 05/11/99 | 05/11/99 | 11/15/00 | 05/11/99 | 05/11/99 | 05/11/99 | 05/11/99 | 05/11/99 | 11/15/00 |
| | | | ug/Kg | | ug/Kg | | | | | | |
| 1,1,1-Trichloroethane | | 6 UD | 6 UD | 6 UD | 6 UD | 4 UD | 4 UD | 4 UD | 4 UD | 7 UD | 7 UD |
| 1,1,2,2-Tetrachloroethane | | 6 UD | 6 UD | 6 UD | 6 UD | 4 UD | 4 UD | 4 UD | 4 UD | 7 UD | 7 UD |
| 1,1,2-Trichloroethane | | 6 UD | 6 UD | 6 UD | 6 UD | 4 UD | 4 UD | 4 UD | 4 UD | 7 UD | 7 UD |
| 1,1-Dichloroethane | | 6 UD | 6 UD | 6 UD | 6 UD | 4 UD | 4 UD | 4 UD | 4 UD | 7 UD | 7 UD |
| 1,1-Dichloroethene | | 6 UD | 6 UD | 6 UD | 6 UD | 4 UD | 4 UD | 4 UD | 4 UD | 7 UD | 7 UD |
| 1,2-Dichloroethane | | 6 UD | 6 UD | 6 UD | 6 UD | 4 UD | 4 UD | 4 UD | 4 UD | 7 UD | 7 UD |
| 1,2-Dichloropropane | | 6 UD | 20 J | 23 UD | 31 | 18 UD | 18 UD | 18 UD | 18 UD | 28 UD | 28 UD |
| 2-Butanone (MEK) | | 12 UD | 12 UD | 12 UD | 11 UD | 9 UD | 9 UD | 9 UD | 9 UD | 14 UD | 14 UD |
| 2-Hexanone | | 12 UD | 12 UD | 12 UD | 11 UD | 9 UD | 9 UD | 9 UD | 9 UD | 14 UD | 14 UD |
| 4-Methyl-2-pentanone (MBK) | | 47 | 35 | 120 | 24 | 24 | 24 | 24 | 24 | 28 UD | 28 UD |
| Acetone | | | | | | 2 J | 1 J | 1 J | 1 J | 7 UD | 7 UD |
| Benzene | | 6 UD | 6 UD | 6 UD | 6 UD | 4 UD | 4 UD | 4 UD | 4 UD | 7 UD | 7 UD |
| Bromodichloromethane | | 6 UD | 6 UD | 6 UD | 6 UD | 4 UD | 4 UD | 4 UD | 4 UD | 7 UD | 7 UD |
| Bromoform | | 6 UD | 6 UD | 6 UD | 6 UD | 4 UD | 4 UD | 4 UD | 4 UD | 7 UD | 7 UD |
| Bromomethane | | 12 UD | 12 UD | 12 UD | 11 UD | 9 UD | 9 UD | 9 UD | 9 UD | 14 UD | 14 UD |
| Carbon disulfide | | 6 UD | 6 UD | 6 UD | 8 | 4 UD | 4 UD | 4 UD | 4 UD | 7 UD | 7 UD |
| Carbon tetrachloride | | 6 UD | 6 UD | 6 UD | 6 UD | 4 UD | 4 UD | 4 UD | 4 UD | 7 UD | 7 UD |
| Chlorobenzene | | 6 UD | 6 UD | 6 UD | 6 UD | 4 UD | 4 UD | 4 UD | 4 UD | 7 UD | 7 UD |
| Chloroethane | | 12 UD | 6 UD | 6 UD | 6 UD | 4 UD | 4 UD | 4 UD | 4 UD | 7 UD | 7 UD |
| Chloroform | | 6 UD | 12 UD | 12 UD | 11 UD | 9 UD | 9 UD | 9 UD | 9 UD | 14 UD | 14 UD |
| Chloronethane | | 12 UD | 12 UD | 12 UD | 11 UD | 9 UD | 9 UD | 9 UD | 9 UD | 14 UD | 14 UD |
| Dibromochloromethane | | 6 UD | 6 UD | 6 UD | 6 UD | 4 UD | 4 UD | 4 UD | 4 UD | 7 UD | 7 UD |
| Ethene, 1,2-dichloro-, (E)- | | 6 UD | 6 UD | 6 UD | 6 UD | 4 UD | 4 UD | 4 UD | 4 UD | 7 UD | 7 UD |
| Ethybenzene | | 6 UD | 7 J | 7 J | 9 UD | 9 UD | 9 UD | 9 UD | 9 UD | 7 UD | 7 UD |
| Methylene chloride | | 3 J | | | | | | | | 2 J | 2 J |
| Styrene | | 6 UD | 6 UD | 6 UD | 6 UD | 4 UD | 4 UD | 4 UD | 4 UD | 7 UD | 7 UD |
| Tetrachloroethene | | 6 UD | 6 UD | 6 UD | 6 UD | 4 UD | 4 UD | 4 UD | 4 UD | 7 UD | 7 UD |
| Toluene | | 400 | 120 | 6 UD | 4 UD | 4 UD | 4 UD | 4 UD | 4 UD | 7 UD | 7 UD |
| Trichloroethene | | 6 UD | 6 UD | 4 J | 4 UD | 4 UD | 4 UD | 4 UD | 4 UD | 7 UD | 7 UD |
| Vinyl chloride | | 12 UD | 12 UD | 11 UD | 9 UD | 9 UD | 9 UD | 9 UD | 9 UD | 14 UD | 14 UD |
| Xylene (total) | | 6 UD | 6 UD | 6 UD | 4 UD | 4 UD | 4 UD | 4 UD | 4 UD | 7 UD | 7 UD |
| cis-1,2-Dichloroethene | | 6 UD | 6 UD | 6 UD | 6 UD | 4 UD | 4 UD | 4 UD | 4 UD | 7 UD | 7 UD |
| cis-1,3-Dichloropropene | | 6 UD | 6 UD | 6 UD | 6 UD | 4 UD | 4 UD | 4 UD | 4 UD | 7 UD | 7 UD |
| trans-1,3-Dichloropropene | | 6 UD | 6 UD | 6 UD | 6 UD | 4 UD | 4 UD | 4 UD | 4 UD | 7 UD | 7 UD |

NOTES: U - Not detected, J - Estimated value, E - outside linear range, D - from diluted analysis, --- - Not analyzed.



Table 2
Honeywell
Tonawanda, NY
Semivolatile Organic Compound Data

| Compound | Sample ID | Outfall #1 | Outfall #1 RE | Storm Sediment Samples | | | Outfall #2 DL | Outfall #2 RE |
|-----------------------------|-----------|------------|---------------|------------------------|------------|-----------|---------------|---------------|
| | | | | Sample Date | Outfall #1 | | | |
| | | | | | 11/15/00 | ug/Kg | Outfall #2 | 05/10/99 |
| Compound | Sample ID | Outfall #1 | Outfall #1 RE | Outfall #1 | 11/15/00 | ug/Kg | Outfall #2 | 05/10/99 |
| | | Units | ug/Kg | Units | ug/Kg | ug/Kg | ug/Kg | ug/Kg |
| 1,2,4-Trichlorobenzene | | 7200 UD | 7200 UD | 31000 UD | 35000 UD | 38000 UD | 76000 UD | 38000 UD |
| 1,2-Dichlorobenzene | | 7200 UD | 7200 UD | 31000 UD | 35000 UD | 38000 UD | 76000 UD | 38000 UD |
| 1,3-Dichlorobenzene | | 7200 UD | 7200 UD | 31000 UD | 35000 UD | 38000 UD | 76000 UD | 38000 UD |
| 1,4-Dichlorobenzene | | 7200 UD | 7200 UD | 160000 UD | 170000 UD | 190000 UD | 380000 UD | 190000 UD |
| 2,4,5-Trichlorophenol | | 36000 UD | 36000 UD | 31000 UD | 35000 UD | 38000 UD | 76000 UD | 38000 UD |
| 2,4,6-Trichlorophenol | | 7200 UD | 7200 UD | 31000 UD | 35000 UD | 38000 UD | 76000 UD | 38000 UD |
| 2,4-Dichlorophenol | | 7200 UD | 7200 UD | 31000 UD | 35000 UD | 38000 UD | 76000 UD | 38000 UD |
| 2,4-Dimethylphenol | | 7200 UD | 7200 UD | 160000 UD | 170000 UD | 190000 UD | 380000 UD | 190000 UD |
| 2,4-Dinitrophenol | | 36000 UD | 36000 UD | 31000 UD | 35000 UD | 38000 UD | 76000 UD | 38000 UD |
| 2,4-Dinitrotoluene | | 7200 UD | 7200 UD | 31000 UD | 35000 UD | 38000 UD | 76000 UD | 38000 UD |
| 2-Chloronaphthalene | | 7200 UD | 7200 UD | 31000 UD | 35000 UD | 38000 UD | 76000 UD | 38000 UD |
| 2-Chlorophenol | | 7200 UD | 7200 UD | 31000 UD | 35000 UD | 38000 UD | 76000 UD | 38000 UD |
| 2-Methylnaphthalene | 750 J | 7200 UD | 7200 UD | 34000 J | 35000 UD | 71000 | 73000 JD | 70000 |
| 2-Methylphenol | | 7200 UD | 7200 UD | 31000 UD | 35000 UD | 38000 UD | 76000 UD | 38000 UD |
| 2-Nitroaniline | | 36000 UD | 36000 UD | 160000 UD | 170000 UD | 190000 UD | 380000 UD | 190000 UD |
| 2-Nitrophenol | | 7200 UD | 7200 UD | 31000 UD | 35000 UD | 38000 UD | 76000 UD | 38000 UD |
| 3,3-Dichlorobenzidine | | 14000 UD | 14000 UD | 63000 UD | 69000 UD | 76000 UD | 150000 UD | 76000 UD |
| 3-Nitroaniline | | 36000 UD | 36000 UD | 160000 UD | 170000 UD | 190000 UD | 380000 UD | 190000 UD |
| 4,6-Dinitro-2-methylphenol | | 36000 UD | 36000 UD | 160000 UD | 170000 UD | 190000 UD | 380000 UD | 190000 UD |
| 4-Bromophenyl phenyl ether | | 7200 UD | 7200 UD | 31000 UD | 35000 UD | 38000 UD | 76000 UD | 38000 UD |
| 4-Chloro-3-methylphenol | | 7200 UD | 7200 UD | 31000 UD | 35000 UD | 38000 UD | 76000 UD | 38000 UD |
| 4-Chloraniline | | 7200 UD | 7200 UD | 31000 UD | 35000 UD | 38000 UD | 76000 UD | 38000 UD |
| 4-Chlorophenyl phenyl ether | | 7200 UD | 7200 UD | 31000 UD | 35000 UD | 38000 UD | 76000 UD | 38000 UD |
| 4-Methylphenol | | 7200 UD | 7200 UD | 31000 UD | 35000 UD | 38000 UD | 76000 UD | 38000 UD |
| 4-Nitroaniline | | 36000 UD | 36000 UD | 160000 UD | 170000 UD | 190000 UD | 380000 UD | 190000 UD |
| 4-Nitrophenol | | 36000 UD | 36000 UD | 160000 UD | 170000 UD | 190000 UD | 380000 UD | 190000 UD |
| Acenaphthene | | 1800 J | 1800 J | 8700 J | 35000 UD | 38000 UD | 76000 UD | 38000 UD |
| Acenaphthylene | | 1500 J | 1400 J | 23000 J | 17000 J | 68000 | 73000 JD | 67000 |
| Anthracene | | 6200 J | 6200 J | 43000 J | 34000 J | 140000 | 150000 D | 140000 |
| Benzo(a)anthracene | | 14000 | 14000 | 130000 | 93000 | 130000 | 140000 D | 140000 |
| Benzo(b)fluoranthene | | 20000 | 18000 | 170000 | 120000 | 170000 | 160000 D | 170000 |
| Benzo(ghi)perylene | | 8800 | 11000 | 53000 | 40000 | 74000 | 50000 JD | 89000 |
| Benzo(k)fluoranthene | | 5900 J | 7100 J | 52000 | 37000 | 62000 | 71000 JD | 58000 |
| Benzo[a]pyrene | | 14000 | 15000 | 120000 | 91000 | 130000 | 130000 D | 140000 |

NOTES: U - Not detected, J - Estimated value, E - outside linear range, D - from diluted analysis, --- - Not analyzed.



Table 2
Honeywell
Tonawanda, NY
Semivolatile Organic Compound Data

| Compound | Sample ID | Outfall #1 | Outfall #1 RE | Outfall #1 | Outfall #1 DUP | Outfall #1 | Outfall #2 | Outfall #2 DL | Outfall #2 RE |
|-----------------------------------|-----------|------------|---------------|------------|----------------|------------|------------|---------------|---------------|
| | | | | | | | | | |
| | | | | | | | | | |
| Bis(2-chloroethoxy)methane | 7200 UD | 7200 UD | 31000 UD | 35000 UD | 38000 UD | 76000 UD | 38000 UD | 38000 UD | 38000 UD |
| Bis(2-chloroethyl)ether | 7200 UD | 7200 UD | 31000 UD | 35000 UD | 38000 UD | 76000 UD | 38000 UD | 38000 UD | 38000 UD |
| Bis(2-chloroisopropyl) ether | 7200 UD | 7200 UD | 31000 UD | 35000 UD | 38000 UD | 76000 UD | 38000 UD | 38000 UD | 38000 UD |
| Bis(2-ethylhexyl)phthalate (BEHP) | 1700 J | 2000 J | 31000 UD | 35000 UD | 38000 UD | 76000 UD | 38000 UD | 38000 UD | 38000 UD |
| Butyl benzyl phthalate | 7200 UD | 7200 UD | 31000 UD | 35000 UD | 38000 UD | 76000 UD | 38000 UD | 38000 UD | 38000 UD |
| Carbazole | 1300 J | 1200 J | 3800 J | 4300 J | 62000 | 66000 JD | 18000 D | 160000 | 160000 |
| Chrysene | 16000 | 17000 | 110000 | 79000 | 170000 | 180000 D | 38000 UD | 38000 UD | 38000 UD |
| Di-n-butyl phthalate | 7200 UD | 7200 UD | 31000 UD | 35000 UD | 38000 UD | 76000 UD | 38000 UD | 38000 UD | 38000 UD |
| Di-n-octyl phthalate | 7200 UD | 7200 UD | 31000 UD | 35000 UD | 38000 UD | 76000 UD | 38000 UD | 38000 UD | 38000 UD |
| Dibenz(a,h)anthracene | 2200 J | 2500 J | 14000 J | 16000 J | 18000 J | 14000 JD | 21000 J | 110000 | 110000 |
| Dibenzofuran | 1100 J | 1100 J | 8700 J | 8800 J | 110000 | 120000 D | 38000 UD | 38000 UD | 38000 UD |
| Diethyl phthalate | 7200 UD | 7200 UD | 31000 UD | 35000 UD | 38000 UD | 76000 UD | 38000 UD | 38000 UD | 38000 UD |
| Dimethyl phthalate | 7200 UD | 7200 UD | 31000 UD | 35000 UD | 38000 UD | 76000 UD | 38000 UD | 38000 UD | 38000 UD |
| Fluoranthene | 27000 | 25000 | 290000 | 230000 | 320000 E | 410000 D | 180000 D | 160000 | 160000 |
| Fluorene | 2700 J | 2600 J | 24000 J | 20000 J | 170000 | 170000 | 380000 UD | 380000 UD | 380000 UD |
| Hexachlorobenzene | 7200 UD | 7200 UD | 31000 UD | 35000 UD | 38000 UD | 76000 UD | 38000 UD | 38000 UD | 38000 UD |
| Hexachlorobutadiene | 7200 UD | 7200 UD | 31000 UD | 35000 UD | 38000 UD | 76000 UD | 38000 UD | 38000 UD | 38000 UD |
| Hexachlorocyclopentadiene | 7200 UD | 7200 UD | 31000 UD | 35000 UD | 38000 UD | 76000 UD | 38000 UD | 38000 UD | 38000 UD |
| Hexachloroethane | 7200 UD | 7200 UD | 31000 UD | 35000 UD | 38000 UD | 76000 UD | 38000 UD | 38000 UD | 38000 UD |
| Indeno(1,2,3-cd)pyrene | 7600 | 9100 | 53000 | 39000 | 63000 | 48000 JD | 75000 | 75000 | 75000 |
| Isophorone | 7200 UD | 7200 UD | 31000 UD | 35000 UD | 38000 UD | 76000 UD | 38000 UD | 38000 UD | 38000 UD |
| N-Nitrosodiphenylamine | 7200 UD | 7200 UD | 31000 UD | 35000 UD | 38000 UD | 76000 UD | 38000 UD | 38000 UD | 38000 UD |
| N-Nitrosodipropylamine | 7200 UD | 7200 UD | 31000 UD | 35000 UD | 38000 UD | 76000 UD | 38000 UD | 38000 UD | 38000 UD |
| Naphthalene | 4900 J | 4900 J | 17000 J | 14000 J | 480000 E | 510000 D | 480000 E | 480000 E | 480000 E |
| Nitrobenzene | 7200 UD | 7200 UD | 31000 UD | 35000 UD | 38000 UD | 76000 UD | 38000 UD | 38000 UD | 38000 UD |
| Pentachlorophenol | 360000 UD | 360000 UD | 160000 UD | 170000 UD | 190000 UD | 380000 UD | 190000 UD | 190000 UD | 190000 UD |
| Phenanthrene | 17000 | 17000 | 190000 | 170000 | 470000 E | 520000 D | 470000 E | 470000 E | 470000 E |
| Phenol | 7200 UD | 7200 UD | 310000 UD | 350000 UD | 380000 UD | 760000 UD | 380000 D | 380000 UD | 380000 UD |
| Pyrene | 38000 | 47000 | 300000 | 230000 | 400000 E | 300000 D | 490000 E | 490000 E | 490000 E |

NOTES: U - Not detected, J - Estimated value, E - outside linear range, D - from diluted analysis, --- - Not analyzed



Table 2
Honeywell
Tonawanda, NY
Semivolatile Organic Compound Data

| Compound | Sample ID | Outfall #2 | Storm Sediment Samples | | | Storm Inlet A RE | Storm Inlet A DL | Storm Inlet A | Storm Inlet B | |
|-----------------------------|-----------|------------|------------------------|---------------|---------|------------------|------------------|---------------|---------------|---------|
| | | | Sample Date | Storm Inlet A | | | | | | |
| | | | | 05/11/99 | ug/Kg | ug/Kg | ug/Kg | ug/Kg | ug/Kg | |
| 1,2,4-Trichlorobenzene | | 42000 UD | 1500 UD | 3100 UD | 1500 UD | 15000 UD | 15000 UD | 15000 UD | 2900 UD | 920 UD |
| 1,2-Dichlorobenzene | | 42000 UD | 1500 UD | 3100 UD | 1500 UD | 15000 UD | 15000 UD | 15000 UD | 2900 UD | 920 UD |
| 1,3-Dichlorobenzene | | 42000 UD | 1500 UD | 3100 UD | 1500 UD | 15000 UD | 15000 UD | 15000 UD | 2900 UD | 920 UD |
| 1,4-Dichlorobenzene | | 42000 UD | 1500 UD | 3100 UD | 1500 UD | 15000 UD | 15000 UD | 15000 UD | 2900 UD | 920 UD |
| 2,4,5-Trichlorophenol | | 210000 UD | 7800 UD | 16000 UD | 7800 UD | 74000 UD | 15000 UD | 15000 UD | 4600 UD | 4600 UD |
| 2,4,6-Trichlorophenol | | 42000 UD | 1500 UD | 3100 UD | 1500 UD | 15000 UD | 15000 UD | 15000 UD | 2900 UD | 920 UD |
| 2,4-Dichlorophenol | | 42000 UD | 1500 UD | 3100 UD | 1500 UD | 15000 UD | 15000 UD | 15000 UD | 2900 UD | 920 UD |
| 2,4-Dimethylphenol | | 42000 UD | 1500 UD | 3100 UD | 1500 UD | 15000 UD | 15000 UD | 15000 UD | 2900 UD | 920 UD |
| 2,4-Dinitrophenol | | 210000 UD | 7800 UD | 16000 UD | 7800 UD | 74000 UD | 15000 UD | 15000 UD | 4600 UD | 4600 UD |
| 2,4-Dinitrotoluene | | 42000 UD | 1500 UD | 3100 UD | 1500 UD | 15000 UD | 15000 UD | 15000 UD | 2900 UD | 920 UD |
| 2-Chloronaphthalene | | 42000 UD | 1500 UD | 3100 UD | 1500 UD | 15000 UD | 15000 UD | 15000 UD | 2900 UD | 920 UD |
| 2-Chlorophenol | | 42000 UD | 1500 UD | 3100 UD | 1500 UD | 15000 UD | 15000 UD | 15000 UD | 2900 UD | 920 UD |
| 2-Methylnaphthalene | | 7900 J | 1500 UD | 3100 UD | 1500 UD | 15000 UD | 15000 UD | 15000 UD | 2900 UD | 520 J |
| 2-Methylphenol | | 42000 UD | 1500 UD | 3100 UD | 1500 UD | 15000 UD | 15000 UD | 15000 UD | 2900 UD | 920 UD |
| 2-Nitroaniline | | 210000 UD | 7800 UD | 16000 UD | 7800 UD | 74000 UD | 15000 UD | 15000 UD | 4600 UD | 4600 UD |
| 2-Nitrophenol | | 42000 UD | 1500 UD | 3100 UD | 1500 UD | 15000 UD | 15000 UD | 15000 UD | 2900 UD | 920 UD |
| 3,3-Dichlorobenzidine | | 30000 UD | 6200 UD | 3100 UD | 6200 UD | 3100 UD | 30000 UD | 30000 UD | 5800 UD | 1900 UD |
| 3-Nitroaniline | | 210000 UD | 7800 UD | 16000 UD | 7800 UD | 74000 UD | 15000 UD | 15000 UD | 4600 UD | 4600 UD |
| 4,6-Dinitro-2-methylphenol | | 210000 UD | 7800 UD | 16000 UD | 7800 UD | 74000 UD | 15000 UD | 15000 UD | 4600 UD | 4600 UD |
| 4-Bromophenyl phenyl ether | | 42000 UD | 1500 UD | 3100 UD | 1500 UD | 15000 UD | 15000 UD | 15000 UD | 2900 UD | 920 UD |
| 4-Chloro-3-methylphenol | | 42000 UD | 1500 UD | 3100 UD | 1500 UD | 15000 UD | 15000 UD | 15000 UD | 2900 UD | 920 UD |
| 4-Chloroaniline | | 42000 UD | 1500 UD | 3100 UD | 1500 UD | 15000 UD | 15000 UD | 15000 UD | 2900 UD | 920 UD |
| 4-Chlorophenyl phenyl ether | | 42000 UD | 1500 UD | 3100 UD | 1500 UD | 15000 UD | 15000 UD | 15000 UD | 2900 UD | 920 UD |
| 4-Methylphenol | | 42000 UD | 1500 UD | 3100 UD | 1500 UD | 15000 UD | 15000 UD | 15000 UD | 2900 UD | 920 UD |
| 4-Nitroaniline | | 210000 UD | 7800 UD | 16000 UD | 7800 UD | 74000 UD | 15000 UD | 15000 UD | 4600 UD | 4600 UD |
| 4-Nitrophenol | | 210000 UD | 7800 UD | 16000 UD | 7800 UD | 74000 UD | 15000 UD | 15000 UD | 4600 UD | 4600 UD |
| Acenaphthene | | 32000 J | 310 J | 31000 UD | 300 J | 5000 UD | 2900 UD | 2900 UD | 470 J | |
| Acenaphthylene | | 38000 J | 550 J | 750 JD | 570 J | 15000 UD | 550 J | 550 J | | 1300 |
| Anthracene | | 110000 | 2000 | 21000 JD | 2000 | 150000 UD | 1100 J | 1100 J | | 3500 |
| Benz(a)anthracene | | 210000 | 4800 | 5200 D | 5000 | 2600 J | 5000 | 5000 | | 9200 |
| Benz(b)fluoranthene | | 250000 | 6500 | 6800 D | 6700 | 5800 J | 7900 | 7900 | | 8900 |
| Benz(g)perylene | | 87000 | 3300 | 3400 D | 3500 | 1600 J | 2800 J | 2800 J | | 100 J |
| Benz(k)fluoranthene | | 110000 | 2400 | 3100 JD | 2200 | 2700 J | 2900 J | 2900 J | | 3700 |
| Benzalpyrene | | 200000 | 4800 | 5400 D | 4800 | 3000 J | 5600 | 5600 | | 1300 |

NOTES: U - Not detected, J - Estimated value, E - outside linear range, D - from diluted analysis, --- - Not analyzed.



Table 2
Honeywell
Tonawanda, NY
Semivolatile Organic Compound Data
Storm Sediment Samples

| Compound | Sample ID | Outfall #2 | Storm Inlet A | Storm Inlet A DL | | Storm Inlet A RE | Storm Inlet A | Storm Inlet B | Storm Inlet B |
|-----------------------------------|-----------|------------|---------------|------------------|----------|------------------|---------------|---------------|---------------|
| | | | | Sample Date | 05/11/99 | | | | |
| | | | | | ug/Kg | ug/Kg | ug/Kg | ug/Kg | |
| Bis(2-chloroethoxy)methane | | 42000 UD | 1500 UD | 3100 UD | 1500 UD | 15000 UD | 15000 UD | 2900 UD | 920 UD |
| Bis(2-chloroethyl)ether | | 42000 UD | 1500 UD | 3100 UD | 1500 UD | 15000 UD | 15000 UD | 2900 UD | 920 UD |
| Bis(2-chloroisopropyl) ether | | 42000 UD | 1500 UD | 3100 UD | 1500 UD | 15000 UD | 15000 UD | 2900 UD | 920 UD |
| Bis(2-ethylhexyl)phthalate (BEHP) | | 7400 J | 340 J | 320 JD | 330 J | 23000 | 720 J | 680 J | 920 UD |
| Butyl benzyl phthalate | | 42000 UD | 1500 UD | 3100 UD | 1500 UD | 15000 UD | 15000 UD | 2900 UD | 920 UD |
| Carbazole | | 18000 J | 400 J | 430 JD | 380 J | 15000 UD | 15000 UD | 2900 UD | 290 J |
| Chrysene | | 210000 | 5300 | 5600 D | 5200 | 5300 J | 6300 | 7700 | |
| Di-n-butyl phthalate | | 42000 UD | 1500 UD | 3100 UD | 1500 UD | 2600 J | 2900 UD | 2900 UD | 110 J |
| Di-n-octyl phthalate | | 42000 UD | 1500 UD | 3100 UD | 1500 UD | 15000 UD | 15000 UD | 2900 UD | 920 UD |
| Dibenzoz(a,h)anthracene | | 24000 J | 990 J | 930 JD | 1000 J | 15000 UD | 870 J | 800 J | |
| Dibenzofuran | | 27000 J | 410 J | 440 JD | 410 J | 15000 UD | 2900 UD | 2900 UD | 290 J |
| Diethyl phthalate | | 42000 UD | 1500 UD | 3100 UD | 1500 UD | 15000 UD | 15000 UD | 2900 UD | 550 J |
| Dimethyl phthalate | | 42000 UD | 1500 UD | 3100 UD | 1500 UD | 15000 UD | 2900 UD | 2900 UD | 920 UD |
| Fluoranthene | | 480000 | 7800 | 9900 D | 7200 | 7200 | 7900 | 19000 D | |
| Fluorene | | 60000 | 1200 J | 1300 JD | 1200 J | 15000 UD | 420 J | 1300 | |
| Hexachlorobenzene | | 42000 UD | 1500 UD | 3100 UD | 1500 UD | 15000 UD | 2900 UD | 2900 UD | 240 J |
| Hexachlorobutadiene | | 42000 UD | 1500 UD | 3100 UD | 1500 UD | 15000 UD | 2900 UD | 2900 UD | 920 UD |
| Hexachlorocyclopentadiene | | 42000 UD | 1500 UD | 3100 UD | 1500 UD | 15000 UD | 2900 UD | 2900 UD | 920 UD |
| Hexachloroethane | | 42000 UD | 1500 UD | 3100 UD | 1500 UD | 15000 UD | 2900 UD | 2900 UD | 220 J |
| Indeno[1,2,3-cd]pyrene | | 87000 | 3000 | 3100 JD | 3200 | 1600 J | 2700 J | 1000 | |
| Iophorone | | 42000 UD | 1500 UD | 3100 UD | 1500 UD | 15000 UD | 2900 UD | 2900 UD | 920 UD |
| N-Nitrosodiphenylamine | | 42000 UD | 1500 UD | 3100 UD | 1500 UD | 15000 UD | 2900 UD | 2900 UD | 920 UD |
| N-Nitrosodipipyramine | | 42000 UD | 1500 UD | 3100 UD | 1500 UD | 15000 UD | 2900 UD | 2900 UD | 920 UD |
| Naphthalene | | 64000 | 180 J | 3100 UD | 180 J | 15000 UD | 2900 UD | 2900 UD | 680 J |
| Nitrobenzene | | 42000 UD | 1500 UD | 3100 UD | 1500 UD | 15000 UD | 2900 UD | 2900 UD | 920 UD |
| Pentachlorophenol | | 210000 UD | 7800 UD | 16000 UD | 7800 UD | 74000 UD | 15000 UD | 4600 UD | |
| Phenanthrene | | 340000 | 7100 | 7800 D | 7100 | 2700 J | 2600 J | 9000 | |
| Phenol | | 42000 UD | 1500 UD | 31000 UD | 1500 UD | 15000 UD | 2900 UD | 2900 UD | 920 UD |
| Pyrene | | 500000 | 13000 E | 14000 D | 14000 E | 11000 J | 8400 | 6600 | |

NOTES: U - Not detected, J - Estimated value, E - outside linear range, D - from diluted analysis, --- - Not analyzed.



Table 3
Honeywell
Tonawanda, NY
Inorganic Data

| Compound | Sample ID | Outfall #1 | Storm Sediment Samples | | | Outfall #2 | Storm Inlet A | |
|-----------|-----------|------------|------------------------|----------------|----------|------------|---------------|---------|
| | | | Sample Date | Outfall #1 DUP | | | | |
| | | | | 11/15/99 | 05/10/99 | 11/15/00 | 05/11/99 | |
| Compound | Sample ID | Outfall #1 | Units | mg/Kg | mg/Kg | mg/Kg | mg/Kg | mg/Kg |
| Aluminum | 24700 | 14800 | | 18600 | 11000 | 18300 | 16300 | 17600 |
| Antimony | | 0.71 UD | | 0.29 UD | 0.21 UD | 0.92 UD | 0.26 J | 0.37 UD |
| Arsenic | 23.5 | 9.0 | | 12.3 | 53.5 | 11.1 | 7.3 | 14.5 |
| Barium | 118 | 69.3 | | 100 | 104 | 132 | 118 | 112 |
| Beryllium | 3.2 J | 1.4 J | | 1.7 J | 2.1 J | 3.3 | 1.0 J | 1.3 J |
| Cadmium | 1.9 J | 0.048 UD | | 0.052 UD | 2.0 J | 0.058 J | 0.59 J | 0.25 J |
| Calcium | 15700 | 10500 | | 15900 | 84600 | 63000 | 39700 | 46600 |
| Chromium | 105 | 38.5 | | 50.2 | 34.9 | 18.4 | 28.2 | 41.7 |
| Cobalt | 10 J | 6.9 J | | 9.3 J | 16.3 J | 6.4 | 13.5 | 10.1 J |
| Copper | 64.3 | 31.2 | | 55.8 | 45.1 | 24.0 | 33.6 | 45.9 |
| Cyanide | 17.4 | 0.95 UD | | 1.0 UD | 4.8 | 0.63 UD | 1.2 UD | 1.1 UD |
| Iron | 133000 | 46700 | | 52600 | 118000 | 21800 | 33800 | 30200 |
| Lead | 87.4 | 52.1 | | 73.0 | 41.3 | 31.1 | 34.8 | 131 |
| Magnesium | 6070 | 5010 | | 7870 | 18100 | 12800 | 10600 | 10400 |
| Manganese | 960 | 379 | | 506 | 1420 | 1970 | 523 | 562 |
| Mercury | 0.53 | 0.18 J | | 0.33 | 0.77 | 0.28 | 0.070 J | 0.063 J |
| Nickel | 32.1 | 22.7 | | 37.8 | 39.9 | 14.5 | 24.7 | 29.8 |
| Potassium | 1440 J | | | 2400 | 733 J | 1880 | 2630 | 3180 |
| Selenium | 6.5 | 1.9 | | 2.4 | 4.4 | 2.8 | 1.3 | 1.9 |
| Silver | 0.44 UD | 0.14 UD | | 0.15 UD | 0.57 UD | 0.15 J | 0.23 UD | 0.16 UD |
| Sodium | 405 J | 152 J | | 228 | 402 J | 632 | 399 | 315 |
| Thallium | 1.7 UD | 0.82 J | | 0.77 UD | 2.2 UD | 0.56 J | 0.89 UD | 0.82 UD |
| Vanadium | 43.2 | 26.7 | | 35.2 | 16.8 J | 19.2 | 34.6 | 31.6 |
| Zinc | 446 | 193 | | 255 | 140 | 110 | 125 | 201 |

NOTES: U - Not detected, J - Estimated value, E - outside linear range, D - from diluted analysis, --- Not analyzed.



Table 3
Honeywell
Tonawanda, NY
Inorganic Data
Storm Sediment Samples

| Compound | Sample ID | Storm Inlet B | |
|-----------|-----------|---------------|----------|
| | | Sample Date | 11/15/00 |
| | Units | mg/Kg | mg/Kg |
| Aluminum | 17300 | 30000 | |
| Antimony | 0.28 UD | 0.46 J | |
| Arsenic | 11.9 | 24.0 | |
| Barium | 104 | 125 | |
| Beryllium | 1.3 J | 2.6 J | |
| Cadmium | 0.70 J | 1.4 J | |
| Calcium | 12000 | 5680 | |
| Chromium | 44.1 | 69.9 | |
| Cobalt | 13.6 | 19.9 | |
| Copper | 54.1 | 103 | |
| Cyanide | 0.88 UD | 1.4 UD | |
| Iron | 35360 | 47660 | |
| Lead | 31.1 | 60.6 | |
| Magnesium | 7570 | 6580 | |
| Manganese | 514 | 509 | |
| Mercury | 0.079 J | 0.13 J | |
| Nickel | 32.4 | 55.7 | |
| Potassium | 1870 | 3980 | |
| Selenium | 1.3 | 3.1 | |
| Silver | 0.18 UD | 0.20 UD | |
| Sodium | 187 | 214 J | |
| Thallium | 0.67 UD | 1.5 J | |
| Vanadium | 32.7 | 53.7 | |
| Zinc | 211 | 605 | |

NOTES: U - Not detected, J - Estimated value, E - outside linear range, D - from diluted analysis, --- - Not analyzed



**O'BRIEN & GERE
Honeywell
Tonawanda, NY**

**Table 4
Volatile Organic Compound Data
Ground Water Samples**

| Compound | Sample ID | NYSDEC Class GA GW Standards ug/L | MW-01 | MW-01 DL | MW-01 | MW-01 DUP | MW-02 | |
|-----------------------------|-----------|--|---------|----------|---------|-----------|--------|-------|
| | | | Units | ug/L | ug/L | ug/L | ug/L | |
| 1,1,1-Trichloroethane | 5* | 0.4 J | 2 UD | 50 U Y | 50 U Y | 50 U Y | 0.5 UD | 0.5 U |
| ,,2,2-Tetrachloroethane | NS | 0.5 UD | 2 UD | 50 U | 50 U | 50 U | 0.5 UD | 0.5 U |
| 1,1,2-Trichloroethane | 1 | 0.5 UD | 2 UD Y | 50 U Y | 50 U Y | 50 U Y | 0.5 UD | 0.5 U |
| 1,1-Dichloroethane | 5* | 0.7 | 0.7 JD | 50 U Y | 50 U Y | 50 U Y | 0.5 UD | 0.5 U |
| 1,1-Dichloroethene | 5* | 0.5 UD | 2 UD | 50 U Y | 50 U Y | 50 U Y | 0.5 UD | 0.5 U |
| 1,2-Dichloroethane | 0.6 | 0.5 UD | 2 UD Y | 50 U Y | 50 U Y | 50 U Y | 0.5 UD | 0.5 U |
| 1,2-Dichloropropane | 1 | 0.5 UD | 2 UD Y | 50 U Y | 50 U Y | 50 U Y | 0.5 UD | 0.5 U |
| 2-Butanone (MEK) | NS | 10 UD | 50 UD | 1000 U | 1000 U | 1000 U | 10 UD | 10 U |
| 2-Hexanone | NS | 5 UD | 25 UD | 500 U | 500 U | 500 U | 5 UD | 5 U |
| 4-Methyl-2-pentanone (MBK) | NS | 5 UD | 25 UD | 500 U | 500 U | 500 U | 5 UD | 5 U |
| Acetone | NS | 13 | 14 JD | 1000 U | 1000 U | 1000 U | 1 J | 10 U |
| Benzene | 1 | 70 E Y | 67 D Y | 46 J Y | 79 Y | 79 Y | 0.5 UD | 0.5 U |
| Bromodichloromethane | NS | 0.5 UD | 2 UD | 50 U | 50 U | 50 U | 0.5 UD | 0.5 U |
| Bromoform | NS | 0.5 UD | 2 UD | 50 U | 50 U | 50 U | 0.5 UD | 0.5 U |
| Bromomethane | 5* | 1 UD | 5 UD | 100 U Y | 100 U Y | 100 U Y | 1 UD | 1 U |
| Carbon disulfide | NS | 11 | 11 D | 15 J | 50 U | 50 U | 0.5 UD | 0.1 J |
| Carbon tetrachloride | 5 | 0.5 UD | 2 UD | 50 U Y | 50 U Y | 50 U Y | 0.5 UD | 0.5 U |
| Chlorobenzene | 5* | 0.5 UD | 2 UD | 50 U Y | 50 U Y | 50 U Y | 0.5 UD | 0.5 U |
| Chloroethane | 5* | 1 UD | 5 UD | 100 U Y | 100 U Y | 100 U Y | 1 UD | 1 U |
| Chloroform | 7 | 0.2 J | 2 UD | 50 U Y | 50 U Y | 50 U Y | 0.5 UD | 0.5 U |
| Chloromethane | NS | 1 UD | 5 UD | 100 U | 100 U | 100 U | 1 UD | 1 U |
| Dibromochloromethane | 5* | 0.5 UD | 2 UD | 50 U Y | 50 U Y | 50 U Y | 0.5 UD | 0.5 U |
| Ethene, 1,2-dichloro-, (E)- | 5* | 0.5 UD | 2 UD | 50 U Y | 50 U Y | 50 U Y | 0.5 UD | 0.5 U |
| Ethylbenzene | 5* | 28 Y | 27 D Y | 14 J Y | 18 J Y | 18 J Y | 0.5 UD | 0.5 U |
| Methylene chloride | 5* | 2 UD | 10 UD Y | 200 U Y | 22 J Y | 22 J Y | 2 UD | 0.3 J |
| Styrene | 5* | 2 | 2 JD | 50 U Y | 50 U Y | 50 U Y | 0.5 UD | 0.5 U |
| Tetrachloroethene | 5* | 0.5 | 0.5 JD | 50 U Y | 50 U Y | 50 U Y | 0.5 UD | 0.5 U |
| Toluene | 5* | 29 Y | 26 D Y | 17 J Y | 50 U Y | 50 U Y | 0.5 UD | 0.5 U |
| Trichloroethene | 5* | 1 | 0.9 JD | 50 U Y | 50 U Y | 50 U Y | 0.5 UD | 0.5 U |
| Vinyl chloride | 2 | 1 UD | 5 UD Y | 100 U Y | 100 U Y | 100 U Y | 1 UD | 1 U |
| Xylene (total) | 5* | 200 E Y | 200 D Y | 86 Y | 45 J Y | 45 J Y | 0.5 UD | 0.5 U |
| cis-1,2-Dichloropropene | 5* | 0.2 J | 2 UD | 50 U Y | 50 U Y | 50 U Y | 0.5 UD | 0.5 U |
| cis-1,3-Dichloropropene | NS | 0.5 UD | 2 UD | 50 U | 50 U | 50 U | 0.5 UD | 0.5 U |
| trans-1,3-Dichloropropene | NS | 0.5 UD | 2 UD | 50 U | 50 U | 50 U | 0.5 UD | 0.5 U |

NOTES:

U - not detected, J - estimated value, E - outside linear range, D - from diluted analysis, Y - exceeds GW Standard, --- - not analyzed.

* - principal organic contaminant standard for ground water is 5 ug/L.



Table 4
Honeywell
Tonawanda, NY
Volatile Organic Compound Data
Ground Water Samples

| Compound | Sample ID | NYSDEC Class GA GW Standards | MW-03 05/12/99 | MW-04 05/12/99 | MW-04 06/09/01 | MW-05 05/12/99 | MW-05 06/07/01 |
|-----------------------------|-------------|------------------------------------|-------------------|-------------------|-------------------|-------------------|-------------------|
| | Sample Date | ug/L | ug/L | ug/L | ug/L | ug/L | ug/L |
| | Units | | | | | | |
| 1,1,1-Trichloroethane | 5* | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U |
| 1,1,2,2-Tetrachloroethane | NS | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U |
| 1,1,2-Trichloroethane | 1 | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U |
| 1,1-Dichloroethane | 5* | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U |
| 1,1-Dichloroethene | 5* | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U |
| 1,2-Dichloroethane | 0.6 | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U |
| 1,2-Dichloropropane | 1 | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U |
| 2-Butanone (MEK) | NS | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U |
| 2-Hexanone | NS | 5 UD | 5 U | 5 UD | 5 U | 5 UD | 5 U |
| 4-Methyl-2-pentanone (MBK) | NS | 5 UD | 5 U | 5 UD | 5 U | 5 UD | 5 U |
| Acetone | NS | 2 J | 6 J | 10 UD | 3 J | 2 J | 10 J |
| Benzene | 1 | 0.5 UD | 0.1 J | 1 | 0.5 U | 0.5 UD | 0.5 U |
| Bromodichloromethane | NS | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U |
| Bromoform | NS | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U |
| Bromomethane | 5* | 1 UD | 1 U | 1 UD | 1 U | 1 UD | 1 U |
| Carbon disulfide | NS | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U |
| Carbon tetrachloride | 5 | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U |
| Chlorobenzene | 5* | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U |
| Chloroethane | 5* | 1 UD | 1 U | 1 UD | 1 U | 1 UD | 1 U |
| Chloroform | 7 | 0.5 UD | 0.2 J | 0.1 J | 0.5 U | 0.5 UD | 0.5 U |
| Chloromethane | NS | 1 UD | 1 U | 1 UD | 1 U | 1 UD | 1 U |
| Dibromochloromethane | 5* | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U |
| Ethene, 1,2-dichloro-, (E)- | 5* | 0.5 UD | 1 | 0.5 UD | 0.5 U | 0.5 | 0.8 |
| Ethybenzene | 5* | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U |
| Methylene chloride | 5* | 2 UD | 2 U | 2 UD | 0.4 J | 2 UD | 2 U |
| Styrene | 5* | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U |
| Tetrachloroethane | 5* | 0.5 UD | 0.3 J | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U |
| Toluene | 5* | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U | 0.5 UD | 0.1 J |
| Trichloroethene | 5* | 0.5 UD | 6 Y | 0.8 | 0.5 U | 4 | 0.7 |
| Vinyl chloride | 2 | 1 UD | 0.7 J | 1 | 1 U | 0.4 J | 0.3 J |
| Xylene (total) | 5* | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U |
| cis-1,2-Dichloroethylene | 5* | 0.5 UD | 20 Y | 2 | 0.5 U | 1 | 2 |
| cis-1,3-Dichloropropene | NS | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U |
| trans-1,3-Dichloropropene | NS | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U | 0.5 UD | 0.5 U |

NOTES:

U - not detected, J - estimated value, E - outside linear range, D - from diluted analysis, Y - exceeds GW Standard, — - not analyzed.

*

- principal organic contaminant standard for ground water is 5 ug/L.



Table 4
Honeywell
Tonawanda, NY
Volatile Organic Compound Data
Ground Water Samples

| Compound | Sample ID | NYSDEC Class GA GW Standards | MW-06 05/12/99 | MW-06 DL. 06/06/01 | MW-06 05/12/99 | MW-07 06/06/01 | MW-07 05/12/99 | MW-08 05/12/99 |
|-----------------------------|-----------|------------------------------------|-------------------|-----------------------|-------------------|-------------------|-------------------|-------------------|
| | | ug/L | ug/L | ug/L | ug/L | ug/L | ug/L | ug/L |
| 1,1,1-Trichloroethane | 5* | 0.5 UD | 2 UD | 50 U Y | 0.5 UD | 5 U | 5 U | 1 |
| 1,1,2,2-Tetrachloroethane | NS | 0.5 UD | 2 UD | 50 U | 0.5 UD | 5 U | 5 U | 0.5 UD |
| 1,1,2-Trichloroethane | 1 | 0.5 UD | 2 UD Y | 50 U Y | 0.5 UD | 5 U Y | 5 U Y | 0.5 UD |
| 1,1-Dichloroethane | 5* | 0.3 J | 2 UD | 50 U Y | 0.4 J | 5 U | 5 U | 9 Y |
| 1,1-Dichloroethene | 5* | 0.5 UD | 2 UD | 50 U Y | 0.5 UD | 5 U | 5 U | 0.5 UD |
| 1,2-Dibromoethane | 0.6 | 0.5 UD | 2 UD Y | 50 U Y | 0.5 UD | 5 U Y | 5 U Y | 0.5 UD |
| 1,2-Dibromoethane | 1 | 0.5 UD | 2 UD Y | 50 U Y | 0.5 UD | 5 U Y | 5 U Y | 0.5 UD |
| 2-Butanone (MEK) | NS | 10 UD | 50 UD | 1000 U | 2 J | 100 U | 100 U | 10 UD |
| 2-Hexanone | NS | 5 UD | 25 UD | 500 U | 5 UD | 50 U | 50 U | 5 UD |
| 4-Methyl-2-pentanone (MIBK) | NS | 5 UD | 25 UD | 500 U | 5 UD | 50 U | 50 U | 5 UD |
| Acetone | NS | 10 UD | 50 UD | 1000 U | 27 | 14 J | 14 J | 10 UD |
| Benzene | 1 | 82 E Y | 81 D Y | 79 Y | 15 Y | 13 Y | 13 Y | 0.5 UD |
| Bromodichloromethane | NS | 0.5 UD | 2 UD | 50 U | 0.5 UD | 5 U | 5 U | 0.5 UD |
| Bromoform | NS | 0.5 UD | 2 UD | 50 U | 0.5 UD | 5 U | 5 U | 0.5 UD |
| Bromomethane | 5* | 1 UD | 5 UD | 100 U Y | 1 UD | 10 U Y | 10 U Y | 1 UD |
| Carbon disulfide | NS | 3 | 14 D | 12 J | 3 | 13 | 13 | 0.5 UD |
| Carbon tetrachloride | 5 | 0.5 UD | 2 UD | 50 U Y | 0.5 UD | 5 U | 5 U | 0.5 UD |
| Chlorobenzene | 5* | 0.5 UD | 2 UD | 50 U Y | 0.5 UD | 5 U | 5 U | 0.5 UD |
| Chloroethane | 5* | 1 UD | 5 UD | 100 U Y | 1 UD | 10 U Y | 10 U Y | 1 UD |
| Chloroform | 7 | 0.2 J | 2 UD | 50 U Y | 0.5 UD | 5 U | 5 U | 0.5 UD |
| Chloromethane | NS | 1 UD | 5 UD | 100 U | 1 UD | 10 U | 10 U | 1 UD |
| Dibromochloromethane | 5* | 0.5 UD | 2 UD | 50 U Y | 0.5 UD | 5 U | 5 U | 0.5 UD |
| Ethene, 1,2-dichloro-, (E)- | 5* | 0.5 UD | 2 UD | 50 U Y | 0.5 UD | 5 U | 5 U | 0.5 UD |
| Ethylbenzene | 5* | 20 Y | 19 D Y | 17 J Y | 1 | 2 J | 2 J | 0.5 UD |
| Methylene chloride | 5* | 2 UD | 10 UD Y | 14 J Y | 2 UD | 2 J | 2 J | 2 UD |
| Styrene | 5* | 0.6 | 0.6 JD | 50 U Y | 0.2 J | 5 U | 5 U | 0.5 UD |
| Tetrachloroethene | 5* | 0.1 J | 2 UD | 50 U Y | 0.5 UD | 5 U | 5 U | 0.5 UD |
| Toluene | 5* | 10 Y | 9 D Y | 50 U Y | 4 | 4 J | 4 J | 0.5 UD |
| Trichloroethene | 5* | 3 | 3 D | 50 U Y | 0.2 J | 5 U | 5 U | 0.2 J |
| Vinyl chloride | 2 | 1 UD | 5 UD Y | 100 U Y | 1 UD | 10 U Y | 10 U Y | 1 UD |
| Xylene (total) | 5* | 54 Y | 50 D Y | 49 J Y | 6 Y | 9 Y | 9 Y | 0.5 UD |
| cis-1,2-Dichloroethene | 5* | 0.7 | 0.6 JD | 50 U Y | 0.5 UD | 5 U | 5 U | 1 |
| cis-1,3-Dichloropropene | NS | 0.5 UD | 2 UD | 50 U | 0.5 UD | 5 U | 5 U | 0.5 UD |
| trans-1,3-Dichloropropene | NS | 0.5 UD | 2 UD | 50 U | 0.5 UD | 5 U | 5 U | 0.5 UD |

NOTES: U - not detected, J - estimated value, E - outside linear range, D - from diluted analysis, Y - exceeds GW Standard, --- - not analyzed.
* - principal organic contaminant standard for ground water is 5 ug/L.



Table 4
Honeywell
Tonawanda, NY
Volatile Organic Compound Data
Ground Water Samples

| Compound | Sample ID | NYSDEC Class GA GW Standards | MW-08 06/05/01 | MW-09 06/06/01 | MW-10 06/05/01 | MW-11 06/07/01 |
|-----------------------------|-----------|------------------------------------|-------------------|-------------------|-------------------|-------------------|
| | | ug/L | ug/L | ug/L | ug/L | ug/L |
| 1,1,1-Trichloroethane | 5* | 1 | 1 | 1 | 5 U | 0.5 U |
| 1,1,2,2-Tetrachloroethane | NS | 0.5 U | 1 U | 5 U | 5 U | 0.5 U |
| 1,1,2-Trichloroethane | 1 | 0.5 U | 1 U | 5 U | 5 U | 0.5 U |
| 1,1-Dichloroethane | 5* | 7 Y | 1 U | 5 U | 5 U | 0.5 U |
| 1,1-Dichloroethene | 5* | 0.5 U | 1 U | 5 U | 5 U | 0.5 U |
| 1,2-Dichloroethane | 0.6 | 0.5 U | 1 U Y | 5 U Y | 5 U Y | 0.5 U |
| 1,2-Dichloropropane | 1 | 0.5 U | 1 U | 5 U Y | 5 U Y | 0.5 U |
| 2-Butanone (MEK) | NS | 10 U | 20 U | 100 U | 100 U | 10 U |
| 2-Hexanone | NS | 5 U | 10 U | 50 U | 50 U | 5 U |
| 4-Methyl-1-pentanone (MIBK) | NS | 5 U | 10 U | 50 U | 50 U | 5 U |
| Acetone | NS | 4 J | 20 U | 100 U | 100 U | 6 J |
| Benzene | 1 | 0.5 U | 5 Y | 16 Y | 16 Y | 0.6 |
| Bromodichloromethane | NS | 0.5 U | 1 U | 5 U | 5 U | 0.5 U |
| Bromoform | NS | 0.5 U | 1 U | 5 U | 5 U | 0.5 U |
| Bromomethane | 5* | 1 U | 2 U | 10 U Y | 10 U Y | 1 U |
| Carbon disulfide | NS | 0.5 U | 1 U | 5 U | 5 U | 0.5 U |
| Carbon tetrachloride | 5 | 0.5 U | 1 U | 5 U | 5 U | 0.5 U |
| Chlorobenzene | 5* | 0.5 U | 1 U | 5 U | 5 U | 0.5 U |
| Chloroethane | 5* | 1 U | 2 U | 10 U Y | 10 U Y | 1 U |
| Chloroform | 7 | 0.5 U | 1 U | 5 U | 5 U | 0.5 U |
| Chloromethane | NS | 1 U | 2 U | 10 U | 10 U | 1 U |
| Dibromochloromethane | 5* | 0.5 U | 1 U | 5 U | 5 U | 0.5 U |
| Ethene, 1,2-dichloro-, (E)- | 5* | 0.5 U | 2 | 5 U | 5 U | 0.2 J |
| Ethylbenzene | 5* | 0.5 U | 7 Y | 1 J | 1 J | 0.5 U |
| Methylene chloride | 5* | 2 U | 0.4 J | 20 U Y | 20 U Y | 0.1 J |
| Styrene | 5* | 0.5 U | 1 U | 5 U | 5 U | 0.5 U |
| Tetrachloroethene | 5* | 0.5 U | 1 U | 5 U | 5 U | 0.5 U |
| Toluene | 5* | 0.5 U | 1 J | 3 J | 3 J | 0.1 J |
| Trichloroethene | 5* | 0.2 J | 2 | 5 U | 5 U | 2 |
| Vinyl chloride | 2 | 1 U | 5 Y | 10 U Y | 10 U Y | 0.3 J |
| Xylene (total) | 5* | 0.5 U | 5 | 5 | 5 | 0.1 J |
| cis-1,2-Dichloroethene | 5* | 0.8 | 8 Y | 5 U | 5 U | 1 |
| cis-1,3-Dichloropropene | NS | 0.5 U | 1 U | 5 U | 5 U | 0.5 U |
| trans-1,3-Dichloropropene | NS | 0.5 U | 1 U | 5 U | 5 U | 0.5 U |

NOTES:

U - not detected, J - estimated value, E - outside linear range, D - from diluted analysis, Y - exceeds GW Standard, --- - not analyzed.
 * - principal organic contaminant standard for ground water is 5 ug/L.

File Number: 1163-22900

Date Printed: 01/16/02 17:08:03
 DBF File: N:\1163\22900\TEMP\DBF
 FXP File: N:\1163\22900\8260MW.FXP

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Table 5
Honeywell
Tonawanda, NY
Semivolatile Organic Compound Data
Ground Water Samples

| Compound | Sample ID | NYSDEC Class GA GW Standards ug/L | MW-01 05/12/99 | MW-01 DL 05/12/99 | MW-01 06/05/01 | MW-01 DUP 06/05/01 | MW-01 DL 06/05/01 | MW-02 05/12/99 |
|-----------------------------|-----------|--|-------------------|----------------------|-------------------|-----------------------|----------------------|-------------------|
| | | ug/L | ug/L | ug/L | ug/L | ug/L | ug/L | ug/L |
| 1,2,4-Trichlorobenzene | 5* | 5 UD | 530 UD Y | 5 U | 5 U | 100 U Y | 100 U Y | 5 UD |
| 1,2-Dichlorobenzene | 3 | 5 UD Y | 530 UD Y | 5 U Y | 5 U Y | 100 U Y | 100 U Y | 5 UD Y |
| 1,3-Dichlorobenzene | 3 | 5 UD Y | 530 UD Y | 5 U Y | 5 U Y | 100 U Y | 100 U Y | 5 UD Y |
| 1,4-Dichlorobenzene | 3 | 5 UD Y | 530 UD Y | 5 U Y | 5 U Y | 100 U Y | 100 U Y | 5 UD Y |
| 2,4,5-Trichlorophenol | NS | 26 UD | 2600 UD | 26 U | 25 U | 510 U | 26 UD | 510 U |
| 2,4,6-Trichlorophenol | NS | 5 UD | 530 UD | 5 U | 5 U | 100 U | 5 UD | 5 UD |
| 2,4-Dichlorophenol | 1 | 5 UD Y | 530 UD Y | 5 U Y | 5 U Y | 100 U Y | 100 U Y | 5 UD Y |
| 2,4-Dimethylphenol | 1 | 17 Y | 530 UD Y | 8 Y | 9 Y | 100 U Y | 100 U Y | 5 UD Y |
| 2,4-Dinitrophenol | 1 | 26 UD Y | 2600 UD Y | 26 U Y | 25 U Y | 510 U | 26 UD Y | 510 U |
| 2,4-Dinitrotoluene | 5* | 5 UD | 530 UD Y | 5 U | 5 U | 100 U Y | 100 U Y | 5 UD |
| 2,6-Dinitrotoluene | 5* | 5 UD | 530 UD Y | 5 U | 5 U | 100 U Y | 100 U Y | 5 UD |
| 2-Chloronaphthalene | NS | 11 UD | 1100 UD | 10 U | 10 U | 200 U | 10 UD | 200 U |
| 2-Chlorophenol | NS | 5 UD | 530 UD | 5 U | 5 U | 100 U | 5 UD | 100 U |
| 2-Methylnaphthalene | NS | 380 E | 400 JD | 200 E | 200 E | 230 D | 10 UD | 230 D |
| 2-Methylphenol | 1 | 9 Y | 530 UD Y | 4 J Y | 5 J Y | 100 U Y | 5 UD | 100 U Y |
| 2-Nitroaniline | 5* | 26 UD Y | 2600 UD Y | 26 U Y | 25 U Y | 510 U Y | 26 UD Y | 510 U Y |
| 2-Nitrophenol | NS | 5 UD | 530 UD | 5 U | 5 U | 100 U | 5 UD | 100 U |
| 3,3-Dichlorobenzidine | 5* | 11 UD Y | 1100 UD Y | 10 U Y | 10 U Y | 200 U Y | 10 UD Y | 200 U Y |
| 3-Nitroaniline | 5* | 26 UD Y | 2600 UD Y | 26 U Y | 25 U Y | 510 U Y | 26 UD Y | 510 U Y |
| 4,6-Dinitro-2-methylphenol | NS | 53 UD | 5300 UD | 51 U | 51 U | 1000 U | 51 UD | 1000 U |
| 4-Bromophenyl phenyl ether | NS | 11 UD | 1100 UD | 10 U | 10 U | 200 U | 10 UD | 200 U |
| 4-Chloro-3-methylphenol | NS | 5 UD | 530 UD | 5 U | 5 U | 100 U | 5 UD | 100 U |
| 4-Chloraniline | 5* | 5 UD | 530 UD Y | 5 U | 5 U | 100 U Y | 5 UD | 100 U Y |
| 4-Chlorophenyl phenyl ether | NS | 11 UD | 1100 UD | 10 U | 10 U | 200 U | 10 UD | 200 U |
| 4-Methylphenol | 1 | 13 Y | 1100 UD Y | 7 J Y | 7 J Y | 200 U Y | 10 UD Y | 200 U Y |
| 4-Nitroaniline | 5* | 26 UD Y | 2600 UD Y | 26 U Y | 25 U Y | 510 U Y | 26 UD Y | 510 U Y |
| 4-Nitrophenol | NS | 26 UD | 2600 UD | 26 U | 25 U | 510 U | 26 UD | 510 U |
| Acenaphthene | NS | 19 | 1100 UD | 12 | 13 | 200 U | 10 UD | 200 U |
| Acenaphthylene | NS | 15 | 1100 UD | 10 J | 10 J | 200 U | 10 UD | 200 U |
| Anthracene | NS | 9 J | 1100 UD | 5 J | 5 J | 200 U | 10 UD | 200 U |
| Benz(a)anthracene | NS | 5 UD | 530 UD | 5 U | 5 U | 100 U | 5 UD | 100 U |
| Benz(b)fluoranthene | NS | 11 UD | 1100 UD | 10 U | 10 U | 200 U | 10 UD | 200 U |
| Benz(gh)perylene | NS | 11 UD | 1100 UD | 10 U | 10 U | 200 U | 10 UD | 200 U |
| Benz(k)fluoranthene | NS | 11 UD | 1100 UD | 10 U | 10 U | 200 U | 10 UD | 200 U |
| Benz[a]pyrene | NS | 11 UD | 1100 UD | 10 U | 10 U | 200 U | 10 UD | 200 U |

NOTES:

U - not detected, J - estimated value, E - outside linear range, D - from diluted analysis, Y - exceeds GW Standard, --- - not analyzed.

* - principal organic contaminant standard for ground water is 5 ug/L.



Table 5
Honeywell
Tonawanda, NY
Semivolatile Organic Compound Data
Ground Water Samples

| Compound | Sample ID | NYSDEC Class GA GW Standards | MW-01 05/12/99 | MW-01 DL. 05/12/99 | MW-01 06/05/01 | MW-01 DUP 06/05/01 | MW-01 DL 06/05/01 | MW-02 05/12/99 |
|-----------------------------------|-----------|------------------------------------|-------------------|-----------------------|-------------------|-----------------------|----------------------|-------------------|
| | | ug/L | ug/L | ug/L | ug/L | ug/L | ug/L | ug/L |
| Bis(2-chloroethoxy)methane | 5* | 5 UD | 530 UD Y | 5 U | 5 U | 100 U Y | 5 U | 5 UD |
| Bis(2-chloroethyl)ether | 1 | 5 UD Y | 530 UD Y | 5 U Y | 5 U Y | 100 U Y | 5 U Y | 5 UD Y |
| Bis(2-chloroisopropyl) ether | NS | 11 UD | 100 UD | 10 U | 10 U | 200 U | 10 U | 10 UD |
| Bis(2-ethylhexyl)phthalate (BEHP) | 5 | 5 UD | 530 UD Y | 5 U | 5 U | 100 U Y | 5 U | 5 UD |
| Butyl benzyl phthalate | NS | 11 UD | 1100 UD | 10 U | 10 U | 200 U | 10 U | 10 UD |
| Carbazole | NS | 92 | 1100 UD | 67 | 73 | 77 J | 77 J | 10 UD |
| Chrysene | NS | 5 UD | 530 UD | 5 U | 5 U | 100 U | 5 U | 5 UD |
| Di-n-butyl phthalate | 50 | 11 UD | 1100 UD Y | 10 U | 10 U | 200 U Y | 10 U | 10 UD |
| Di-n-octyl phthalate | NS | 11 UD | 1100 UD | 10 U | 10 U | 200 U | 10 U | 10 UD |
| Dibenz(a,h)anthracene | NS | 11 UD | 1100 UD | 10 U | 10 U | 200 U | 10 U | 10 UD |
| Dibenzofuran | NS | 44 | 530 UD | 25 | 27 | 29 J | 29 J | 5 UD |
| Diethyl phthalate | NS | 11 UD | 1100 UD | 10 U | 10 U | 200 U | 10 U | 10 UD |
| Dimethyl phthalate | NS | 11 UD | 1100 UD | 10 U | 10 U | 200 U | 10 U | 10 UD |
| Fluoranthene | NS | 5 J | 1100 UD | 3 J | 3 J | 200 U | 200 U | 10 UD |
| Fluorene | NS | 68 | 1100 UD | 4 J | 44 | 46 J | 46 J | 10 UD |
| Hexachlorobenzene | 0.04 | 5 UD Y | 530 UD Y | 5 U Y | 5 U Y | 100 U Y | 5 U | 5 UD Y |
| Hexachlorobutadiene | 0.5 | 5 UD Y | 530 UD Y | 5 U Y | 5 U Y | 100 U Y | 5 U | 5 UD Y |
| Hexachlorocyclopentadiene | 5* | 11 UD Y | 1100 UD Y | 10 U Y | 10 U Y | 200 U Y | 10 U | 10 UD Y |
| Hexachloroethane | 5* | 5 UD | 530 UD Y | 5 U | 5 U | 100 U Y | 5 U | 5 UD |
| Indeno(1,2,3-cd)pyrene | NS | 11 UD | 1100 UD | 10 U | 10 U | 200 U | 10 U | 10 UD |
| Isophorone | NS | 11 UD | 1100 UD | 10 U | 10 U | 200 U | 10 U | 10 UD |
| N-Nitrosodiphenylamine | NS | 11 UD | 1100 UD | 10 U | 10 U | 200 U | 10 U | 10 UD |
| N-Nitrosodipropylamine | NS | 11 UD | 1100 UD | 10 U | 10 U | 200 U | 10 U | 10 UD |
| Naphthalene | NS | 2500 E | 4600 D | 1300 E | 1300 E | 2500 D | 10 U | 10 UD |
| Nitrobenzene | 0.4 | 5 UD Y | 530 UD Y | 5 U Y | 5 U Y | 100 U Y | 5 U | 5 UD Y |
| Pentachlorophenol | 1 | 26 UD Y | 2600 UD Y | 26 U Y | 25 U Y | 510 U Y | 26 UD Y | 26 UD Y |
| Phenanthrene | NS | 42 | 1100 UD | 30 | 31 | 33 J | 33 J | 10 UD |
| Phenol | 1 | 4 J Y | 530 UD Y | 4 J Y | 4 J Y | 100 U Y | 5 U | 5 UD Y |
| Pyrene | NS | 4 J | 1100 UD | 3 J | 3 J | 200 U | 10 U | 10 UD |

NOTES:

U - not detected, J - estimated value, E - outside linear range, D - from diluted analysis, Y - exceeds GW Standard, — - not analyzed.

* - principal organic contaminant standard for ground water is 5 ug/L.



Table 5
Honeywell
Tonawanda, NY
Semivolatile Organic Compound Data
Ground Water Samples

| Compound | Sample ID | NYSDEC Class GA GW Standards | MW-02 RE 05/12/99 | MW-02 05/09/01 | MW-03 05/12/99 | MW-03 06/09/01 | MW-04 05/12/99 | MW-04 06/09/01 |
|-----------------------------|-----------|------------------------------------|----------------------|-------------------|-------------------|-------------------|-------------------|-------------------|
| | | ug/L | ug/L | ug/L | ug/L | ug/L | ug/L | ug/L |
| 1,2,4-Trichlorobenzene | 5* | 5 UD | 5 U | 5 UD | 5 U | 5 UD | 5 U | 5 U |
| 1,2-Dichlorobenzene | 3 | 5 UD Y | 5 U Y | 5 UD Y | 5 U Y | 5 UD Y | 5 U Y | 5 U Y |
| 1,3-Dichlorobenzene | 3 | 5 UD Y | 5 U Y | 5 UD Y | 5 U Y | 5 UD Y | 5 U Y | 5 U Y |
| 1,4-Dichlorobenzene | 3 | 5 UD Y | 5 U Y | 5 UD Y | 5 U Y | 5 UD Y | 5 U Y | 5 U Y |
| 2,4,5-Trichlorophenol | NS | 26 UD | 25 U | 25 UD | 26 U | 25 UD | 26 U | 26 U |
| 2,4,6-Trichlorophenol | NS | 5 UD | 5 U | 5 UD | 5 U | 5 UD | 5 U | 5 U |
| 2,4-Dichlorophenol | 1 | 5 UD Y | 5 U Y | 5 UD Y | 5 U Y | 5 UD Y | 5 U Y | 5 U Y |
| 2,4-Dimethylphenol | 1 | 5 UD Y | 5 U Y | 5 UD Y | 5 U Y | 5 UD Y | 5 U Y | 5 U Y |
| 2,4-Dinitrophenol | 1 | 26 UD Y | 25 U Y | 25 UD Y | 26 U Y | 25 UD Y | 26 U Y | 26 U Y |
| 2,4-Dinitrotoluene | 5* | 5 UD | 5 U | 5 UD | 5 U | 5 UD | 5 U | 5 U |
| 2,6-Dinitrotoluene | 5* | 5 UD | 5 U | 5 UD | 5 U | 5 UD | 5 U | 5 U |
| 2-Chloronaphthalene | NS | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U | 10 U |
| 2-Chlorophenol | NS | 5 UD | 5 U | 5 UD | 5 U | 5 UD | 5 U | 5 U |
| 2-Methylnaphthalene | NS | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U | 10 U |
| 2-Methylphenol | 1 | 5 UD Y | 5 U Y | 5 UD Y | 5 U Y | 5 UD Y | 5 U Y | 5 U Y |
| 2-Nitroaniline | 5* | 26 UD Y | 25 U Y | 25 UD Y | 26 U Y | 25 UD Y | 26 U Y | 26 U Y |
| 2-Nitrophenol | NS | 5 UD | 5 U | 5 UD | 5 U | 5 UD | 5 U | 5 U |
| 3,3-Dichlorobenzidine | 5* | 10 UD Y | 10 U Y | 10 UD Y | 10 U Y | 10 UD Y | 10 U Y | 10 U Y |
| 3-Nitroaniline | 5* | 26 UD Y | 25 U Y | 25 UD Y | 26 U Y | 25 UD Y | 26 U Y | 26 U Y |
| 4,6-Dinitro-2-methylphenol | NS | 51 UD | 51 U | 50 UD | 51 U | 50 UD | 52 U | 52 U |
| 4-Bromophenyl phenyl ether | NS | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U | 10 U |
| 4-Chloro-3-methylphenol | NS | 5 UD | 5 U | 5 UD | 5 U | 5 UD | 5 U | 5 U |
| 4-Chloraniline | 5* | 5 UD | 5 U | 5 UD | 5 U | 5 UD | 5 U | 5 U |
| 4-Chlorophenyl phenyl ether | NS | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U | 10 U |
| 4-Methylphenol | 1 | 10 UD Y | 10 U Y | 10 UD Y | 10 U Y | 10 UD Y | 10 U Y | 10 U Y |
| 4-Nitroaniline | 5* | 26 UD Y | 25 U Y | 25 UD Y | 26 U Y | 25 UD Y | 26 U Y | 26 U Y |
| 4-Nitrophenol | NS | 26 UD | 25 U | 25 UD | 26 U | 25 UD | 26 U | 26 U |
| Acenaphthene | NS | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U | 10 U |
| Acenaphthylene | NS | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U | 10 U |
| Anthracene | NS | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U | 10 U |
| Benzoaanthracene | NS | 5 UD | 5 U | 5 UD | 5 U | 5 UD | 5 U | 5 U |
| Benzofluoranthene | NS | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U | 10 U |
| Benzohperylene | NS | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U | 10 U |
| Benzokfluoranthene | NS | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U | 10 U |
| Benzolapryene | NS | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U | 10 U |

NOTES: U - not detected, J - estimated value, E - outside linear range, D - from diluted analysis, Y - exceeds GW Standard, --- - not analyzed.

* - principal organic contaminant standard for ground water is 5 ug/L.



Table 5
Honeywell
Tonawanda, NY
Semivolatile Organic Compound Data
Ground Water Samples

| Compound | Sample ID | NYSDEC Class GA GW Standards ug/L | MW-02 RE 05/12/99 ug/L | MW-02 06/09/01 ug/L | MW-03 05/12/99 ug/L | MW-03 06/09/01 ug/L | MW-04 05/12/99 ug/L | MW-04 06/09/01 ug/L |
|-----------------------------------|-----------|--|------------------------------|---------------------------|---------------------------|---------------------------|---------------------------|---------------------------|
| Bis(2-chloroethoxy)methane | 5* | 5 UD | 5 U | 5 UD | 5 U | 5 UD | 5 UD | 5 U |
| Bis(2-chloroethyl)ether | 1 | 5 UD Y | 5 U Y | 5 UD Y | 5 U Y | 5 UD Y | 5 UD Y | 5 U Y |
| Bis(2-chloroisopropyl) ether | NS | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 UD | 10 U |
| Bis(2-ethylhexyl)phthalate (BEHP) | 5 | 2 J | 2 J | 3 J | 14 Y | 3 J | 10 Y | 10 Y |
| Butyl benzyl phthalate | NS | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 UD | 10 U |
| Carbamate | NS | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 UD | 10 U |
| Chrysene | NS | 5 UD | 5 U | 5 UD | 5 U | 5 UD | 5 UD | 5 U |
| Di-n-butyl phthalate | 50 | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 UD | 10 U |
| Di-n-octyl phthalate | NS | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 UD | 10 U |
| Dibenz(a,h)anthracene | NS | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 UD | 10 U |
| Dibenzofuran | NS | 5 UD | 5 U | 5 UD | 5 U | 5 UD | 5 UD | 5 U |
| Diethyl phthalate | NS | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 UD | 10 U |
| Dimethyl phthalate | NS | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 UD | 10 U |
| Fluoranthene | NS | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 UD | 10 U |
| Fluorene | NS | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 UD | 10 U |
| Hexachlorobenzene | 0.04 | 5 UD Y | 5 U Y | 5 UD Y | 5 U Y | 5 UD Y | 5 UD Y | 5 U Y |
| Hexachlorobutadiene | 0.5 | 5 UD Y | 5 U Y | 5 UD Y | 5 U Y | 5 UD Y | 5 UD Y | 5 U Y |
| Hexachlorocyclopentadiene | 5* | 10 UD Y | 10 U Y | 10 UD Y | 10 U Y | 10 UD Y | 10 UD Y | 10 U Y |
| Hexachloroethane | 5* | 5 UD | 5 U | 5 UD | 5 U | 5 UD | 5 UD | 5 U |
| Indeno(1,2,3-cd)pyrene | NS | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 UD | 10 U |
| Isophorone | NS | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 UD | 10 U |
| N-Nitrosodiphenylamine | NS | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 UD | 10 U |
| N-Nitrosodiphenylamine | NS | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 UD | 10 U |
| Naphthalene | NS | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 UD | 10 U |
| Nitrobenzene | 0.4 | 5 UD Y | 5 U Y | 5 UD Y | 5 U Y | 5 UD Y | 5 UD Y | 5 U Y |
| Pentachlorophenol | 1 | 26 UD Y | 25 UD Y | 25 UD Y | 26 U Y | 25 UD Y | 25 UD Y | 26 U Y |
| Phenanthrene | NS | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 UD | 10 U |
| Phenol | 1 | 5 UD Y | 5 U Y | 5 UD Y | 5 U Y | 5 UD Y | 5 UD Y | 5 U Y |
| Pyrene | NS | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 UD | 10 U |

NOTES: U - not detected, J - estimated value, E - outside linear range, D - from diluted analysis, Y - exceeds GW Standard, --- - not analyzed.
* - principal organic contaminant standard for ground water is 5 ug/L.



Table 5
Honeywell
Tonawanda, NY
Semivolatile Organic Compound Data
Ground Water Samples

| Compound | Sample ID NYSDEC Class GA GW Standards | MW-05 05/12/99 | MW-05 RE | | MW-06 05/12/99 | MW-06 DL | | MW-06 RE 05/12/99 | |
|-----------------------------|---|-------------------|----------|--------|-------------------|----------|----------|----------------------|--|
| | | | Units | ug/L | MW-05 | | ug/L | | |
| | | | | | ug/L | ug/L | | | |
| 1,2,4-Trichlorobenzene | 5* | 6 UD Y | 6 UD Y | 6 UD Y | 6 U Y | 5 UD | 100 UD Y | 5 UD | |
| 1,2-Dichlorobenzene | 3 | 6 UD Y | 6 UD Y | 6 UD Y | 6 U Y | 5 UD Y | 100 UD Y | 5 UD Y | |
| 1,3-Dichlorobenzene | 3 | 6 UD Y | 6 UD Y | 6 UD Y | 6 U Y | 5 UD Y | 100 UD Y | 5 UD Y | |
| 1,4-Dichlorobenzene | 3 | 6 UD Y | 6 UD Y | 6 UD Y | 6 U Y | 5 UD Y | 100 UD Y | 5 UD Y | |
| 2,4,5-Trichlorophenol | NS | 29 UD | 29 UD | 32 U | 25 UD | 500 UD | 25 UD | 25 UD | |
| 2,4,6-Trichlorophenol | NS | 6 UD | 6 UD | 6 U | 5 UD | 100 UD | 5 UD | 5 UD | |
| 2,4-Dichlorophenol | 1 | 6 UD Y | 6 UD Y | 6 UD Y | 6 U Y | 5 UD Y | 100 UD Y | 5 UD Y | |
| 2,4-Dimethylphenol | 1 | 6 UD Y | 6 UD Y | 6 UD Y | 6 U Y | 2 J Y | 100 UD Y | --- | |
| 2,4-Dinitrophenol | 1 | 29 UD Y | 29 UD Y | 32 U Y | 25 UD Y | 500 UD Y | 25 UD Y | 25 UD Y | |
| 2,4-Dinitrotoluene | 5* | 6 UD Y | 6 UD Y | 6 UD Y | 6 U Y | 5 UD | 100 UD Y | 5 UD | |
| 2,6-Dinitrotoluene | 5* | 6 UD Y | 6 UD Y | 6 UD Y | 6 U Y | 5 UD | 100 UD Y | 5 UD | |
| 2-Chloronaphthalene | NS | 12 UD | 12 UD | 13 U | 10 UD | 200 UD | 10 UD | 10 UD | |
| 2-Chlorophenol | NS | 6 UD | 6 UD | 6 U | 5 UD | 100 UD | 5 UD | 5 UD | |
| 2-Methylnaphthalene | NS | 12 UD | 12 UD | 13 U | 230 E | 260 D | 230 E | 230 E | |
| 2-Methylphenol | 1 | 6 UD Y | 6 UD Y | 6 UD Y | 6 U Y | 5 UD Y | 100 UD Y | 5 UD Y | |
| 2-Nitroaniline | 5* | 29 UD Y | 29 UD Y | 32 U Y | 25 UD Y | 500 UD Y | 25 UD Y | 25 UD Y | |
| 2-Nitrophenol | NS | 6 UD | 6 UD | 6 U | 5 UD | 100 UD | 5 UD | 5 UD | |
| 3,3-Dichlorobenzidine | 5* | 12 UD Y | 12 UD Y | 13 U Y | 10 UD Y | 200 UD Y | 10 UD Y | 10 UD Y | |
| 3-Nitroaniline | 5* | 29 UD Y | 29 UD Y | 32 U Y | 25 UD Y | 500 UD Y | 25 UD Y | 25 UD Y | |
| 4,6-Dinitro-2-methylphenol | NS | 59 UD | 59 UD | 64 U | 50 UD | 1000 UD | 50 UD | 50 UD | |
| 4-Bromophenyl phenyl ether | NS | 12 UD | 12 UD | 13 U | 10 UD | 200 UD | 10 UD | 10 UD | |
| 4-Chloro-3-methylphenol | NS | 6 UD | 6 UD | 6 U | 5 UD | 100 UD | 5 UD | 5 UD | |
| 4-Chloraniline | 5* | 6 UD Y | 6 UD Y | 6 U Y | 5 UD | 100 UD Y | 5 UD | 5 UD | |
| 4-Chlorophenyl phenyl ether | NS | 12 UD | 12 UD | 13 U | 10 UD | 200 UD | 10 UD | 10 UD | |
| 4-Methylphenol | 1 | 12 UD Y | 12 UD Y | 13 U Y | 10 UD Y | 200 UD Y | 10 UD Y | 10 UD Y | |
| 4-Nitroaniline | 5* | 29 UD Y | 29 UD Y | 32 U Y | 25 UD Y | 500 UD Y | 25 UD Y | 25 UD Y | |
| 4-Nitrophenol | NS | 29 UD | 29 UD | 32 U | 25 UD | 500 UD | 25 UD | 25 UD | |
| Acenaphthene | NS | 12 UD | 12 UD | 13 U | 14 | 200 UD | 15 | 15 | |
| Acenaphthylene | NS | 12 UD | 12 UD | 13 U | 12 | 200 UD | 12 | 12 | |
| Anthracene | NS | 12 UD | 12 UD | 13 U | 9 J | 200 UD | 9 J | 9 J | |
| Benz(a)anthracene | NS | 6 UD | 6 UD | 6 U | 6 | 100 UD | 6 | 6 | |
| Benz(b)fluoranthene | NS | 12 UD | 12 UD | 13 U | 5 J | 200 UD | 6 J | 6 J | |
| Benz(gi)perylene | NS | 12 UD | 12 UD | 13 U | 10 UD | 200 UD | 10 UD | 10 UD | |
| Benz(k)fluoranthene | NS | 12 UD | 12 UD | 13 U | 2 J | 200 UD | 10 UD | 10 UD | |
| Benz(a)pyrene | NS | 12 UD | 12 UD | 13 U | 5 J | 200 UD | 5 J | 5 J | |

NOTES:

U-not detected.

J-estimated value.

E-outside linear range.

D-from diluted analysis.

Y-exceeds GW Standard.

---not analyzed.

*-principal organic contaminant standard for ground water is 5 ug/L.



Table 5
Honeywell
Tonawanda, NY
Semivolatile Organic Compound Data
Ground Water Samples

| Compound | Sample ID | NYSDEC Class GA GW Standards | MW-05 05/12/99 | MW-05 RE 05/12/99 | MW-05 06/07/01 | MW-06 05/12/99 | MW-06 DL 05/12/99 | MW-06 RE 05/12/99 |
|-----------------------------------|-----------|------------------------------------|-------------------|----------------------|-------------------|-------------------|----------------------|----------------------|
| | | ug/L | ug/L | ug/L | ug/L | ug/L | ug/L | ug/L |
| Bis(2-chloroethoxy)methane | 5* | 6 UD Y | 6 UD Y | 6 UD Y | 6 UD Y | 5 UD | 100 UD Y | 5 UD |
| Bis(2-chloroethyl)ether | 1 | 6 UD Y | 6 UD Y | 6 UD Y | 6 UD Y | 5 UD Y | 100 UD Y | 5 UD Y |
| Bis(2-chloroisopropyl) ether | NS | 12 UD | 12 UD | 13 U | 10 UD | 200 UD | 100 UD Y | 10 UD |
| Bis(2-ethylhexyl)phthalate (BEHP) | 5 | 6 UD Y | 6 UD Y | 4 J | 5 UD | 100 UD Y | 5 UD | 5 UD |
| Butylbenzyl phthalate | NS | 12 UD | 12 UD | 13 U | 10 UD | 200 UD | 100 UD Y | 10 UD |
| Carbazole | NS | 12 UD | 12 UD | 13 U | 15 | 200 UD | 100 UD Y | 15 |
| Chrysene | NS | 6 UD | 6 UD | 6 U | 8 | 100 UD | 100 UD Y | 8 |
| Di-n-butyl phthalate | 50 | 12 UD | 12 UD | 12 UD | 10 UD | 200 UD Y | 200 UD Y | 10 UD |
| Di-n-octyl phthalate | NS | 12 UD | 12 UD | 13 U | 10 UD | 200 UD | 100 UD Y | 10 UD |
| Dibenz(a,h)anthracene | NS | 12 UD | 12 UD | 12 UD | 10 UD | 200 UD | 200 UD | 10 UD |
| Dibenzofuran | NS | 6 UD | 6 UD | 6 U | 9 | 100 UD | 100 UD Y | 10 |
| Diethyl phthalate | NS | 12 UD | 12 UD | 13 U | 10 UD | 200 UD | 200 UD | 10 UD |
| Dimethyl phthalate | NS | 12 UD | 12 UD | 13 U | 10 UD | 200 UD | 200 UD | 10 UD |
| Fluoranthene | NS | 12 UD | 12 UD | 13 U | 12 | 200 UD | 200 UD | 11 |
| Fluorene | NS | 12 UD | 12 UD | 13 U | 43 | 47 JD | 47 JD | 43 |
| Hexachlorobenzene | 0.04 | 6 UD Y | 6 UD Y | 6 UD Y | 5 UD | 100 UD Y | 100 UD Y | 5 UD Y |
| Hexachlorobutadiene | 0.5 | 6 UD Y | 6 UD Y | 6 UD Y | 5 UD Y | 100 UD Y | 100 UD Y | 5 UD Y |
| Hexachlorocyclopentadiene | 5* | 12 UD Y | 12 UD Y | 13 U Y | 10 UD Y | 200 UD Y | 200 UD Y | 10 UD Y |
| Hexachloroethane | 5* | 6 UD Y | 6 UD Y | 6 UD Y | 5 UD | 100 UD Y | 100 UD Y | 5 UD |
| Indeno(1,2,3-cd)pyrene | NS | 12 UD | 12 UD | 13 U | 10 UD | 200 UD | 200 UD | 10 UD |
| Isophorone | NS | 12 UD | 12 UD | 13 U | 10 UD | 200 UD | 200 UD | 10 UD |
| N-Nitrosodiphenylamine | NS | 12 UD | 12 UD | 13 U | 10 UD | 200 UD | 200 UD | 10 UD |
| N-Nitrosodipropylamine | NS | 12 UD | 12 UD | 13 U | 10 UD | 200 UD | 200 UD | 10 UD |
| Naphthalene | 0.4 | 6 UD Y | 6 UD Y | 6 UD Y | 5 UD Y | 100 UD Y | 100 UD Y | 5 UD Y |
| Nitrobenzene | 0.4 | 29 UD Y | 29 UD Y | 32 UD Y | 25 UD Y | 500 UD Y | 500 UD Y | 25 UD Y |
| Pentachlorophenol | 1 | 6 UD Y | 6 UD Y | 6 UD Y | 5 UD Y | 100 UD Y | 100 UD Y | 50 |
| Phenanthrene | NS | 12 UD | 12 UD | 13 U | 50 | 55 JD | 55 JD | 50 |
| Phenol | 1 | 6 UD Y | 6 UD Y | 6 UD Y | 5 UD Y | 200 UD | 100 UD Y | 5 UD Y |
| Pyrene | NS | 12 UD | 12 UD | 13 U | 15 | 200 UD | 200 UD | 18 |

NOTES: U - not detected, J - estimated value, E - outside linear range, D - from diluted analysis, Y - exceeds GW Standard, — - not analyzed
* - principal organic contaminant standard for ground water is 5 ug/L.



Table 5
Honeywell
Tonawanda, NY
Semivolatile Organic Compound Data
Ground Water Samples

| Compound | Sample ID | NYSDEC Class GA GW Standards | MW-06 06/06/01 | MW-06 DL 06/06/01 | MW-07 05/12/99 | MW-07 DL 05/12/99 | MW-07 RE 05/12/99 | MW-07 06/06/01 |
|-----------------------------|-----------|------------------------------------|-------------------|----------------------|-------------------|----------------------|----------------------|-------------------|
| | | ug/L | ug/L | ug/L | ug/L | ug/L | ug/L | ug/L |
| 1,2,4-Trichlorobenzene | 5* | 5 U | 51 U Y | 5 U D | 20 U D Y | 5 U D | 5 U | 5 U |
| 1,2-Dichlorobenzene | 3 | 5 U Y | 51 U Y | 5 U D Y | 20 U D Y | 5 U D Y | 5 U Y | 5 U Y |
| 1,3-Dichlorobenzene | 3 | 5 U Y | 51 U Y | 5 U D Y | 20 U D Y | 5 U D Y | 5 U Y | 5 U Y |
| 1,4-Dichlorobenzene | 3 | 5 U Y | 51 U Y | 5 U D Y | 20 U D Y | 5 U D Y | 5 U Y | 5 U Y |
| 2,4,5-Trichlorophenol | NS | 25 U | 250 U | 26 U D | 100 U D | 26 U D | 26 U | 26 U |
| 2,4,6-Trichlorophenol | NS | 5 U | 51 U | 5 U D | 20 U D | 5 U D | 5 U | 5 U |
| 2,4-Dichlorophenol | 1 | 5 U Y | 51 U Y | 5 U D Y | 20 U D Y | 5 U D Y | 5 U Y | 5 U Y |
| 2,4-Dimethylphenol | 1 | 5 U Y | 51 U Y | 3 J Y | 20 U D Y | 3 J Y | 1 J | 1 J |
| 2,4-Dinitrophenol | 1 | 25 U Y | 250 U Y | 26 U D Y | 100 U D Y | 26 U D Y | 26 U Y | 26 U Y |
| 2,4-Dinitrotoluene | 5* | 5 U | 51 U Y | 5 U D | 20 U D Y | 5 U D | 5 U | 5 U |
| 2,6-Dinitrotoluene | NS | 10 U | 100 U | 10 U D | 40 U D | 10 U D | 10 U | 10 U |
| 2-Chloronaphthalene | NS | 5 U | 51 U | 5 U D | 20 U D | 5 U D | 5 U | 5 U |
| 2-Chlorophenol | NS | 180 E | 210 D | 61 | 64 D | 61 | 26 | 26 |
| 2-Methylnaphthalene | NS | 5 U Y | 51 U Y | 2 J Y | 20 U D Y | 2 J Y | 5 U Y | 5 U Y |
| 2-Methylphenol | 1 | 5 U Y | 25 U Y | 250 U Y | 26 U D Y | 100 U D Y | 26 U D Y | 26 U Y |
| 2-Nitroaniline | 5* | 5 U | 51 U | 5 U D | 20 U D | 5 U D | 5 U | 5 U |
| 2-Nitrophenol | NS | 10 U Y | 100 U Y | 10 U D Y | 40 U D Y | 10 U D Y | 10 U Y | 10 U Y |
| 3,3-Dichlorobenzidine | 5* | 25 U Y | 250 U Y | 26 U D Y | 100 U D Y | 26 U D Y | 26 U Y | 26 U Y |
| 3-Nitroaniline | 5* | 51 U | 510 U | 51 U D | 200 U D | 51 U D | 51 U | 51 U |
| 4,6-Dinitro-2-methylphenol | NS | 10 U | 100 U | 10 U D | 40 U D | 10 U D | 10 U | 10 U |
| 4-Bromophenyl phenyl ether | NS | 5 U | 51 U | 5 U D | 20 U D | 5 U D | 5 U | 5 U |
| 4-Chloro-2-methylphenol | 5* | 5 U | 51 U Y | 5 U D | 20 U D Y | 5 U D | 5 U | 5 U |
| 4-Chloroaniline | NS | 10 U | 100 U | 10 U D | 40 U D | 10 U D | 10 U | 10 U |
| 4-Chlorophenyl phenyl ether | 1 | 10 U Y | 100 U Y | 3 J Y | 40 U D Y | 3 J Y | 2 J Y | 2 J Y |
| 4-Nitroaniline | 5* | 25 U | 250 U Y | 26 U D Y | 100 U D Y | 26 U D Y | 26 U Y | 26 U Y |
| 4-Nitrophenol | NS | 25 U | 250 U | 26 U D | 100 U D | 26 U D | 26 U | 26 U |
| Acenaphthene | NS | 11 | 111 J | 3 J | 40 U D | 3 J | 1 J | 1 J |
| Acenaphthylene | NS | 3 J | 100 U | 3 J | 40 U D | 3 J | 10 U | 10 U |
| Anthracene | NS | 5 J | 100 U | 5 J | 40 U D | 5 J | 10 U | 10 U |
| Benzof[a]anthracene | NS | 5 U | 51 U | 7 | 20 U D | 7 | 5 U | 5 U |
| Benzof[b]fluoranthene | NS | 10 U | 100 U | 11 | 11 JD | 11 | 10 U | 10 U |
| Benzof[b]phenylene | NS | 10 U | 100 U | 3 J | 40 U D | 4 J | 10 U | 10 U |
| Benzof[k]fluoranthene | NS | 10 U | 100 U | 4 J | 40 U D | 5 J | 10 U | 10 U |
| Benzalpyrene | NS | 10 U | 100 U | 7 J | 40 U D | 8 J | 10 U | 10 U |

NOTES: U - not detected, J - estimated value, E - outside linear range, D - exceeds GW Standard, — - not analyzed.

* - principal organic contaminant standard for ground water is 5 ug/L.



Table 5
Honeywell
Tonawanda, NY
Semivolatile Organic Compound Data
Ground Water Samples

| Compound | Sample ID | NYSDEC Class GA GW Standards | MW-06 06/06/01 | MW-06 DL. 06/06/01 | MW-07 05/12/99 | MW-07 DL. 05/12/99 | MW-07 RE 05/12/99 | MW-07 06/06/01 |
|-----------------------------------|-----------|------------------------------------|-------------------|-----------------------|-------------------|-----------------------|----------------------|-------------------|
| | | ug/L | ug/L | ug/L | ug/L | ug/L | ug/L | ug/L |
| Bis(2-chloroethyl)ether | 5* | 5 U | 51 U Y | 5 U D | 20 U D Y | 5 U D | 5 U | 5 U |
| Bis(2-chloroisopropyl) ether | 1 NS | 5 U | 51 U Y | 5 U D Y | 20 U D Y | 5 U D Y | 5 U Y | 10 U |
| Bis(2-ethylhexyl)phthalate (BEHP) | 5 NS | 10 U | 100 U | 10 U D | 40 U D | 10 U D | 10 U | 10 U |
| Butyl benzyl phthalate | NS | 5 U | 51 U Y | 2 J | 20 U D Y | 3 J | 5 U | 5 U |
| Carbazole | NS | 10 U | 100 U | 10 U D | 40 U D | 10 U D | 10 U | 10 U |
| Chrysene | NS | 15 | 16 J | 3 J | 40 U D | 3 J | 1 J | 1 J |
| Di-n-butyl phthalate | 50 NS | 10 U | 100 U Y | 10 U D | 40 U D | 10 U D | 10 U | 10 U |
| Di-n-octyl phthalate | NS | 10 U | 100 U | 10 U D | 40 U D | 10 U D | 10 U | 10 U |
| Dibenzof[a,h]anthracene | NS | 10 U | 100 U | 10 U D | 40 U D | 10 U D | 10 U | 10 U |
| Dibenzofuran | NS | 5 J | 51 U | 3 J | 20 U D | 3 J | 5 U | 5 U |
| Diethyl phthalate | NS | 10 U | 100 U | 10 U D | 40 U D | 10 U D | 10 U | 10 U |
| Dimethyl phthalate | NS | 10 U | 100 U | 10 U D | 40 U D | 10 U D | 10 U | 10 U |
| Fluoranthene | NS | 3 J | 100 U | 1 J | 12 J D | 1 J | 10 U | 10 U |
| Fluorene | NS | 34 | 35 J | 1 J | 11 J D | 1 J | 3 J | 3 J |
| Hexachlorobenzene | 0.04 | 5 U Y | 51 U Y | 5 U D Y | 20 U D Y | 5 U D Y | 5 U Y | 5 U Y |
| Hexachlorobutadiene | 0.5 | 5 U Y | 51 U Y | 5 U D Y | 20 U D Y | 5 U D Y | 5 U Y | 5 U Y |
| Hexachlorocyclopentadiene | 5* | 10 U Y | 100 U Y | 10 U D Y | 40 U D Y | 10 U D Y | 10 U Y | 10 U Y |
| Hexachloroethane | 5* | 5 U | 51 U Y | 5 U D | 20 U D Y | 5 U D | 5 U | 5 U |
| Indeno[1,2,3-ed]pyrene | NS | 10 U | 100 U | 3 J | 40 U D | 4 J | 10 U | 10 U |
| Isophorone | NS | 10 U | 100 U | 10 U D | 40 U D | 10 U D | 10 U | 10 U |
| N-Nitrodiphenylamine | NS | 10 U | 100 U | 10 U D | 40 U D | 10 U D | 10 U | 10 U |
| N-Nitrosodipropylamine | NS | 10 U | 100 U | 10 U D | 40 U D | 10 U D | 10 U | 10 U |
| Naphthalene | NS | 720 E | 1100 D | 280 E | 330 D | 280 E | 220 E | 220 E |
| Nitrobenzene | 0.4 | 5 U Y | 51 U Y | 5 U D Y | 20 U D Y | 5 U D Y | 5 U Y | 5 U Y |
| Pentachlorophenol | 1 | 25 U Y | 250 U Y | 26 U D Y | 100 U D Y | 26 U D Y | 26 U Y | 26 U Y |
| Phenanthrene | NS | 35 | 37 J | 18 | 19 J D | 18 | 2 J | 2 J |
| Phenol | 1 | 1 J | 51 U Y | 2 J Y | 20 U D Y | 2 J Y | 2 J Y | 2 J Y |
| Pyrene | NS | 3 J | 100 U | 14 | 12 J D | 15 | 10 U | 10 U |

NOTES: U - not detected, J - estimated value, E - outside linear range, D - from diluted analysis, Y - exceeds GW Standard, --- - not analyzed.
 * - principal organic contaminant standard for ground water is 5 ug/L.



Table 5
Honeywell
Tonawanda, NY
Semivolatile Organic Compound Data

| Compound | Sample ID | NYSDEC Class GA | MW-07 DL | MW-08 | | | MW-09 06/05/01 | MW-10 06/05/01 | MW-11 06/07/01 | | | |
|-----------------------------|-----------|--------------------|----------|-------------|--------------|--------|-------------------|-------------------|-------------------|--|--|--|
| | | | | Sample Date | GW Standards | ug/L | | | | | | |
| | | | | | | | | | | | | |
| 1,2,4-Trichlorobenzene | 5* | | 10 U Y | 5 UD | 5 U | 5 U | 5 U | 5 U | 5 U | | | |
| 1,2-Dichlorobenzene | 3 | | 10 U Y | 5 UD Y | 5 U Y | 5 U Y | 5 U Y | 5 U Y | 5 U Y | | | |
| 1,3-Dichlorobenzene | 3 | | 10 U Y | 5 UD Y | 5 U Y | 5 U Y | 5 U Y | 5 U Y | 5 U Y | | | |
| 1,4-Dichlorobenzene | 3 | | 10 U Y | 5 UD Y | 5 U Y | 5 U Y | 5 U Y | 5 U Y | 5 U Y | | | |
| 2,4,5-Trichlorophenol | NS | | 51 U | 26 UD | 26 U | 25 U | 26 U | 26 U | 27 U | | | |
| 2,4,6-Trichlorophenol | NS | | 10 U | 5 UD | 5 U | 5 U | 5 U | 5 U | 5 U | | | |
| 2,4-Dichlorophenol | 1 | | 10 U Y | 5 UD Y | 5 U Y | 5 U Y | 5 U Y | 5 U Y | 5 U Y | | | |
| 2,4-Dimethylphenol | 1 | | 10 U Y | 5 UD Y | 5 U Y | 5 U Y | 5 U Y | 5 U Y | 5 U Y | | | |
| 2,4-Dinitrophenol | 1 | | 51 U Y | 26 UD Y | 26 U Y | 25 U Y | 26 U Y | 26 U Y | 27 U Y | | | |
| 2,4-Dinitrotoluene | 5* | | 10 U Y | 5 UD | 5 U | 5 U | 5 U | 5 U | 5 U | | | |
| 2,6-Dinitrotoluene | 5* | | 10 U Y | 5 UD | 5 U | 5 U | 5 U | 5 U | 5 U | | | |
| 2-Chloronaphthalene | NS | | 20 U | 10 UD | 10 U | 10 U | 10 U | 10 U | 11 U | | | |
| 2-Chlorophenol | NS | | 10 U | 5 UD | 5 U | 5 U | 5 U | 5 U | 5 U | | | |
| 2-Methylnaphthalene | NS | | 27 D | 10 UD | 10 U | 10 U | 10 U | 15 | 11 U | | | |
| 2-Methylphenol | 1 | | 10 U Y | 5 UD Y | 5 U Y | 5 U Y | 5 U Y | 5 U Y | 5 U Y | | | |
| 2-Nitroaniline | 5* | | 51 U Y | 26 UD Y | 26 U Y | 25 U Y | 26 U Y | 26 U Y | 27 U Y | | | |
| 2-Nitrophenol | NS | | 10 U | 5 UD | 5 U | 5 U | 5 U | 5 U | 5 U | | | |
| 3,3-Dichlorobenzidine | 5* | | 20 U Y | 10 UD Y | 10 U Y | 10 U Y | 10 U Y | 10 U Y | 11 U Y | | | |
| 3-Nitroaniline | 5* | | 51 U Y | 26 UD Y | 26 U Y | 25 U Y | 26 U Y | 26 U Y | 27 U Y | | | |
| 4,6-Dinitro-2-methylphenol | NS | | 100 U | 52 UD | 51 U | 51 U | 51 U | 51 U | 55 U | | | |
| 4-Bromophenyl phenyl ether | NS | | 20 U | 10 UD | 10 U | 10 U | 10 U | 10 U | 11 U | | | |
| 4-Chloro-3-methylphenol | NS | | 10 U | 5 UD | 5 U | 5 U | 5 U | 5 U | 5 U | | | |
| 4-Chloroaniline | 5* | | 10 U Y | 5 UD | 5 U | 5 U | 5 U | 5 U | 5 U | | | |
| 4-Chlorophenyl phenyl ether | NS | | 20 U | 10 UD | 10 U | 10 U | 10 U | 10 U | 11 U | | | |
| 4-Methylphenol | 1 | | 20 U Y | 10 UD Y | 10 U Y | 10 U Y | 10 U Y | 10 U Y | 11 U Y | | | |
| 4-Nitroaniline | 5* | | 51 U Y | 26 UD Y | 26 U Y | 25 U Y | 26 U Y | 26 U Y | 27 U Y | | | |
| 4-Nitrophenol | NS | | 51 U | 26 UD | 26 U | 25 U | 26 U | 26 U | 27 U | | | |
| Acenaphthene | NS | | 20 U | 10 UD | 10 U | 10 U | 10 U | 10 U | 11 U | | | |
| Acenaphthylene | NS | | 20 U | 10 UD | 10 U | 10 U | 10 U | 10 U | 11 U | | | |
| Anthracene | NS | | 20 U | 10 UD | 10 U | 10 U | 10 U | 10 U | 11 U | | | |
| Benz(a)anthracene | NS | | 10 U | 5 UD | 5 U | 5 U | 5 U | 5 U | 5 U | | | |
| Benz(b)fluoranthene | NS | | 20 U | 10 UD | 10 U | 10 U | 10 U | 10 U | 11 U | | | |
| Benz(b)perylene | NS | | 20 U | 10 UD | 10 U | 10 U | 10 U | 10 U | 11 U | | | |
| Benz(k)fluoranthene | NS | | 20 U | 10 UD | 10 U | 10 U | 10 U | 10 U | 11 U | | | |
| Benzo(a)pyrene | NS | | 20 U | 10 UD | 10 U | 10 U | 10 U | 10 U | 11 U | | | |

NOTES: U - not detected, J - estimated value, E - outside linear range, D - from diluted analysis, Y - exceeds GW Standard, --- - not analyzed.

* - principal organic contaminant standard for ground water is 5 ug/L.



**Honeywell
Tonawanda, NY**

Table 5
Semivolatile Organic Compound Data
Ground Water Samples

| Compound | Sample ID | NYSDEC Class GA GW Standards | MW-07 DL | MW-08 | MW-09 | MW-10 | MW-11 |
|-----------------------------------|-----------|------------------------------------|------------------|------------------|------------------|------------------|------------------|
| | | | 06/06/01 ug/L | 05/12/99 ug/L | 06/06/01 ug/L | 06/05/01 ug/L | 06/07/01 ug/L |
| Bis(2-chloroethyl)ether | 5* | 10 U Y | 5 UD | 5 U | 5 U | 5 U | 5 U |
| Bis(2-chloroisopropyl) ether | 1 | 10 U Y | 5 UD Y | 5 U Y | 5 U Y | 5 U Y | 5 U Y |
| Bis(2-chloroisopropyl) ether | NS | 20 U | 10 UD | 10 U | 10 U | 10 U | 11 U |
| Bis(2-ethylhexyl)phthalate (BEHP) | 5 | 10 U Y | 5 UD | 5 U | 5 U | 5 U | 12 Y |
| Butyl benzyl phthalate | NS | 20 U | 10 UD | 10 U | 10 U | 10 U | 11 U |
| Carbazole | NS | 20 U | 10 UD | 10 U | 1 J | 13 | 11 U |
| Chrysene | NS | 10 U | 5 UD | 5 U | 5 U | 5 U | 5 U |
| Di-n-butyl phthalate | 50 | 20 U | 10 UD | 10 U | 10 U | 10 U | 11 U |
| Di-n-octyl phthalate | NS | 20 U | 10 UD | 10 U | 10 U | 10 U | 11 U |
| Dibenz(a,b)anthracene | NS | 20 U | 10 UD | 10 U | 10 U | 10 U | 11 U |
| Dibenzofuran | NS | 10 U | 5 UD | 5 U | 5 U | 5 U | 5 U |
| Diethyl phthalate | NS | 20 U | 10 UD | 10 U | 10 U | 10 U | 11 U |
| Dimethyl phthalate | NS | 20 U | 10 UD | 10 U | 10 U | 10 U | 11 U |
| Fluoranthene | NS | 20 U | 10 UD | 10 U | 10 U | 3 J | 11 U |
| Fluorene | NS | 3 J | 10 UD | 10 U | 10 U | 19 | 11 U |
| Hexachlorobenzene | 0.04 | 10 U Y | 5 UD Y | 5 U Y | 5 U Y | 5 U Y | 5 U Y |
| Hexachlorobutadiene | 0.5 | 10 U Y | 5 UD Y | 5 U Y | 5 U Y | 5 U Y | 5 U Y |
| Hexachlorocyclopentadiene | 5* | 20 U | 10 UD Y | 10 U Y | 10 U Y | 10 U Y | 11 U Y |
| Hexachloroethane | 5* | 10 U Y | 5 UD | 5 U | 5 U | 5 U | 5 U |
| Indeno(1,2,3-cd)pyrene | NS | 20 U | 10 UD | 10 U | 10 U | 10 U | 11 U |
| Iophorone | NS | 20 U | 10 UD | 10 U | 10 U | 10 U | 11 U |
| N-Nitrosodiphenylamine | NS | 20 U | 10 UD | 10 U | 10 U | 10 U | 11 U |
| N-Nitrosodipropylamine | NS | 20 U | 10 UD | 10 U | 10 U | 10 U | 11 U |
| Naphthalene | NS | 250 D | 10 UD | 10 U | 10 U | 100 | 11 U |
| Nitrobenzene | 0.4 | 10 U Y | 5 UD Y | 5 U Y | 5 U Y | 5 U Y | 5 U Y |
| Pentachlorophenol | 1 | 51 U Y | 26 UD Y | 26 U Y | 25 U Y | 26 U Y | 27 U Y |
| Phenanthrene | NS | 20 U | 10 UD | 10 U | 10 U | 17 | 11 U |
| Phenol | 1 | 10 U Y | 5 UD Y | 5 U Y | 5 U Y | 1 J | 5 U Y |
| Pyrene | NS | 20 U | 10 UD | 10 U | 10 U | 2 J | 11 U |

NOTES: U - not detected, J - estimated value, E - outside linear range, D - from diluted analysis, Y - exceeds GW Standard, -- - not analyzed.

* - principal organic contaminant standard for ground water is 5 ug/L.



Table 6
Honeywell
Tonawanda, NY
Metals Data

Ground Water Samples

| Compound | Sample ID NYSDEC Class GA GW Standards mg/L | MW-01 05/12/99 mg/L | MW-01 | MW-01 DUP | MW-02 | MW-02 |
|-----------|---|---------------------------|-----------|------------------|------------------|------------------|
| | | | Units | 06/05/01 mg/L | 05/12/99 mg/L | 06/09/01 mg/L |
| | NS | 32.2 | 22.6 | 3.12 | 0.0709 J | 9.34 |
| Aluminum | NS | 0.003 | 0.0016 UD | 0.0027 J | 0.0018 J | 0.0022 J Y |
| Antimony | 0.025 | 0.0019 UD | 0.0016 U | 0.0016 U | 0.0037 J | 0.0076 |
| Barium | 1 | 0.0163 J | 0.0122 J | 0.0123 J | 0.102 | 0.0644 J |
| Beryllium | NS | 0.0039 J | 0.0029 J | 0.0029 J | 0.00016 J | 0.00023 J |
| Cadmium | 5 | 0.0058 | 0.00095 J | 0.0012 J | 0.00042 UD | 0.00042 UD |
| Calcium | NS | 273 | 294 | 290 | 512 | 504 |
| Chromium | 0.050 | 0.156 Y | 0.131 Y | 0.131 Y | 0.0121 | 0.0047 J |
| Cobalt | NS | 0.0016 UD | 0.00093 U | 0.00093 U | 0.0016 UD | 0.00093 U |
| Copper | 0.2 | 0.0059 J | 0.0012 J | 0.0017 J | 0.0112 | 0.0119 J |
| Cyanide | 0.2 | 0.54 Y | 0.50 Y | 0.39 Y | 3.4 Y | 2.10 Y |
| Iron | 0.3 | 371 Y | 307 Y | 300 Y | 4.4 Y | 9.15 Y |
| Lead | 0.025 | 0.0011 UD | 0.00066 U | 0.00066 U | 0.0166 | 0.0016 J |
| Magnesium | NS | 78.0 | 100 | 98.1 | 119 | 136 |
| Manganese | 0.3 | 4.34 Y | 4.52 Y | 4.43 Y | 0.786 Y | 2.69 Y |
| Mercury | 0.0007 | 0.00011 UD | 0.00018 U | 0.00018 U | 0.00014 J | 0.00018 U |
| Nickel | 0.1 | 0.0455 J | 0.0400 J | 0.0414 J | 0.0174 J | 0.0046 J |
| Potassium | NS | 5.13 | 4.37 J | 4.38 J | 16.3 | 13.8 |
| Selenium | 0.05 | 0.0145 | 0.0047 J | 0.0057 | 0.0036 UD | 0.0026 J |
| Silver | 0.05 | 0.0010 UD | 0.00073 U | 0.00073 U | 0.0010 UD | 0.00073 U |
| Sodium | 20 | 30.8 Y | 37.4 Y | 36.5 Y | 87.1 Y | 112 Y |
| Thallium | NS | 0.0038 UD | 0.0036 U | 0.0036 U | 0.0038 UD | 0.0036 U |
| Vanadium | NS | 0.0126 J | 0.0075 J | 0.0075 J | 0.0052 J | 0.0021 J |
| Zinc | NS | 0.351 | 0.277 | 0.281 | 0.0331 | 0.00997 U |
| | | | | | | 0.117 |

NOTES: U - not detected, J - estimated value, E - outside linear range, D - from diluted analysis, Y - exceeds GW Standard, -- - not analyzed.
 * - principal organic contaminant standard for ground water is 5 µg/L.



Table 6
Honeywell
Tonawanda, NY
Metals Data

Ground Water Samples

| Compound | Sample ID | NYSDEC Class GA GW Standards | MW-03 | MW-04 | MW-04 | MW-05 | MW-05 | MW-06 |
|-----------|-----------|------------------------------------|-------------|------------|------------|------------|-----------|----------|
| | | | 06/09/01 | 05/12/99 | 06/09/01 | 05/12/99 | 06/07/01 | 05/12/99 |
| Units | mg/L | mg/L | mg/L | mg/L | mg/L | mg/L | mg/L | mg/L |
| Aluminum | NS | 0.0118 U | 4.4 | 0.0118 U | 12.4 | 0.0118 U | 10.7 | |
| Antimony | 0.003 | 0.0018 J | 0.0016 UD | 0.0024 J | 0.0016 UD | 0.0036 J Y | 0.0018 J | |
| Arsenic | 0.025 | 0.0016 U | 0.0034 J | 0.0078 | 0.0082 | 0.0020 J | 0.0029 J | |
| Barium | 1 | 0.0360 J | 0.135 | 0.0572 J | 0.200 | 0.0626 J | 0.0313 J | |
| Beryllium | NS | 0.000076 U | 0.00029 J | 0.000076 U | 0.00061 J | 0.000076 U | 0.0012 J | |
| Cadmium | 5 | 0.00024 U | 0.00042 UD | 0.00024 U | 0.00042 UD | 0.00024 U | 0.00020 J | |
| Calcium | NS | 115 | 240 | 290 | 161 | 170 | 546 | |
| Chromium | 0.050 | 0.0015 J | 0.0162 | 0.0036 J | 0.0657 Y | 0.0037 J | 0.109 Y | |
| Cobalt | NS | 0.00093 U | 0.0031 J | 0.00093 U | 0.0089 J | 0.0026 J | 0.0094 J | |
| Copper | 0.2 | 0.00089 J | 0.0089 J | 0.00049 U | 0.105 | 0.0056 J | 0.0270 | |
| Cyanide | 0.2 | 0.36 Y | 0.78 Y | 0.63 Y | 0.46 Y | 0.53 Y | 4.9 Y | |
| Iron | 0.3 | 0.228 | 7.6 Y | 8.00 Y | 24.5 Y | 0.549 Y | 50.6 Y | |
| Lead | 0.025 | 0.00087 J | 0.0618 Y | 0.00066 U | 0.212 Y | 0.0024 J | 0.0238 | |
| Magnesium | NS | 14.8 | 58.7 | 43.8 | 166 | 182 | 80.8 | |
| Manganese | 0.3 | 0.363 Y | 1.44 Y | 1.70 Y | 3.98 Y | 3.57 Y | 20.4 Y | |
| Mercury | 0.0007 | 0.00018 U | 0.000111 UD | 0.00018 U | 0.00038 | 0.00018 U | 0.00035 | |
| Nickel | 0.1 | 0.00097 J | 0.0061 J | 0.0021 J | 0.150 Y | 0.0447 J | 0.0594 | |
| Potassium | NS | 5.69 | 2.34 J | 3.25 J | 3.82 J | 2.66 J | 30.4 | |
| Selenium | 0.05 | 0.0018 U | 0.0036 UD | 0.0018 U | 0.0036 UD | 0.0020 J | 0.0180 UD | |
| Silver | 0.05 | 0.00073 U | 0.0010 UD | 0.00073 U | 0.0012 J | 0.00073 U | 0.0010 UD | |
| Sodium | 20 | 9.77 | 49.1 Y | 56.0 Y | 67.0 Y | 66.5 Y | 54.5 Y | |
| Thallium | NS | 0.0036 U | 0.0045 J | 0.0036 U | 0.0038 UD | 0.0036 U | 0.0190 UD | |
| Vanadium | NS | 0.00041 J | 0.0069 J | 0.0011 J | 0.0279 J | 0.0029 J | 0.185 | |
| Zinc | NS | 0.0032 J | 0.0651 | 0.00097 U | 0.108 | 0.00097 U | 0.152 | |

NOTES: U - not detected, J - estimated value, E - outside linear range D - from diluted analysis, Y - exceeds GW Standard, -- not analyzed.

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FXP File: N:\1163\22900\METALSMW.FXP



Table 6
Honeywell
Tonawanda, NY
Metals Data

Ground Water Samples

| Compound | Sample ID | NYSDEC Class GA GW Standards | MW-06 | MW-07 | MW-07 | MW-08 | MW-08 |
|-----------|-----------|------------------------------------|-------------|------------|------------|------------|------------|
| | | | Sample Date | 06/06/01 | 05/12/99 | 06/06/01 | 05/12/99 |
| | | | Units | mg/L | mg/L | mg/L | mg/L |
| Aluminum | NS | 13.2 | 37.6 | 18.6 | 35.6 | 7.34 | 0.0162 J |
| Antimony | 0.003 | 0.0019 J | 0.0073 J Y | 0.0081 J Y | 0.0016 UD | 0.0014 U | 0.0014 U |
| Arsenic | 0.025 | 0.0016 U | 0.0373 Y | 0.0440 Y | 0.0092 | 0.0021 J | 0.0021 J |
| Barium | 1 | 0.0112 J | 0.0730 J | 0.0081 J | 0.321 | 0.0317 J | 0.0648 J |
| Beryllium | NS | 0.0017 J | 0.0052 | 0.0027 J | 0.0018 J | 0.000076 U | 0.000076 U |
| Cadmium | 5 | 0.00024 U | 0.0083 | 0.0024 J | 0.00063 J | 0.00024 U | 0.00024 U |
| Calcium | NS | 490 | 355 | 309 | 140 | 91.2 | 246 |
| Chromium | 0.050 | 0.0661 Y | 0.301 Y | 0.0220 Y | 0.0626 Y | 0.0668 J | 0.0031 J |
| Cobalt | NS | 0.00093 U | 0.0213 J | 0.0297 J | 0.0161 J | 0.0093 U | 0.00093 U |
| Copper | 0.2 | 0.00049 U | 0.130 | 0.0096 J | 0.0429 | 0.0017 J | 0.00091 J |
| Cyanide | 0.2 | 1.80 Y | 3.6 Y | 0.75 Y | 0.52 Y | 0.40 Y | 1.10 Y |
| Iron | 0.3 | 77.6 Y | 383 Y | 277 Y | 49.6 Y | 2.04 Y | 2.96 Y |
| Lead | 0.025 | 0.0013 J | 0.158 Y | 0.143 Y | 0.015 Y | 0.00066 U | 0.00084 J |
| Magnesium | NS | 53.1 | 56.6 | 34.1 | 190 | 198 | 40.9 |
| Manganese | 0.3 | 0.961 Y | 4.55 Y | 2.15 Y | 0.822 Y | 0.105 | 2.72 Y |
| Mercury | 0.0007 | 0.00018 U | 0.00014 Y | 0.00018 U | 0.00011 UD | 0.00018 U | 0.00018 U |
| Nickel | 0.1 | 0.0029 J | 0.351 Y | 0.259 Y | 0.0507 | 0.030 J | 0.0046 J |
| Potassium | NS | 9.36 | 8.86 | 7.10 | 12.0 | 3.92 J | 18.9 |
| Selenium | 0.05 | 0.0018 U | 0.0152 | 0.0041 J | 0.0036 UD | 0.0018 U | 0.0021 J |
| Silver | 0.05 | 0.00073 U | 0.0010 UD | 0.00073 U | 0.0010 UD | 0.00073 U | 0.00073 U |
| Sodium | 20 | 30.0 Y | 18.6 | 12.8 | 11.8 Y | 131 Y | 36.0 Y |
| Thallium | NS | 0.0036 U | 0.0038 UD | 0.0036 U | 0.0038 UD | 0.0036 U | 0.0036 U |
| Vanadium | NS | 0.0141 J | 0.0560 | 0.0154 J | 0.0607 | 0.0211 J | 0.00072 J |
| Zinc | NS | 0.0623 | 0.832 | 0.597 | 0.146 | 0.0097 U | 0.00097 U |

NOTES: U - not detected, J - estimated value, E - outside linear range, D - from diluted analysis, Y - exceeds GW Standard, -- - not analyzed
 * - principal organic contaminant standard for ground water is 5 ug/L.



Table 6
Honeywell
Tonawanda, NY
Metals Data

Ground Water Samples

| Compound | Sample ID Sample Date Units | NYSDEC Class GA GW Standards mg/L | MW-10 | | MW-10 DL | | MW-11 | |
|-----------|-----------------------------------|--|------------|---------|-------------|-----------|------------|------|
| | | | 06/05/01 | | 06/05/01 | | 06/07/01 | |
| | | | mg/L | mg/L | mg/L | mg/L | mg/L | mg/L |
| Aluminum | NS | 5.90 | | 6.09 D | | 6.0128 J | | |
| Antimony | 0.003 | | 0.0032 J Y | | 0.0076 JD Y | | 0.0014 U | |
| Arsenic | 0.025 | | 0.0016 U | | 0.0080 UD | | 0.0016 J | |
| Barium | 1 | | 0.0288 J | | 0.0326 JD | | 0.332 | |
| Beryllium | NS | | 0.0023 J | | 0.0022 JD | | 0.000076 U | |
| Cadmium | 5 | | 0.0030 J | | 0.0034 JD | | 0.00024 U | |
| Calcium | NS | 491 | | 492 D | | 85.8 | | |
| Chromium | 0.050 | | 0.0285 | | 0.0534 D Y | | 0.0062 J | |
| Cobalt | NS | | 0.00093 U | | 0.0046 UD | | 0.00093 U | |
| Copper | 0.2 | | 0.0056 J | | 0.0058 JD | | 0.0043 J | |
| Cyanide | 0.2 | | 0.10 | | -- | | 0.022 | |
| Iron | 0.3 | | 88.5 E Y | | 879 D Y | | 0.0312 J | |
| Lead | 0.025 | | 0.00066 U | | 0.0033 UD | | 0.00066 U | |
| Magnesium | NS | 204 | | 211 D | | 177 | | |
| Manganese | 0.3 | | 36.6 E Y | | 42.3 D Y | | 0.914 Y | |
| Mercury | 0.0007 | | 0.00018 U | | -- | | 0.00018 U | |
| Nickel | 0.1 | | 0.00071 U | | 0.0036 UD | | 0.0142 J | |
| Potassium | NS | 16.0 | | 14.7 JD | | 12.5 | | |
| Selenium | 0.05 | | 0.0350 | | 0.0320 D | | 0.0048 J | |
| Silver | 0.05 | | 0.00073 U | | 0.0036 UD | | 0.00073 U | |
| Sodium | 20 | | 23.1 Y | | 20.2 D Y | | 34.4 Y | |
| Thallium | NS | | 0.0036 U | | 0.0180 UD | | 0.0036 U | |
| Vanadium | NS | | 0.0049 J | | 0.0062 JD | | 0.0024 J | |
| Zinc | NS | 0.358 | | 0.368 D | | 0.00097 U | | |

NOTES: U - not detected, J - estimated value, E - outside linear range, D - from diluted analysis, Y - exceeds GW Standard, -- - not analyzed
* - principal organic contaminant standard for ground water is 5 ug/L



Table 7
AlliedSignal Inc
Tonawanda, NY
Volatile Organic Compound Data

| Compound | Storm Water Samples | | | | | | SW Outfall #1 05/10/1999 ug/L | SW Outfall #1 12/11/2001 ug/L | SW Outfall #2 05/10/1999 ug/L | | | |
|-----------------------------|---------------------|------------|------------|------------|------------|------|-------------------------------------|-------------------------------------|-------------------------------------|--|--|--|
| | Sample ID | | SW Inlet A | | SW Inlet B | | | | | | | |
| | Sample Date | 05/11/1999 | 12/11/2001 | 05/11/1999 | 12/11/2001 | | | | | | | |
| Units | ug/L | ug/L | ug/L | ug/L | ug/L | ug/L | | | | | | |
| 1,1,1-Trichloroethane | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.1 J | | | |
| 1,1,2,2-Tetrachloroethane | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | | | |
| 1,1,2-Trichloroethane | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | | | |
| 1,1-Dichloroethane | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | | | |
| 1,1-Dichloroethene | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | | | |
| 1,2-Dichloroethane | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | | | |
| 1,2-Dichloropropane | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | | | |
| 2-Butanone (MEK) | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U | 10 UD | | | |
| 2-Hexanone | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 10 U | 5 UD | | | |
| 4-Methyl-2-pentanone (MIBK) | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 10 U | 5 UD | | | |
| Acetone | 2 J | 10 U | 4 J | 10 U | 4 J | 10 U | 4 J | 3 J | 6 J | | | |
| Benzene | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | | | |
| Bromodichloromethane | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.2 J | | | |
| Bromoform | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | | | |
| Bromomethane | 1 UD | 10 U | 1 UD | 10 U | 1 UD | 10 U | 1 UD | 10 U | 1 UD | | | |
| Carbon disulfide | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.2 J | 0.8 J | 0.5 UD | | | |
| Carbon tetrachloride | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | | | |
| Chlorobenzene | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | | | |
| Chloroethane | 1 UD | 10 U | 1 UD | 10 U | 1 UD | 10 U | 1 UD | 10 U | 1 UD | | | |
| Chloroform | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.3 J | | | |
| Chloromethane | 1 UD | 10 U | 1 UD | 10 U | 1 UD | 10 U | 1 UD | 10 U | 1 UD | | | |
| Dibromochloromethane | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | | | |
| Ethane, 1,2-dichloro, (E)- | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | | | |
| Ethybenzene | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.4 J | 10 U | 0.4 J | 0.6 J | 0.5 UD | | | |
| Methylene chloride | 2 UD | 1 J | 2 UD | 1 J | 2 UD | 1 J | 2 UD | 1 J | 2 UD | | | |
| Styrene | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | | | |
| Tetrachloroethene | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.4 J | | | |
| Toluene | 0.1 J | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.9 | 2 J | 0.5 UD | | | |
| Trichloroethene | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.9 | 10 U | 0.9 | 10 U | 0.2 J | | | |
| Vinyl chloride | 1 UD | 10 U | 1 UD | 10 U | 1 UD | 10 U | 1 UD | 10 U | 0.5 J | | | |
| Xylene (total) | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 2 | 2 J | 0.5 UD | | | |
| cis-1,2-Dichloroethene | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 8 | 0.8 J | 1 | | | |
| cis-1,3-Dichloropropylene | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | | | |
| trans-1,3-Dichloropropene | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | 10 U | 0.5 UD | | | |

NOTES: U - Not detected. J - Estimated value. --- - Not analyzed.



Table 7
AlliedSignal Inc
Tonawanda, NY
Volatile Organic Compound Data
Storm Water Samples

| Compound | Sample ID | SW Outfall #2 |
|-----------------------------|-------------|---------------|
| | Sample Date | 12/11/2001 |
| | Units | ug/L |
| 1,1,1-Trichloroethane | 10 U | |
| 1,1,2,2-Tetrachloroethane | 10 U | |
| 1,1,2-Trichloroethane | 10 U | |
| 1,1-Dichloroethane | 10 U | |
| 1,1-Dichloroethene | 10 U | |
| 1,2-Dibromoethane | 10 U | |
| 1,2-Dichloropropane | 10 U | |
| 2-Butanone (MEK) | 10 U | |
| 2-Hexanone | 10 U | |
| 4-Methyl-2-pentanone (MIBK) | 10 U | |
| Acetone | 3 J | |
| Benzene | 10 U | |
| Bromodichloromethane | 10 U | |
| Bromoform | 10 U | |
| Bromomethane | 10 U | |
| Carbon disulfide | 10 U | |
| Carbon tetrachloride | 10 U | |
| Chlorobenzene | 10 U | |
| Chloroethane | 10 U | |
| Chloroform | 10 U | |
| Chloromethane | 10 U | |
| Dibromodichloromethane | 10 U | |
| Ethene, 1,2-dichloro- (E)- | 10 U | |
| Ethylbenzene | 10 U | |
| Methylene chloride | 1 J | |
| Styrene | 10 U | |
| Tetrachloroethene | 10 U | |
| Toluene | 10 U | |
| Trichloroethene | 10 U | |
| Vinyl chloride | 0.5 J | |
| Xylene (total) | 10 U | |
| dis-1,2-Dichloroethene | 0.7 J | |
| cis-1,3-Dichloropropene | 10 U | |
| trans-1,3-Dichloropropene | 10 U | |

NOTES: U - Not detected. J - Estimated value. --- Not analyzed



Table 8
AlliedSignal Inc
Tonawanda, NY
Semivolatile Organic Compound Data

| Compound | Sample ID | SW Inlet A | | SW Inlet B | | SW Outfall #1 | | SW Outfall #2 | |
|-----------------------------|-----------|-------------|------------|------------|------------|---------------|-------|---------------|-------|
| | | Sample Date | 05/11/1999 | 05/11/1999 | 12/11/2001 | ug/L | ug/L | 12/11/2001 | ug/L |
| | | | 12/11/2001 | | ug/L | | ug/L | | |
| 1,2,4-Trichlorobenzene | | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 5 UD | 10 U | 5 UD |
| 1,2-Dichlorobenzene | | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 5 UD | 10 U | 5 UD |
| 1,3-Dichlorobenzene | | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 5 UD | 10 U | 5 UD |
| 1,4-Dichlorobenzene | | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 5 UD | 10 U | 5 UD |
| 2,4,5-Trichlorophenol | | 25 UD | 25 U | 25 UD | 26 U | 25 UD | 25 U | 25 U | 25 UD |
| 2,4,6-Trichlorophenol | | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 5 UD | 10 U | 5 UD |
| 2,4-Dichlorophenol | | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 5 UD | 10 U | 5 UD |
| 2,4-Dimethylphenol | | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 5 UD | 10 U | 5 UD |
| 2,4-Dinitrophenol | | 25 UD | 25 U | 25 UD | 26 U | 25 UD | 25 U | 25 U | 25 UD |
| 2,4-Dinitrotoluene | | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 5 UD | 10 U | 5 UD |
| 2,6-Dinitrotoluene | | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 5 UD | 10 U | 5 UD |
| 2-Chloronaphthalene | | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 UD | 10 U | 10 UD |
| 2-Chlorophenol | | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 5 UD | 10 U | 5 UD |
| 2-Methylnaphthalene | | 10 UD | 10 U | 10 UD | 10 U | 4 J | 6 J | 10 UD | 5 UD |
| 2-Methylphenol | | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 5 UD | 10 U | 5 UD |
| 2-Nitroaniline | | 25 UD | 25 U | 25 UD | 26 U | 25 UD | 25 U | 25 U | 25 UD |
| 2-Nitrophenol | | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 5 UD | 10 U | 5 UD |
| 3,3-Dichlorobenzidine | | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 UD | 10 U | 10 UD |
| 3-Nitroaniline | | 25 UD | 25 U | 25 UD | 26 U | 25 UD | 25 U | 25 U | 25 UD |
| 4,6-Dinitro-2-methylphenol | | 50 UD | 25 U | 30 UD | 26 U | 51 UD | 25 U | 25 U | 50 UD |
| 4-Bromophenyl phenyl ether | | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 UD | 10 U | 10 UD |
| 4-Chloro-3-methylphenol | | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 5 UD | 10 U | 5 UD |
| 4-Chloraniline | | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 5 UD | 10 U | 5 UD |
| 4-Chlorophenyl phenyl ether | | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 UD | 10 U | 10 UD |
| 4-Methylphenol | | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 UD | 10 U | 10 UD |
| 4-Nitroaniline | | 25 UD | 25 U | 25 UD | 26 U | 25 UD | 25 U | 25 U | 25 UD |
| 4-Nitrophenol | | 25 UD | 25 U | 25 UD | 26 U | 25 UD | 25 U | 25 U | 25 UD |
| Acenaphthene | | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 UD | 10 U | 10 UD |
| Acenaphthylene | | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 UD | 10 U | 10 UD |
| Anthracene | | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 UD | 10 U | 10 UD |
| Benz(a)anthracene | | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 5 UD | 10 U | 5 UD |
| Benz(b)fluoranthene | | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 UD | 10 U | 10 UD |
| Benz(g)fluoranthene | | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 UD | 10 U | 10 UD |
| Benz(k)fluoranthene | | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 UD | 10 U | 10 UD |
| Benzo(a)pyrene | | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 UD | 10 U | 10 UD |

NOTES: U - Not detected. J - Estimated value. -- - Not analyzed.



Table 8
AlliedSignal Inc
Tonawanda, NY
Semivolatile Organic Compound Data

| Compound | Sample ID | SW Inlet A | | SW Inlet B | | SW Outfall #1 | | SW Outfall #2 | |
|-----------------------------------|-----------|-------------|------------|------------|------|---------------|------|---------------|-------|
| | | Sample Date | 05/11/1999 | 05/11/1999 | | 05/10/1999 | | 12/11/2001 | |
| | | | | ug/L | ug/L | ug/L | ug/L | ug/L | ug/L |
| Bis(2-chloroethoxy)methane | | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 10 U |
| Bis(2-chloroethyl)ether | | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 10 U |
| Bis(2-chloroisopropyl) ether | | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U |
| Bis(2-ethylhexyl)phthalate (BEHP) | | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 10 U |
| Butyl benzyl phthalate | | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U |
| Carbazole | | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U |
| Chrysene | | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 10 U |
| Di-n-butyl phthalate | | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U |
| Di-n-octyl phthalate | | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U |
| Dibenz(a,h)anthracene | | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U |
| Dibenzofuran | | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 10 U |
| Diethyl phthalate | | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U |
| Dimethyl phthalate | | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U |
| Fluoranthene | | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U |
| Fluorene | | 10 UD | 10 U | 5 UD | 10 U | 5 UD | 10 U | 2 J | 10 UD |
| Hexachlorobenzene | | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 10 U |
| Hexachlorobutadiene | | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 10 U |
| Hexachlorocyclopentadiene | | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U |
| Hexachloroethane | | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 10 U |
| Indeno(1,2,3-cd)pyrene | | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U |
| Iophorone | | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U |
| N-Nitrosodiphenylamine | | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U |
| N-Nitrosodipropylamine | | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U |
| Naphthalene | | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U | 25 | 10 UD |
| Nitrobenzene | | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 10 U |
| Pentachlorophenol | | 25 UD | 25 U | 25 UD | 25 U | 25 UD | 25 U | 25 UD | 25 U |
| Phenanthrene | | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U | 2 J | 10 UD |
| Phenol | | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 10 U | 5 UD | 10 U |
| Pyrene | | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U | 10 UD | 10 U |

NOTES: U - Not detected. J - Estimated value. -- - Not analyzed.



Table 8
AlliedSignal Inc
Tonawanda, NY
Semivolatile Organic Compound Data
Storm Water Samples

| Compound | Sample ID | SW Outfall #2 |
|-----------------------------|-------------|---------------|
| | Sample Date | 12/11/2001 |
| | Units | ug/L |
| 1,2,4-Trichlorobenzene | 10 U | |
| 1,2-Dichlorobenzene | 10 U | |
| 1,3-Dichlorobenzene | 10 U | |
| 1,4-Dichlorobenzene | 10 U | |
| 2,4,5-Trichlorophenol | 26 U | |
| 2,4,6-Trichlorophenol | 10 U | |
| 2,4-Dichlorophenol | 10 U | |
| 2,4-Dimethylphenol | 10 U | |
| 2,4-Dinitrophenol | 26 U | |
| 2,4-Dinitrotoluene | 10 U | |
| 2,6-Dinitrotoluene | 10 U | |
| 2-Chloronaphthalene | 10 U | |
| 2-Chlorophenol | 10 U | |
| 2-Methylnaphthalene | 10 U | |
| 2-Methylphenol | 10 U | |
| 2-Nitroaniline | 26 U | |
| 2-Nitrophenol | 10 U | |
| 3,3-Dichlorobenzidine | 10 U | |
| 3-Nitroaniline | 26 U | |
| 4,6-Dinitro-2-methylphenol | 26 U | |
| 4-Bromophenyl phenyl ether | 10 U | |
| 4-Chloro-3-methylphenol | 10 U | |
| 4-Chloraniline | 10 U | |
| 4-Chlorophenyl phenyl ether | 10 U | |
| 4-Methylphenol | 10 U | |
| 4-Nitroaniline | 26 U | |
| 4-Nitrophenol | 26 U | |
| Acenaphthene | 10 U | |
| Acenaphthylene | 10 U | |
| Anthracene | 10 U | |
| Benz(a)anthracene | 10 U | |
| Benz(b)fluoranthene | 10 U | |
| Benz(gi)perylene | 10 U | |
| Benz(k)fluoranthene | 10 U | |
| Benz(a)pyrene | 10 U | |

NOTES: U - Not detected J - Estimated value -- - Not analyzed.



Table 8
AlliedSignal Inc
Tonawanda, NY
Semivolatile Organic Compound Data
Storm Water Samples

| Compound | Sample ID | SW Outfall #2 |
|-----------------------------------|-------------|---------------|
| | Sample Date | 12/11/2001 |
| | Units | ug/L |
| Bis(2-chloroethoxy)methane | 10 U | |
| Bis(2-chloroethyl)ether | 10 U | |
| Bis(2-chloroisopropyl) ether | 10 U | |
| Bis(2-ethylhexyl)phthalate (BEHP) | 10 U | |
| Butyl benzyl phthalate | 10 U | |
| Carbazole | 10 U | |
| Chrysene | 10 U | |
| Di-n-butyl phthalate | 10 U | |
| Di-n-octyl phthalate | 10 U | |
| Dibenzoz(a,h)anthracene | 10 U | |
| Dibenzofuran | 10 U | |
| Diethyl phthalate | 10 U | |
| Dimethyl phthalate | 10 U | |
| Fluoranthene | 10 U | |
| Fluorene | 10 U | |
| Hexachlorobenzene | 10 U | |
| Hexachlorobutadiene | 10 U | |
| Hexachlorocyclopentadiene | 10 U | |
| Hexachloroethane | 10 U | |
| Indeno(1,2,3-cd)pyrene | 10 U | |
| Isophorone | 10 U | |
| N-Nitrosodiphenylamine | 10 U | |
| N-Nitrosodipropyamine | 10 U | |
| Naphthalene | 22 | |
| Nitrobenzene | 10 U | |
| Pentachlorophenol | 26 U | |
| Phenanthrene | 10 U | |
| Phenol | 10 U | |
| Pyrene | 10 U | |

NOTES: U - Not detected J - Estimated value -- - Not analyzed.



Table 9
AlliedSignal Inc
Tonawanda, NY
Metals Data

| Compound | Storm Water Samples | | | | | | SW Outfall #1 05/10/1999 mg/L | SW Outfall #2 12/11/2001 mg/L | | |
|-----------|---------------------|------------|------------|------------|------------|------------|-------------------------------------|-------------------------------------|--|--|
| | Sample ID | | SW Inlet A | | SW Inlet B | | | | | |
| | Sample Date | 05/11/1999 | 12/11/2001 | 05/11/1999 | 12/11/2001 | 05/11/1999 | | | | |
| Compound | Units | mg/L | mg/L | mg/L | mg/L | mg/L | mg/L | mg/L | | |
| Aluminum | 0.555 | 2.51 | 0.109 | 0.910 | 0.856 | 1.04 | 0.290 | | | |
| Antimony | 0.0016 UD | 0.0600 U | 0.0028 J | 0.0680 U | 0.0016 UD | 0.0600 U | 0.0035 J | | | |
| Arsenic | 0.0040 J | 0.01 U | 0.0019 UD | 0.01 U | 0.0019 UD | 0.01 U | 0.0019 UD | | | |
| Barium | 0.0524 J | 0.200 U | 0.0368 J | 0.200 U | 0.0467 J | 0.200 U | 0.0380 J | | | |
| Beryllium | 0.00013 UD | 0.0050 U | 0.00026 J | 0.0050 U | 0.00033 J | 0.0050 U | 0.00015 J | | | |
| Cadmium | 0.00042 UD | 0.0050 U | 0.00055 J | 0.0050 U | 0.0012 J | 0.0050 U | 0.00081 J | | | |
| Calcium | 123 | 156 | 71.4 | 83.4 | 158 | 182 | 72.5 | | | |
| Chromium | 0.0074 J | 0.0103 | 0.0054 J | 0.01 U | 0.0153 | 0.0178 | 0.0063 J | | | |
| Cobalt | 0.0016 UD | 0.0500 U | 0.0029 J | 0.0500 U | 0.0016 UD | 0.0500 U | 0.0027 J | | | |
| Copper | 0.0022 J | 0.0250 U | 0.0020 J | 0.0250 U | 0.0010 J | 0.0250 U | 0.0062 J | | | |
| Cyanide | 0.01 UD | 0.01 | 0.01 UD | 0.01 U | 0.01 UD | 0.12 | 0.01 UD | | | |
| Iron | 2.31 | 8.99 | 1.67 | 2.21 | 57.5 | 77.1 | 1.11 | | | |
| Lead | 0.0016 J | 0.0080 | 0.0011 UD | 0.0030 U | 0.0011 UD | 0.0030 U | 0.0030 J | | | |
| Magnesium | 34.7 | 45.1 | 14.2 | 19.0 | 44.7 | 51.2 | 13.8 | | | |
| Manganese | 0.995 | 1.24 | 1.53 | 0.652 | 1.59 | 1.92 | 0.743 | | | |
| Mercury | 0.00011 UD | 0.00020 U | 0.00011 UD | 0.00020 U | 0.00011 UD | 0.00020 U | 0.00011 UD | | | |
| Nickel | 0.0013 UD | 0.0400 U | 0.0033 J | 0.0400 U | 0.0066 J | 0.0400 U | 0.0034 J | | | |
| Potassium | 4.93 J | 5.00 U | 3.55 J | 5.00 U | 6.4 | 5.02 | 3.61 J | | | |
| Selenium | 0.0036 UD | 0.0050 U | 0.0036 UD | 0.0050 U | 0.0036 UD | 0.0050 U | 0.0036 UD | | | |
| Silver | 0.0010 UD | 0.01 U | 0.0010 UD | 0.01 U | 0.0010 UD | 0.01 U | 0.0010 UD | | | |
| Sodium | 43.2 | 43.2 | 23.5 | 19.5 | 41.2 | 39.0 | 23.4 | | | |
| Thallium | 0.0038 UD | 0.01 U | 0.0073 J | 0.01 U | 0.0061 J | 0.01 U | 0.0065 J | | | |
| Vanadium | 0.0018 J | 0.0500 U | 0.0060 UD | 0.0500 U | 0.0063 J | 0.0500 U | 0.0060 UD | | | |
| Zinc | 0.0193 | 0.418 | 0.0174 | 0.0200 U | 0.0577 | 0.0200 U | 0.0295 | | | |

NOTES: U - Not detected J - Estimated value. — - Not analyzed.



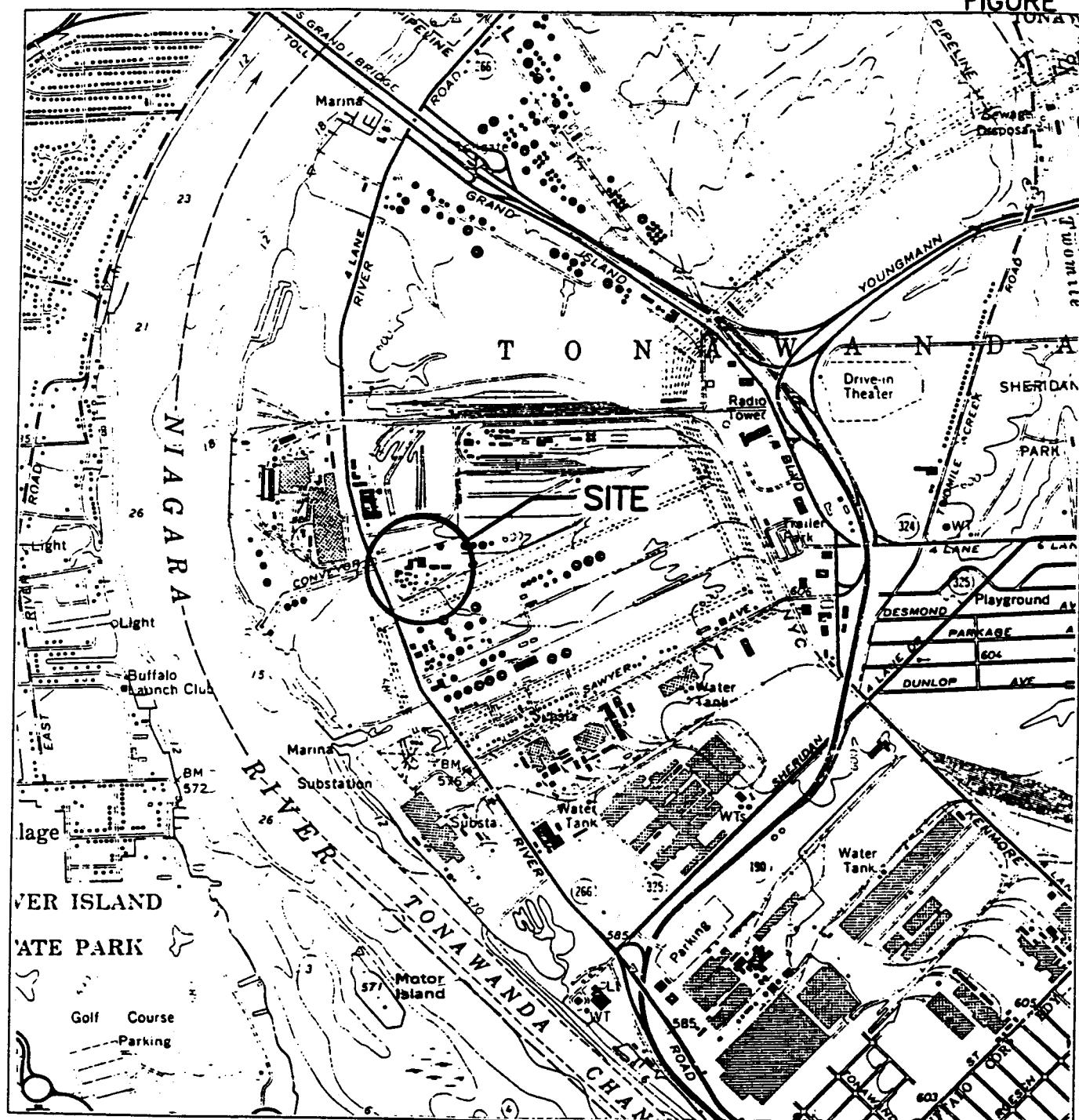
Table 9
AlliedSignal Inc
Tonawanda, NY
Metals Data
Storm Water Samples

| Compound | Sample ID | SW Outfall #2 |
|-----------|-------------|---------------|
| | Sample Date | 12/1/2001 |
| | Units | mg/L |
| Aluminum | | 1.51 |
| Antimony | | 0.0600 U |
| Arsenic | | 0.01 U |
| Barium | | 0.200 U |
| Beryllium | | 0.0050 U |
| Cadmium | | 0.0050 U |
| Calcium | | 84.9 |
| Chromium | | 0.01 U |
| Cobalt | | 0.0500 U |
| Copper | | 0.0250 U |
| Cyanide | | 0.02 |
| Iron | | 2.25 |
| Lead | | 0.0098 |
| Magnesium | | 18.2 |
| Manganese | | 0.466 |
| Mercury | | 0.00020 U |
| Nickel | | 0.0400 U |
| Potassium | | 5.00 U |
| Selenium | | 0.0050 U |
| Silver | | 0.01 U |
| Sodium | | 23.4 |
| Thallium | | 0.01 U |
| Vanadium | | 0.0500 U |
| Zinc | | 0.0200 U |

NOTES:

U - Not detected. J - Estimated value. --- Not analyzed

FIGURE 1



ADAPTED FROM: BUFFALO NW, N.Y.-ONT. 7.5 MIN. SERIES QUAD. (TOPO.)



FORMER ALLIED SPECIALTY
CHEMICAL SITE
TONAWANDA, NEW YORK

SITE LOCATION MAP

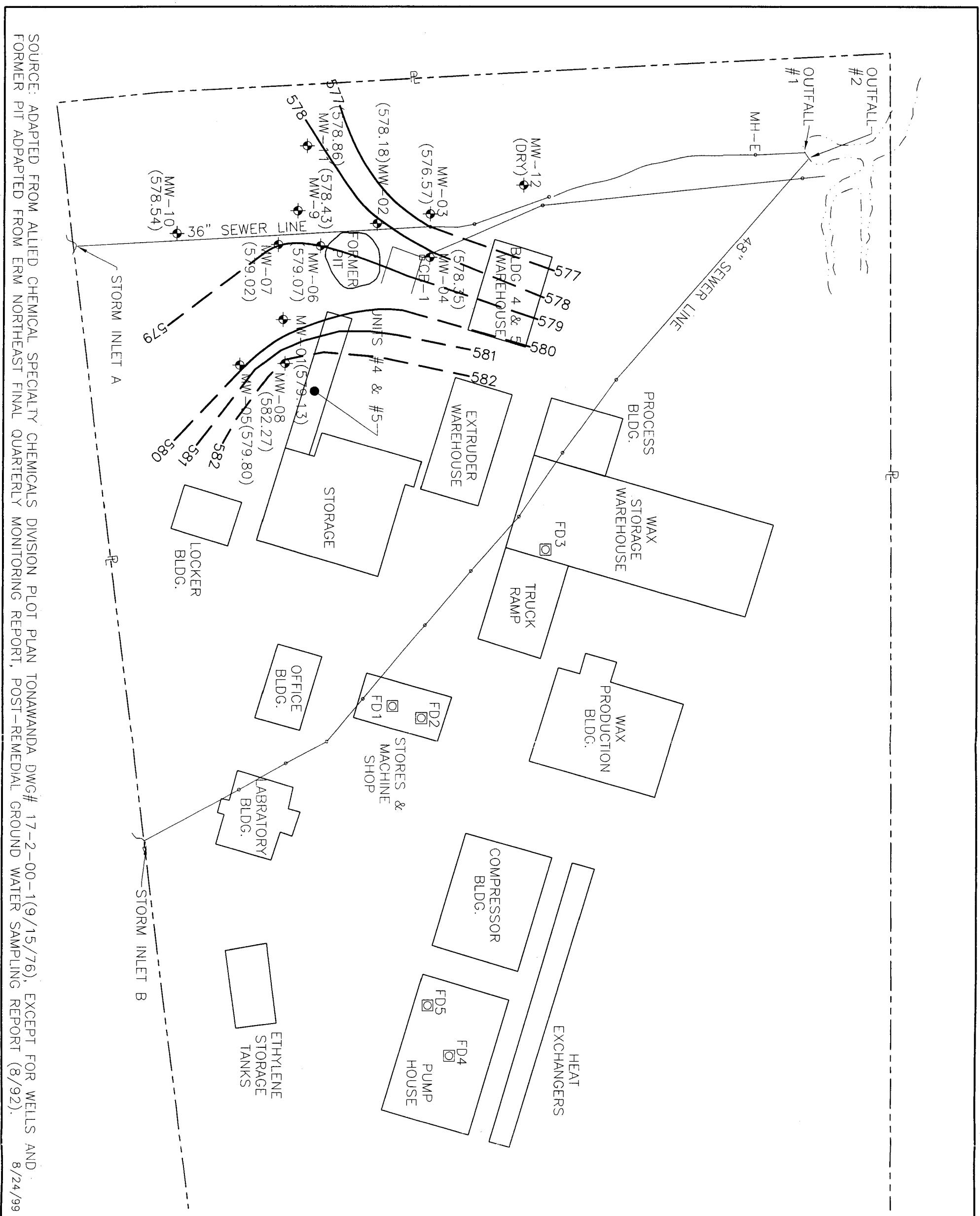


FILE NO. 1163.22900-001
DATE: 2/18/99

2000 0 2000 4000
1" = 2000'

O'Brien & Gere
ENGINEERS INC.

FIGURE 2



GROUND WATER
ELEVATIONS
(JUNE 5, 2001)

1"=80' 80' 0' 80'

AUGUST 1999
1163-26080.001

SOURCE: ADAPTED FROM ALLIED CHEMICAL SPECIALTY CHEMICALS DIVISION PLOT PLAN TONAWANDA DWG# 17-2-00-1(9/15/76), EXCEPT FOR WELLS AND FORMER PIT ADAPTED FROM ERM NORTHEAST FINAL QUARTERLY MONITORING REPORT, POST-REMEDIAL GROUND WATER SAMPLING REPORT (8/92). 8/24/99

**Laboratory Case Narratives
Sediment, Ground Water, and Storm
Water Samples**

SEDIMENT SAMPLES

ANALYTICAL PACKAGE

for

**Honeywell
Tonawanda, NY
Soil Samples**

Soil samples collected: November 15, 2000

Volume 1 of 4

Prepared for: O'Brien & Gere Engineers, Inc.
5000 Brittonfield Parkway
P.O. Box 4873
Syracuse, NY 13221

Prepared by: O'Brien & Gere Laboratories, Inc.
5000 Brittonfield Parkway
Suite 300, P.O. Box 4942
Syracuse, NY 13221

Authorized



Date

12-19-00

Reviewed



Date

12/20/00

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Sample Data Summary Package

NEW YORK STATE DEPARTMENT OF ENVIRONMENTAL CONSERVATION

**SAMPLE IDENTIFICATION AND
ANALYTICAL SUMMARY**

| Customer Sample Code | Laboratory Sample Code | Analytical Requirements | | | | | |
|----------------------|------------------------|-------------------------|---------------------|------------------|---------------------|---------|--------------------------|
| | | 'VOA GC/MS Method # | 'BNA GC/MS Method # | 'VOA GC Method # | 'Pest PCBs Method # | 'Metals | 'Other |
| Storm Inlet B | R6023 | 8260B | 8270C | | | 6010B | 9010B/ 9014 2540-G |
| Storm Inlet A | R6024 | 8260B | 8270C | | | 6010B | 9010B/ 9014 2540-G |
| Storm Inlet A | R6024MS | 8260B | 8270C | | | 6010B | 9010B/ 9014 |
| Storm Inlet A | R6024MSD | 8260B | 8270C | | | 6010B | 9010B/ 9014 |
| Storm Inlet A | R6024D | | | | | 6010B | 9010B/ 9014 2540-G |
| Outfall #1 | R6025 | 8260B | 8270C | | | 6010B | 9010B/ 9014 2540-G |
| Outfall #2 | R6026 | 8260B | 8270C | | | 6010B | 9010B/ 9014 2540-G |
| BLIND DUP | R6027 | 8260B | 8270C | | | 6010B | 9010B/ 9014 2540-G |
| QC Trip Blank | R6028 | 8260B | | | | | |
| Storage Blank | R6029 | 8260B | | | | | |

NEW YORK STATE DEPARTMENT OF ENVIRONMENTAL CONSERVATION

**SAMPLE PREPARATION AND ANALYSIS SUMMARY
INORGANIC ANALYSIS**

| Laboratory Sample ID | Matrix | Metals Requested | Date Rec'd at Lab | Date Analyzed |
|----------------------|--------|------------------|-------------------|---------------|
| R6023 | SOIL | TCL-LIST | 11/16/00 | 12/01/00 |
| R6024 | SOIL | TCL-LIST | 11/16/00 | 12/01/00 |
| R6024D | SOIL | TCL-LIST | 11/16/00 | 12/01/00 |
| R6024MS | SOIL | TCL-LIST | 11/16/00 | 12/01/00 |
| R6024MSD | SOIL | TCL-LIST | 11/16/00 | 12/01/00 |
| R6025 | SOIL | TCL-LIST | 11/16/00 | 12/01/00 |
| R6026 | SOIL | TCL-LIST | 11/16/00 | 12/01/00 |
| R6027 | SOIL | TCL-LIST | 11/16/00 | 12/01/00 |

NEW YORK STATE DEPARTMENT OF ENVIRONMENTAL CONSERVATION

SAMPLE PREPARATION AND ANALYSIS SUMMARY
INORGANIC ANALYSIS

| Laboratory Sample ID | Matrix | Metals Requested | Date Rec'd at Lab | Date Analyzed |
|----------------------|--------|------------------|-------------------|---------------|
| R6023 | SOIL | TCL-LIST | 11/16/00 | 12/01/00 |
| R6024 | SOIL | TCL-LIST | 11/16/00 | 12/01/00 |
| R6024D | SOIL | TCL-LIST | 11/16/00 | 12/01/00 |
| R6024MS | SOIL | TCL-LIST | 11/16/00 | 12/01/00 |
| R6024MSD | SOIL | TCL-LIST | 11/16/00 | 12/01/00 |
| R6025 | SOIL | TCL-LIST | 11/16/00 | 12/01/00 |
| R6026 | SOIL | TCL-LIST | 11/16/00 | 12/01/00 |
| R6027 | SOIL | TCL-LIST | 11/16/00 | 12/01/00 |

Project Management Case Narrative

INTRODUCTION/ANALYTICAL RESULTS

This report summarizes the laboratory results for samples from the Honeywell project located in Tonawanda, NY. New York State Department of Environmental Conservation forms are included in the Sample Data Package.

CONDITION UPON RECEIPT/CHAIN OF CUSTODY

The cooler(s) were received intact. When the cooler(s) were received by the laboratory, the sample custodian(s) opened and inspected the shipment(s) for damage, custody inconsistencies and proper preservation. Chain of custodies documenting receipt are presented in the chain of custody section. Each sample was assigned a unique laboratory number and a custody file created. The samples were placed in a secured walk-in cooler and signed in and out by the chemists performing the tests. The sign out record, or lab chronicle, is presented in the chain of custody section.

No discrepancies were noted upon receipt. Cooler temperatures were 4°C.

METHODOLOGY

The following methods were used to perform the analyses:

| PARAMETER | METHOD | REFERENCE |
|-----------------------|------------|-----------|
| Volatile Organics | 8260B | 1 |
| Semivolatile Organics | 8270C | 1 |
| ICP Metals | 6010B | 1 |
| Mercury | 7471A | 1 |
| Cyanide | 9010B/9014 | 1 |
| Percent Total Solids | 2540-G | 2 |

- 1) New York State Department of Environmental Conservation Analytical Services Protocol, October 1995.
- 2) Standard Methods for the Examination of Water and Wastewater, 18th Edition, 1992.

QUALITY CONTROL

QA/QC results are summarized in the Sample Data Package.

RAW DATA

The raw data is organized in the New York State Department of Environmental Conservation Analytical Services Protocol Category B order of data requirements.

GC/MS Volatile Organics Case Narrative

Client: O'Brien & Gere Engineers, Inc.
Job Number: 3435.061.076
Package #: 7435
Methodology: 8260B

Analyzed/Reviewed by (Date/Initials): SG 12-19-00

Supervisor/Reviewed by (Date/Initials): AD 12-19-00

QA/QC Review (Date/Initials): JAT 12/19/00

File Name in G/ Drive: A:\V7435.NAR

GC/MS Volatile Organics

The GC/MS Volatile instruments used a J&W DB-VRX, 60 m x 0.25 mm ID capillary column and a Vocabr 3000 trap.

Holding Times and Sample Preservation

All samples were prepared and analyzed within the method and/or QAPP specified holding time requirements. Water samples had a pH of less than 2.

Laboratory Control Sample

The following compound(s) did not meet laboratory control sample recovery criteria:

| LCS No. | Compound | Corrective Action |
|-----------|---------------|-------------------|
| L112500S3 | Chloromethane | 1 |
| | Chloroethane | 1 |
| | Bromomethane | 2 |

1. This compound exceeded the upper control limit and was not detected in the associated samples. No corrective action was taken.
2. The recovery was below the lower control limit. The compound was not detected in the associated samples. The analyte was not a primary target for the associated samples, which were dilutions or reanalyses for other target analytes.

MS/MSD/MSB

All spike recovery and RPD data met method and/or project specific QC criteria.

Surrogate

The following sample(s) did not meet surrogate recovery criteria:

| Sample Description | Sample # | Surrogate | Corrective Action |
|--------------------|----------|------------|-------------------|
| Storm Inlet A | R6024 | Toluene-d8 | 1 |

GC/MS Volatile Organics Case Narrative - Page 2

Client: O'Brien & Gere Engineers, Inc.
Job Number: 3435.061.076
Package #: 7435
Methodology: 8260B

1. The surrogate recovery was confirmed by MS/MSD analysis. No corrective action was taken.

Internal Standards

The internal standard area for the following sample(s) did not meet abundance criteria:

| Sample Description | Sample # | Internal Standard | Corrective Action |
|--------------------|----------|---------------------|-------------------|
| Storm Inlet A | R6024 | Fluorobenzene | 1 |
| | | Chlorobenzene | 1 |
| | | 1,4-Dichlorobenzene | 1 |
| Outfall #1 | R6025 | 1,4-Dichlorobenzene | 2 |

1. The sample, MS and MSD showed similar results. Matrix interference is suspected as the cause of the internal standard failure. No corrective action was taken.
2. The internal standard area met criteria in the reanalysis. No further corrective action was taken. Both results are reported due to calibration excursions associated with the second analysis.

Calibrations

The following continuing calibration compound(s) exceeded method percent drift and/or RRF criteria:

| Calibration Date | Instrument | Compound | %D | RRF | Corrective Action |
|------------------|------------|--------------|------|-----|-------------------|
| 11/25/00 | MS#3 | Bromomethane | 69.4 | | 1 |

1. Due to system performance the compound does not meet method/QAPP criteria. There were no positive hits for the compound in associated samples (that are reanalyses and dilution only). No corrective action was taken.

Preparation Blanks

All preparation blanks met method and/or project specific QC criteria.

Miscellaneous

[R6021] was reanalyzed due to suspected c/o of xylene and ethylbenzene from the previous sample. Both results are reported due to the calibration excursion associated with the second analysis.

[R6024MSD] was analyzed outside the 12 hrs tune clock by 1 minute. No corrective action was taken.

GC/MS Semivolatile Organics Case Narrative

Client: O'Brien & Gere Engineers, Inc.
Job Number: 3435.061.076
Package #: 7435
Methodology: 8270C

Analyzed/Reviewed by (Date/Initials): M.A 12-19-00

Supervisor/Reviewed by (Date/Initials): (A) 12-19-00

QA/QC Review (Date/Initials): J.M 12/19/00

File Name in G/ Drive: K:\WPWIN60\WPDOCS\7435SV.NAR

GC/MS Semivolatile Organics

The GC/MS Semivolatile instruments used a J&W DB-5MS, 30 m X 0.25 mm ID capillary column.

Holding Times

All samples were prepared and analyzed within the method and/or QAPP specified holding time requirements.

Laboratory Control Sample

The following compound(s) did not meet laboratory control sample recovery criteria:

| LCS No. | Compound | Corrective Action |
|-----------|----------------------|-------------------|
| L111700S1 | 4-Chloroaniline | 1 |
| | Benzo[a]pyrene | 2 |
| | Benzo[g,h,i]perylene | 3 |

1. The LCS was reanalyzed with similar results. The compound failing to meet acceptance criteria was inspected down to the MDL in the associated samples with no positive results detected. Results may be biased low. Both sets of data are included. No further corrective action was taken.
2. The LCS was reanalyzed with similar results. The compound was detected in the associated samples and results may be biased low. Both sets of data included. No further corrective action was taken.
3. A GPC problem affected recoveries for this compound. Results are biased low. No corrective action was taken.

MS/MSD/MSB

The following compound(s) did not meet matrix spike/matrix spike duplicate and/or matrix spike blank percent recovery and/or RPD criteria:

GC/MS Semivolatile Organics Case Narrative - Page 2

Client: O'Brien & Gere Engineers, Inc.
Job Number: 3435.061.076
Package: 7435
Methodology 8270C

| Sample Description | Sample # | Compound | % REC | RPD | Corrective Action |
|--------------------|------------|--------------------|-------|-----|-------------------|
| MSB01 | PS111700S1 | 2,4-Dinitrotoluene | X | | 1 |
| Storm Inlet A | R6024 | several | X | X | 2 |

1. The compound's recovery was above the UCL. No corrective action was taken.
2. The extract was diluted for high levels of matrix interference. Recoveries may have been affected. With the exception of 2,4-Dinitrotoluene, the associated matrix spike blank showed acceptable results. All failing compounds met percent recovery criteria in the associated LCS. No corrective action was taken.

Surrogates

The following sample(s) did not meet surrogate recovery criteria:

| Sample Description | Sample # | Surrogate | Corrective Action |
|--------------------|------------|------------------------|-------------------|
| MSB01 | PS111700S1 | 2-Chlorophenol-d4 | 1 |
| | | 1,2-Dichlorobenzene-d4 | 1 |

1. The control limits for surrogates 2-chlorophenol-d4 and/or 1,2-dichlorobenzene-d4 are advisory only. No corrective action is required.

Internal Standard Areas

The internal standard area for the following sample(s) did not meet abundance criteria:

| Sample Description | Sample # | Internal Standard | Corrective Action |
|--------------------|----------|-------------------|-------------------|
| Storm Inlet A | R6024 | Perylene-d12 | 1 |
| | R6024MSD | Perylene-d12 | 1 |

1. Although the MS met acceptance criteria, the sample, MS and MSD all contained high levels of matrix interference and showed similar results. No further corrective action was taken.

Calibrations

All calibrations and calibration verifications met method and/or project specific QC criteria.

Preparation Blanks

All preparation blanks met method and/or project specific QC criteria.

Miscellaneous

Elevated detection limits for sample extract Storm Inlet A[R6024] are due matrix interference.

GC/MS Semivolatile Organics Case Narrative - Page 3
Client: O'Brien & Gere Engineers, Inc.
Job Number: 3435.061.076
Package: 7435
Methodology 8270C

A GPC problem affected recoveries for Benzo[g,h,i]perylene and results are biased low for all samples.

The prep blank[PB111701S1] was lost during concentration of the extract. A second prep blank was created using the same solvent lot and surrogate.

Trace Metals Case Narrative

Client: O'Brien & Gere Engineers, Inc.
Job Number: 3435.061.076
Package #: 7435
Methodology: ICP metals- 6010B

Analyzed/Reviewed by (Date/Initials): MT 12-11-00

Supervisor/Reviewed by (Date/Initials): MT 12-11-00

QA/QC Review (Date/Initials): JH 12/13/00

File Name in G/ Drive: G:\NARRATIV\7435OBGE.ICP

Trace Metals

Holding Times

All samples were prepared and analyzed within the method and/or QAPP specified holding time requirements.

Laboratory Control Sample

All spike recoveries met method and/or project specific QC criteria.

MS/MSD AND MS/MSD RPD

The following analytes did not meet matrix spike/matrix spike duplicate percent recovery and/or MS/MSD RPD criteria:

| Sample Description | Sample # | Analyte | % REC | RPD | Corrective Action |
|--------------------|----------|-----------|-------|-----|-------------------|
| Storm Inlet A | R6024 | Aluminum | X | | 1 |
| | | Antimony | X | | 1 |
| | | Iron | X | | 1 |
| | | Lead | X | X | 1 |
| | | Manganese | X | | 1 |
| | | Zinc | X | | 1 |

1. A post-digestion spike was performed as required. No further corrective action was taken.

Sample Duplicate

The following analyte did not meet sample duplicate RPD criteria:

| Sample Description | Sample # | Analyte | RPD | Corrective Action |
|--------------------|----------|---------|-----|-------------------|
| Storm Inlet A | R6024 | Calcium | X | 1 |

1. The failing RPD is attributed to matrix complexity. No corrective action was taken.

Trace Metals Case Narrative - Page 2

Client:

O'Brien & Gere Engineers, Inc.

Job Number:

3435.061.076

Package #:

7435

Methodology:

ICP metals- 6010B

ICP Serial Dilution

The following analyte did not meet ICP serial dilution recovery criteria:

| Sample Description | Sample # | Analyte | Corrective Action |
|--------------------|----------|---------|-------------------|
| Storm Inlet A | R6024 | Sodium | 1 |

1. No corrective action was required.

Calibrations

All calibrations and calibration verifications met method and/or project specific QC criteria.

Preparation Blanks

All preparation blanks met method and/or project specific QC criteria.

Trace Metals Case Narrative

Client: O'Brien & Gere Engineers, Inc.
Job Number: 3435.061.076
Package #: 7435
Methodology: Mercury - 7471A

Analyzed/Reviewed by (Date/Initials): 12-11-00 my

Supervisor/Reviewed by (Date/Initials): 12-11-00 m

QA/QC Review (Date/Initials): JM 12/13/00

File Name in G/ Drive: G:\NARRATIV\7435OBGE.HG

Trace Metals

Holding Times

All samples were prepared and analyzed within the method and/or QAPP specified holding time requirements.

Laboratory Control Sample

All spike recoveries met method and/or project specific QC criteria.

MS/MSD AND MS/MSD RPD

The following analyte did not meet matrix spike/matrix spike duplicate percent recovery and/or MS/MSD RPD criteria:

| Sample Description | Sample # | Analyte | % REC | RPD | Corrective Action |
|--------------------|----------|---------|-------|-----|-------------------|
| Storm Inlet A | R6024 | Mercury | X | X | 1 |

1. MS/MSD recoveries were reasonable, suspect matrix complexity. No corrective action was taken.

Sample Duplicate

All sample duplicate RPD data met method and/or project specific QC criteria.

Calibrations

All calibrations and calibration verifications met method and/or project specific QC criteria.

Preparation Blanks

All preparation blanks met method and/or project specific QC criteria.

Wet Chemistry Case Narrative

Client: O'Brien & Gere Engineers
Job Number: 3435.61.76
Package #: 7435
Methodology: %Total Solids - 2540-G
Total cyanide-CLP-B - 9010B/9014

Analyzed/Reviewed by (Date/Initials): MS 12/8/02

Supervisor/Reviewed by (Date/Initials): MT 12/8/02

QA/QC Review (Date/Initials): JH 12/13/02

File Name in G/ Drive: G:\NARRATIV\7435OBGE.WC

Wet Chemistry

There were no excursions to note. All QC results were within established control limits.

GROUND WATER SAMPLES

ANALYTICAL PACKAGE

for

**Honeywell
Tonawanda, NY
Water Samples**

Water samples collected: June 5, 6, 7, and 9, 2001

Volume 1 of 4

Prepared for:

O'Brien & Gere Engineers, Inc.
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Prepared by:

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Date

Thomas J. Wilson
7.11.01

Reviewed

Date

Joseph C. Harr
7/12/01

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Sample Data Summary Package

NEW YORK STATE DEPARTMENT OF ENVIRONMENTAL CONSERVATION

**SAMPLE IDENTIFICATION AND
ANALYTICAL SUMMARY**

| Customer Sample Code | Laboratory Sample Code | Analytical Requirements | | | | | |
|----------------------|------------------------|-------------------------|---------------------|------------------|---------------------|-------------|--------|
| | | 'VOA GC/MS Method # | 'BNA GC/MS Method # | 'VOA GC Method # | 'Pest PCBs Method # | 'Metals | 'Other |
| MW-10 | S6749 | 8260B | 8270C | | | 6010B/7470A | 335.2 |
| MW-8 | S6750 | 8260B | 8270C | | | 6010B/7470A | 335.2 |
| MW-1 | S6751 | 8260B | 8270C | | | 6010B/7470A | 335.2 |
| MW-7 | S6752 | 8260B | 8270C | | | 6010B/7470A | 335.2 |
| MW-7 | S6752MS | 8260B | 8270C | | | 6010B/7470A | 335.2 |
| MW-7 | S6752MSD | 8260B | 8270C | | | 6010B/7470A | 335.2 |
| MW-7 | S6752D | | | | | 6010B/7470A | 335.2 |
| MW-6 | S6753 | 8260B | 8270C | | | 6010B/7470A | 335.2 |
| MW-9 | S6754 | 8260B | 8270C | | | 6010B/7470A | 335.2 |
| X-1 | S6755 | 8260B | 8270C | | | 6010B/7470A | 335.2 |
| QC Trip Blank | S6756 | 8260B | | | | | |
| MW-11 | S6817 | 8260B | 8270C | | | 6010B/7470A | 335.2 |
| MW-5 | S6818 | 8260B | 8270C | | | 6010B/7470A | 335.2 |
| QC Trip Blank | S6819 | 8260B | | | | | |
| Storage Blank | S6820 | 8260B | | | | | |
| MW-2 | S6833 | 8260B | 8270C | | | 6010B/7470A | 335.2 |

NEW YORK STATE DEPARTMENT OF ENVIRONMENTAL CONSERVATION

**SAMPLE IDENTIFICATION AND
ANALYTICAL SUMMARY**

| Customer Sample Code | Laboratory Sample Code | Analytical Requirements | | | | | |
|----------------------------|------------------------------|------------------------------|------------------------------|---------------------------|------------------------------|-----------------|--------|
| | | 'VOA GC/MS Method # | 'BNA GC/MS Method # | 'VOA GC Method # | 'Pest PCBs Method # | 'Metals | 'Other |
| MW-4 | S6834 | 8260B | 8270C | | | 6010B/ 7470A | 335.2 |
| MW-3 | S6835 | 8260B | 8270C | | | 6010B/ 7470A | 335.2 |
| QC Trip Blank | S6836 | 8260B | | | | | |

NEW YORK STATE DEPARTMENT OF ENVIRONMENTAL CONSERVATION

**SAMPLE PREPARATION AND ANALYSIS SUMMARY
SEMIVOLATILE(BNA) ANALYSES**

| LABORATORY SAMPLE ID | MATRIX | DATE COLLECTED | DATE REC'D at LAB | DATE EXTRACTED | DATE ANALYZED |
|-------------------------|--------|-------------------|----------------------|-------------------|------------------|
| S6749 | Water | 6/5/01 | 6/7/01 | 6/11/01 | 6/18/01 |
| S6750 | Water | 6/5/01 | 6/7/01 | 6/11/01 | 6/18/01 |
| S6751 | Water | 6/5/01 | 6/7/01 | 6/11/01 | 6/18,19/01 |
| S6752 | Water | 6/6/01 | 6/7/01 | 6/11/01 | 6/18,19/01 |
| S6752MS | Water | 6/6/01 | 6/7/01 | 6/11/01 | 6/18/01 |
| S6752MSD | Water | 6/6/01 | 6/7/01 | 6/11/01 | 6/18/01 |
| S6753 | Water | 6/6/01 | 6/7/01 | 6/11/01 | 6/19/01 |
| S6754 | Water | 6/6/01 | 6/7/01 | 6/11/01 | 6/20/01 |
| S6755 | Water | 6/6/01 | 6/7/01 | 6/11/01 | 6/19/01 |
| S6817 | Water | 6/7/01 | 6/9/01 | 6/13/01 | 6/19/01 |
| S6818 | Water | 6/7/01 | 6/9/01 | 6/13/01 | 6/19/01 |
| S6833 | Water | 6/9/01 | 6/11/01 | 6/13/01 | 6/19/01 |
| S6834 | Water | 6/9/01 | 6/11/01 | 6/13/01 | 6/19/01 |
| S6835 | Water | 6/9/01 | 6/11/01 | 6/13/01 | 6/19/01 |
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NEW YORK STATE DEPARTMENT OF ENVIRONMENTAL CONSERVATION

**SAMPLE PREPARATION AND ANALYSIS SUMMARY
INORGANIC ANALYSIS**

| Laboratory Sample ID | Matrix | Metals Requested | Date Rec'd at Lab | Date Analyzed |
|----------------------|--------|------------------|-------------------|-------------------|
| S6749 | Water | TCL Metals | 06/07/01 | 06/20/01,07/02/01 |
| S6750 | Water | TCL Metals | 06/07/01 | 06/20/01,07/02/01 |
| S6751 | Water | TCL Metals | 06/07/01 | 06/20/01,07/02/01 |
| S6752 | Water | TCL Metals | 06/07/01 | 06/20/01,07/02/01 |
| S6752D | Water | TCL Metals | 06/07/01 | 06/20/01,07/02/01 |
| S6752MS | Water | TCL Metals | 06/07/01 | 06/20/01,07/02/01 |
| S6752MSD | Water | TCL Metals | 06/07/01 | 06/20/01,07/02/01 |
| S6753 | Water | TCL Metals | 06/07/01 | 06/20/01,07/02/01 |
| S6754 | Water | TCL Metals | 06/07/01 | 06/20/01,07/02/01 |
| S6755 | Water | TCL Metals | 06/07/01 | 06/20/01,07/02/01 |
| S6817 | Water | TCL Metals | 06/09/01 | 06/20/01,07/02/01 |
| S6818 | Water | TCL Metals | 06/09/01 | 06/20/01,07/02/01 |
| S6833 | Water | TCL Metals | 06/11/01 | 06/20/01,07/02/01 |
| S6834 | Water | TCL Metals | 06/11/01 | 06/20/01,07/02/01 |
| S6835 | Water | TCL Metals | 06/11/01 | 06/20/01,07/02/01 |

Project Management Case Narrative

INTRODUCTION/ANALYTICAL RESULTS

This report summarizes the laboratory results for samples from the Honeywell project located in Tonawanda, NY. New York State Department of Environmental Conservation forms are included in the Sample Data Package.

CONDITION UPON RECEIPT/CHAIN OF CUSTODY

The cooler(s) were received intact. When the cooler(s) were received by the laboratory, the sample custodian(s) opened and inspected the shipment(s) for damage, custody inconsistencies and proper preservation. Chain of custodies documenting receipt are presented in the chain of custody section. Each sample was assigned a unique laboratory number and a custody file created. The samples were placed in a secured walk-in cooler and signed in and out by the chemists performing the tests. The sign out record, or lab chronicle, is presented in the chain of custody section.

Discrepancies noted upon receipt are documented on the case file forms included in the chain of custody section. Cooler temperatures were 3°C, 4°C, and 5°C.

METHODOLOGY

The following methods were used to perform the analyses:

| PARAMETER | METHOD | REFERENCE |
|-----------------------|------------|-----------|
| Volatile Organics | 8260B | 1 |
| Semivolatile Organics | 8270C | 1 |
| ICP Metals | 6010B | 1 |
| Mercury | 7471A | 1 |
| Cyanide | 9010B/9014 | 1 |

- 1) New York State Department of Environmental Conservation Analytical Services Protocol, October 1995.

QUALITY CONTROL

QA/QC results are summarized in the Sample Data Package.

RAW DATA

The raw data is organized in the New York State Department of Environmental Conservation Analytical Services Protocol Category B order of data requirements.

GC/MS Volatile Organics Case Narrative

Client: O'Brien & Gere Engineers, Inc.
Job Number: 3435.061.076
Package: 9162, 9180, 9181, 9186
Methodology: 8260B

Analyzed/Reviewed by (Date/Initials): AJL 7-5-01

Supervisor/Reviewed by (Date/Initials): (A) 7-5-01

QA/QC Review (Date/Initials): M 7/6/01

File Name in G/ Drive: C:\MY DOCUMENTS\9162V.NAR.DOC

GC/MS Volatile Organics

The GC/MS Volatile instruments used a J&W DB-VRX, 75 m x 0.45 mm ID capillary column and a Vocarb 3000 trap.

Holding Times and Sample Preservation

All samples were prepared and analyzed within the method and/or QAPP specified holding time requirements. Samples had a pH of less than 2.

Laboratory Control Sample

All spike recoveries met method and/or project specific QC criteria.

MS/MSD

The following compound(s) did not meet matrix spike/matrix spike duplicate percent recovery and/or RPD criteria:

| Sample Description | Sample # | Compound | % REC | RPD | Corrective Action |
|--------------------|------------|--------------------|-------|-----|-------------------|
| MSB01 | PS061201W2 | Bromomethane | X | | 1 |
| MW-7 | S6752 | Bromomethane | X | X | 1 |
| | | Styrene | X | | 2 |
| | | Methylene chloride | | X | 3 |

1. This compound exceeded the upper control limit and was not detected in associated samples. The recovery met acceptance criteria in the LCS. No corrective action was taken.
2. The percent recovery was marginally below the lower control limit and was not detected in associated samples. The LCS met acceptance criteria for this compound. No corrective action was taken.
3. The RPD was marginally above the upper control limit. The compound met acceptance criteria in the LCS. No corrective action was taken.

GC/MS Volatile Organics Case Narrative - page 2

Client: O'Brien & Gere Engineers, Inc.
Job Number: 3435.061.076
Package: 9162, 9180, 9181, 9186
Methodology: 8260B

Surrogate

The following sample(s) did not meet surrogate recovery criteria:

| Sample Description | Sample # | Surrogate | Corrective Action |
|--------------------|-----------|------------|-------------------|
| LCS02 | L061301W2 | Toluene-d8 | 1 |

1. This surrogate was marginally above the upper control limit. All surrogates met criteria in the duplicate LCS as well as all associated samples. No corrective action was taken.

Internal Standards

All internal standard areas met method and/or project specific QC criteria.

Calibrations

The following continuing calibration compound(s) exceeded method percent drift and/or RRF criteria:

| Calibration Date | Instrument | Compound | %D | RRF | Corrective Action |
|------------------|------------|---------------------------|-------|-----|-------------------|
| 6/12/01 | MS#2 | 1,1,2,2-Tetrachloroethane | -36.9 | | 1 |

1. The method allows two compounds to fail as long as their percent difference is less than 40%. No corrective action was taken.

Preparation Blanks

All preparation blanks met method and/or project specific QC criteria.

Miscellaneous

Some samples were diluted due to high concentrations of non-target compounds. PQL's are elevated for these samples.

GC/MS Semivolatile Organics Case Narrative

Client: O'Brien & Gere Engineers, Inc.
Job Number: 3435.061.076
Package #: 9162, 9180, 9186
Methodology: 8270C

Analyzed/Reviewed by (Date/Initials): MAD for DTP 7-5-01

Supervisor/Reviewed by (Date/Initials): (@)TGS-01

QA/QC Review (Date/Initials): JW 7/9/01

File Name in G/ Drive: Q:\My Documents\9162sv.nar.doc

GC/MS Semivolatile Organics

The GC/MS Semivolatile instruments used a J&W DB-5MS, 30 m X 0.25 mm ID capillary column.

Holding Times

All samples were prepared and analyzed within the method and/or QAPP specified holding time requirements. Sample MW-2 [S6833] and samples extracted on 6/11/01 were re-extracted outside holding time due to QA/QC excursions and will be reported in a future SDG.

Laboratory Control Sample

The following compound(s) did not meet laboratory control sample recovery criteria:

| LCS No. | Compound | Corrective Action |
|-----------|------------------------|-------------------|
| L061101W2 | Numerous | 1 |
| L061301W2 | Hexachlorobutadiene | 2 |
| | N-Nitrosodiphenylamine | 2 |
| D061301W2 | Hexachlorobutadiene | 2 |
| | N-Nitrosodiphenylamine | 2 |

1. The LCS was reanalyzed and still did not meet criteria. All samples associated with this LCS will be re-extracted over the holding time. The re-extracts will be reported in a future SDG.
2. The recovery was marginally below the lower control limit in the primary and duplicate LCS. Associated samples were inspected for concentrations down to the MDL and none were detected. No further corrective action was taken.

MS/MSD/MSB

The following compound(s) did not meet matrix spike/matrix spike duplicate percent recovery and/or RPD criteria:

GC/MS Semivolatile Organics Case Narrative -page 2

Client: O'Brien & Gere Engineers, Inc.
 Job Number: 3435.061.076
 Package #: 9162, 9180, 9186
 Methodology: 8270C

| Sample Description | Sample # | Compound | % REC | RPD | Corrective Action |
|--------------------|------------|-------------------------|-------|-----|-------------------|
| MSB01 | PS061101W2 | 4-Chloro-3-methylphenol | X | | 1 |
| | | Pentachlorophenol | X | | 1 |

1. The recoveries were above the upper control limit and were not detected in the associated samples. No corrective action was taken.

Surrogates

The following sample(s) did not meet surrogate recovery criteria:

| Sample Description | Sample # | Surrogate | Corrective Action |
|--------------------|----------|------------------|-------------------|
| MW-8 | S6750 | 2-Fluorobiphenyl | 1 |
| MW-2 | S6833 | Several | 2 |

1. Two of the three base/neutral and/or acid extractable surrogate recoveries passed acceptance criteria. Laboratory or QAPP SOP requires no corrective action.
2. Sample will be re-extracted over the holding time and results will be reported in a future SDG.

Internal Standard Areas

All internal standard areas met method and/or project specific QC criteria.

Calibrations

The following continuing calibration compound(s) exceeded method percent drift and/or RRF criteria:

| Calibration Date | Instrument | Compound | %D | RRF | Corrective Action |
|------------------|------------|------------------------|-------|-----|-------------------|
| 6/19/01 | MS#5 | Indeno[1,2,3-cd]pyrene | -33.6 | | 1 |
| | | Dibenz[a,h]anthracene | -34.4 | | 1 |
| | | Benzo[g,h,i]perylene | -39.3 | | 1 |

1. The method allows four compounds to fail as long as their percent difference is less than 40%.

Preparation Blanks

All preparation blanks met method and/or project specific QC criteria.

Trace Metals Case Narrative

Client:

O'Brien & Gere Engineers, Inc.

Job Number:

3435.061.076

Package #:

9162,9180,9186

Methodology:

ICP metals - 6010B

Analyzed/Reviewed by (Date/Initials):

17-02-01 CT

Supervisor/Reviewed by (Date/Initials):

7-2-01 mwt

QA/QC Review (Date/Initials):

1/3/01

File Name in G/ Drive:

G:\NARRATIV\9162OBGEng.icp.wpd

Trace Metals

Holding Times

All samples were prepared and analyzed within the method and/or QAPP specified holding time requirements.

Laboratory Control Sample

All spike recoveries met method and/or project specific QC criteria.

MS/MSD AND MS/MSD RPD

The following analytes did not meet matrix spike/matrix spike duplicate percent recovery criteria:

| Sample Description | Sample # | Analyte | % REC | Corrective Action |
|--------------------|----------|----------|-------|-------------------|
| MW-7 | S6752 | Aluminum | X | 2 |
| | | Iron | X | 2 |
| | | Lead | X | 2 |
| | | Thallium | X | 1 |

1. A post-digestion spike was performed as required. No further corrective action was taken.
2. The concentration of the analyte in the sample was four times greater than the concentration of the spike added. No corrective action was required.

Sample Duplicate

All sample duplicate RPD data met method and/or project specific QC criteria.

ICP Serial Dilution

All percent differences met method and/or project specific QC criteria.

Trace Metals Case Narrative - Page 2

Client: O'Brien & Gere Engineers, Inc.
Job Number: 3435.061.076
Package #: 9162,9180,9186
Methodology: ICP metals - 6010B

Calibrations

All calibrations and calibration verifications met method and/or project specific QC criteria.

Preparation Blanks

All preparation blanks met method and/or project specific QC criteria.

Trace Metals Case Narrative

Client: O'Brien & Gere Engineers, Inc.
Job Number: 3435.061.076
Package #: 9162,9180,9186
Methodology: Mercury - 7470A

Analyzed/Reviewed by (Date/Initials): MT 6-29-01

Supervisor/Reviewed by (Date/Initials): mt 6-29-01

QA/QC Review (Date/Initials): JW 6/20/01

File Name in G/ Drive: G:\NARRATIV\9162obge.hg.wpd

Trace Metals

There were no excursions to note. All QC results were within established control limits.

Wet Chemistry Case Narrative

Client: O'Brien & Gere Engineers, Inc.
Job Number: 3435.061.076
Package #: 9162,9180,9186
Methodology: Total cyanide - EPA - 335.2

Analyzed/Reviewed by (Date/Initials): 06/21/01 ZK

Supervisor/Reviewed by (Date/Initials): 6/21/01 NY

QA/QC Review (Date/Initials): JH 6/26/01

File Name in G/ Drive: G:\NARRATIV\9162OBG.WC

Wet Chemistry

Holding Times

All samples were prepared and analyzed within the method and/or QAPP specified holding times.

Laboratory Control Sample

All spike recoveries met method and/or project specified QC criteria.

MS/MSD AND MS/MSD RPD

The following analyte did not meet matrix spike/matrix spike duplicate percent recovery criteria:

| Sample Description | Sample # | Analyte | % REC | RPD | Corrective Action |
|--------------------|----------|---------------|-------|-----|-------------------|
| MW-7 | S6752 | Total cyanide | X | | 1 |

1. The concentration of the analyte in the sample was much greater than the concentration of the spike added. No further corrective action was taken.

Sample Duplicate

All sample duplicate RPD data met method and/or project specific QC criteria.

Calibrations

All calibrations and calibration verifications met method and/or project specific QC criteria.

Preparation Blanks

All preparation blanks met method and/or project specific QC criteria.

STORM WATER SAMPLES

ANALYTICAL PACKAGE

for

**O'Brien & Gere Engineers, Inc.
Honeywell - Tonawanda Coke
Tonawanda, NY
Water Samples**

Water samples collected: December 11, 2001

Volume 1 of 3

Prepared for: O'Brien & Gere Engineers, Inc.
5000 Brittonfield Parkway
P.O. Box 4873
Syracuse, NY 13221

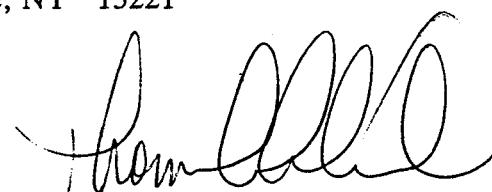
Prepared by: O'Brien & Gere Laboratories, Inc.
5000 Brittonfield Parkway
Suite 300, P.O. Box 4942
Syracuse, NY 13221

Authorized

Date

Reviewed

Date


Tom Hause

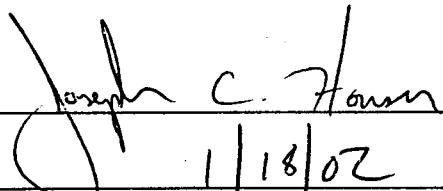
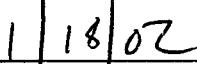
11/18/02

Joseph C. Hause

11/18/02

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Sample Data Summary Package

NEW YORK STATE DEPARTMENT OF ENVIRONMENTAL CONSERVATION

**SAMPLE IDENTIFICATION AND
ANALYTICAL SUMMARY**

| Customer Sample Code | Laboratory Sample Code | Analytical Requirements | | | | | |
|----------------------|------------------------|-------------------------|--------------------|-----------------|--------------------|-----------------------------|--------------|
| | | VOA GC/MS Method # | BNA GC/MS Method # | VOA GC Method # | Pest PCBs Method # | Metals | Other |
| Inlet A | T6812 | 95-1 | 95-2 | | | 200.7 CLP-M* & 245.1 CLP-M* | 335.2 CLP-M* |
| Inlet B | T6813 | 95-1 | 95-2 | | | 200.7 CLP-M* & 245.1 CLP-M* | 335.2 CLP-M* |
| Inlet B | T6813MS | 95-1 | 95-2 | | | 200.7 CLP-M* & 245.1 CLP-M* | 335.2 CLP-M* |
| Inlet B | T6813MSD | 95-1 | 95-2 | | | | |
| Inlet B | T6813D | | | | | 200.7 CLP-M* & 245.1 CLP-M* | 335.2 CLP-M* |
| Outfall 2 | T6814 | 95-1 | 95-2 | | | 200.7 CLP-M* & 245.1 CLP-M* | 335.2 CLP-M* |
| Outfall 1 | T6815 | 95-1 | 95-2 | | | 200.7 CLP-M* & 245.1 CLP-M* | 335.2 CLP-M* |
| X-1 | T6816 | 95-1 | 95-2 | | | 200.7 CLP-M* & 245.1 CLP-M* | 335.2 CLP-M* |
| Storage Blank | T6817 | 95-1 | | | | | |
| QC Trip Blank | T6818 | 95-1 | | | | | |

NEW YORK STATE DEPARTMENT OF ENVIRONMENTAL CONSERVATION

SAMPLE PREPARATION AND ANALYSIS SUMMARY
VOLATILE(VOA) ANALYSES

| LABORATORY SAMPLE ID | MATRIX | DATE COLLECTED | DATE REC'D at LAB | LOW LEVEL MED LEVEL | DATE ANALYZED |
|-------------------------|--------|-------------------|----------------------|------------------------|------------------|
| T6812 | Water | 12/11/01 | 12/12/01 | Low | 12/13/01 |
| T6813 | Water | 12/11/01 | 12/12/01 | Low | 12/13/01 |
| T6813MS | Water | 12/11/01 | 12/12/01 | Low | 12/13/01 |
| T6813MSD | Water | 12/11/01 | 12/12/01 | Low | 12/13/01 |
| T6814 | Water | 12/11/01 | 12/12/01 | Low | 12/13/01 |
| T6815 | Water | 12/11/01 | 12/12/01 | Low | 12/13/01 |
| T6816 | Water | 12/11/01 | 12/12/01 | Low | 12/13/01 |
| T6817 | Water | 12/12/01 | 12/12/01 | Low | 12/13/01 |
| T6818 | Water | 12/11/01 | 12/12/01 | Low | 12/14/01 |
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NEW YORK STATE DEPARTMENT OF ENVIRONMENTAL CONSERVATION

**SAMPLE PREPARATION AND ANALYSIS SUMMARY
INORGANIC ANALYSIS**

| Laboratory Sample ID | Matrix | Metals Requested | Date Rec'd at Lab | Date Analyzed |
|----------------------|--------|------------------|-------------------|------------------------|
| T6812 | WATER | CLP | 12/12/01 | 12/15/01 & 01/02,13/02 |
| T6813 | WATER | CLP | 12/12/01 | 12/15/01 & 01/02,13/02 |
| T6813D | WATER | CLP | 12/12/01 | 12/15/01 & 01/02,13/02 |
| T6813MS | WATER | CLP | 12/12/01 | 12/15/01 & 01/02,13/02 |
| T6814 | WATER | CLP | 12/12/01 | 12/15/01 & 01/02,13/02 |
| T6815 | WATER | CLP | 12/12/01 | 12/15/01 & 01/02,13/02 |
| T6816 | WATER | CLP | 12/12/01 | 12/15/01 & 01/02,13/02 |

Project Management Case Narrative

INTRODUCTION/ANALYTICAL RESULTS

This report summarizes the laboratory results for samples from Honeywell, Tonawanda Coke located in Tonawanda, NY. New York State Department of Environmental Conservation forms are included in the Sample Data Summary Package and in the Sample Data Package.

CONDITION UPON RECEIPT/CHAIN OF CUSTODY

The cooler(s) were received intact. When the cooler(s) were received by the laboratory, the sample custodian(s) opened and inspected the shipment(s) for damage, custody inconsistencies and proper preservation. Chain of custodies documenting receipt are presented in the chain of custody section. Each sample was assigned a unique laboratory number and a custody file created. The samples were placed in a secured walk-in cooler and signed in and out by the chemists performing the tests. The sign out record, or lab chronicle, is presented in the chain of custody section.

No discrepancies were noted upon receipt. Cooler temperatures were 3°C.

METHODOLOGY

The following methods were used to perform the analyses:

| PARAMETER | METHOD | REFERENCE |
|-----------------------|--------------|-----------|
| Volatile Organics | 95-1 | 1 |
| Semivolatile Organics | 95-2 | 1 |
| ICP Metals | 200.7 CLP-M* | 1 |
| Mercury | 245.1 CLP-M* | 1 |
| Cyanide | 335.2 CLP-M* | 1 |

- 1) New York State Department of Environmental Conservation Analytical Services Protocol, October 1995.

QUALITY CONTROL

QA/QC results are summarized in the Sample Data Summary Package and are also included in the raw data.

RAW DATA

The raw data is organized in the New York State Department of Environmental Conservation Analytical Services Protocol Superfund order of data requirements.

GC/MS Volatile Organics Case Narrative

Client: O' Brien & Gere Engineers, Inc.
Job Number: 3435.044.76391
Package #: 725
Methodology: ASP 95-1

Analyzed/Reviewed by (Date/Initials): SG 12-31-01

Supervisor/Reviewed by (Date/Initials): (2) 1-4-02

QA/QC Review (Date/Initials): 1/6/02

GC/MS Volatile Organics

The GC/MS Volatile instruments used a Restek Rtx-502.2, 60 m x 0.25 mm ID capillary column and a Vocarb 3000 trap.

There were no excursions to note. All QC results were within established control limits.

Holding Times and Sample Preservation

All samples were prepared and analyzed within the method and/or QAPP specified holding time requirements. Samples had a pH of less than 2.

Laboratory Control Sample

All spike recoveries met method and/or project specific QC criteria.

MS/MSD/MSB

All spike recovery and RPD data met method and/or project specific QC criteria.

Surrogate Standards

All surrogate recoveries met method and/or project specific QC criteria.

Internal Standards

All internal standard areas met method and/or project specific QC criteria.

Calibrations

All calibrations and calibration verifications met method and/or project specific QC criteria.

Preparation Blanks

All preparation blanks met method and/or project specific QC criteria.

GC/MS Semivolatile Organics Case Narrative

Client: O'Brien & Gere Engineers, Inc.
Job Number: 3435.044.76391
Package #: 725
Methodology: 95-2

Analyzed/Reviewed by (Date/Initials): ASL 1-8-02

Supervisor/Reviewed by (Date/Initials): (AU) 1-8-02

QA/QC Review (Date/Initials): JH 1/9/02

File Name in G/ Drive: C:\My Documents\725sv.nar.doc

GC/MS Semivolatile Organics

The GC/MS Semivolatile instruments used a J&W DB-5MS, 30 m X 0.25 mm ID capillary column.

Holding Times

All samples were prepared and analyzed within the method and/or QAPP specified holding time requirements.

Laboratory Control Sample

The following compound(s) did not meet laboratory control sample recovery criteria:

| LCS No. | Compound | Corrective Action |
|-----------|------------------------|-------------------|
| L121301W2 | 4-Nitroaniline | 1 |
| | 3,3'-Dichlorobenzidine | 1 |

1. The LCS was reanalyzed and still did not meet criteria. Both sets of data are included. The analyte was not detected in associated samples. No further corrective action was taken.

MS/MSD/MSB

The following compound(s) did not meet matrix spike/matrix spike duplicate/matrix spike blank percent recovery and/or RPD criteria:

| Sample Description | Sample # | Compound | % REC | RPD | Corrective Action |
|--------------------|------------|--------------------|-------|-----|-------------------|
| MSB01 | PS121301W2 | 4-Nitrophenol | X | | 1 |
| Inlet B | T6813 | 4-Nitrophenol | X | | 1 |
| | | Pentachlorophenol | X | | 1 |
| | | 2,4-Dinitrotoluene | X | | 1 |

1. The associate LCS met acceptance criteria. No corrective action was taken.

GC/MS Semivolatile Organics Case Narrative - Page 2

Client: O'Brien & Gere Engineers, Inc.
Job Number: 3435.044.76391
Package #: 725
Methodology: 95-2

Surrogates

All surrogate recoveries met method and/or project specific QC criteria.

Internal Standard Areas

All internal standard areas met method and/or project specific QC criteria.

Calibrations

The following continuing calibration compound(s) exceeded method percent drift and/or RRF criteria:

| Calibration Date | Instrument | Compound | %D | RRF | Corrective Action |
|------------------|------------|---------------------------|----|-----|-------------------|
| 12/15/01 | MS#5 | 2-Methylphenol | X | | 1 |
| 12/19/01 | MS#5 | 2-Methylphenol | X | | 1 |
| | | N-Nitrosodi-n-propylamine | X | | 1 |
| | | 2-Fluorophenol | X | | 1 |

1. The method allows four compounds to fail as long as their percent difference is less than 40%. No corrective action was taken.

Preparation Blanks

All preparation blanks met method and/or project specific QC criteria.

Trace Metals Case Narrative

Client: O'Brien & Gere Engineers, Inc.
Job Number: 3435.044.76391
Package #: 725
Methodology: ICP metals - 200.7 CLP-M*
Mercury - 245.1 CLP-M*
Total cyanide - 335.2 CLP-M*

Analyzed/Reviewed by (Date/Initials): 1-15-02 mrt

Supervisor/Reviewed by (Date/Initials): 1-15-02 mrt

QA/QC Review (Date/Initials): JM 1/17/02

File Name in G/ Drive: G:\NARRATIV\725OBGEcoke.clp.wpd

Trace Metals

Holding Times

All samples were prepared and analyzed within the method and/or QAPP specified holding time requirements.

Laboratory Control Sample

All spike recoveries met method and/or project specific QC criteria.

Matrix Spike

All spike recovery and RPD data met method and/or project specific QC criteria.

Sample Duplicate

All sample duplicate RPD data met method and/or project specific QC criteria.

ICP Serial Dilution

The following analyte did not meet ICP serial dilution recovery criteria:

| Sample Description | Sample # | Analyte | Corrective Action |
|--------------------|----------|-----------|-------------------|
| Inlet B | T6813 | Potassium | 1 |

- Form I's were flagged with an "E" accordingly.

Calibrations

All calibrations and calibration verifications met method and/or project specific QC criteria.

Preparation Blanks

All preparation blanks met method and/or project specific QC criteria.

Appendix B

Test Boring Logs

| OBRIEN & GERE ENGINEERS, INC. | | | | | | TEST BORING LOG | REPORT OF BORING MW-12 | | | |
|---|-----|--------------|-----------|------------------|-----------|---|--|------------------|-------------------------|---------------------------|
| Client: Honeywell | | | | | | Sampler: 2-inch Split Spoon | Page 1 of Location: | | | |
| Proj. Loc: Tonawanda, NY | | | | | | Hammer: 140-lb | Start Date: 5/24/01 End Date: 5/24/01 | | | |
| File No.: 1163/26080 | | | | | | Fall: 30-inch | Screen Riser | = | \ | Grout Sand Pack Bentonite |
| Boring Company: SJB Services Foreman: Matt Mattehis OBG Geologist: DJ Carnevale | | | | | | | | | | |
| Depth Below Grade | No. | Depth (feet) | Blows /6" | Penetr/ Recovery | "N" Value | Sample Description | Stratum Change General Descript | Equip. Installed | Field Testing PID (ppm) | UV |
| 0 | | 2 | 2-3 | 2/2 | 5 | Brownish Gray (5YR 4/1), damp, SILT, some gravel fill | | | 14.4 | |
| | | | 2-2 | | | | | | | |
| 2 | | 4 | 4-6 | 2/2 | 10 | Moderate reddish brown (10R 4/6), damp SILT, some gravel fill | | | 8.3 | |
| | | | 4-4 | | | | | | | |
| 4 | | 6 | 2-1 | 2/2 | 4 | Black (N1), moist, gravelly FILL to ~5 ft, to moist, medium gray (N5), clayey SILT to 6 ft | | | 11.4 | |
| | | | 3-3 | | | | | | | |
| 6 | | 8 | 2-3 | 2/2 | 8 | Medium gray (N5), moist, clayey SILT to ~7.5 ft, moderate reddish brown (10R 4/6), damp SILT, little clay to 8 ft | | | 5.4 | |
| | | | 5-5 | | | | | | | |
| 8 | | 10 | 2-5 | 2/2 | 14 | Moderate reddish brown (10R 4/6), moist SILT, little clay to 10 ft | | | 8.2 | |
| | | | 9-12 | | | | | | | |
| Well construction: 2-in PVC 0.010-in slot screen from 4.5 ft to 9.5 ft; sandpack from 2.5 ft to 9.5 ft; bentonite seal from 1.5 ft to 2.5 ft Cement well pad around protective casing from grade to 1.5 ft | | | | | | | | | | |

Ground Water Sampling Field Logs

O'Brien & Gere Engineers, Inc.

Low Flow Ground Water Sampling Log

| | | | | | |
|---------------|-----------------------|-------------------|------------------|-----------|----------------------|
| Date | <u>4/5/01</u> | Personnel | <u>T. Prawel</u> | Weather | <u>Sunny 65° +/-</u> |
| Site Name | <u>Tonawanda Coke</u> | Evacuation Method | <u>Low Flow</u> | Well # | <u>mw-1</u> |
| Site Location | <u>Tonawanda NY</u> | Sampling Method | <u>Low Flow</u> | Project # | <u>26080</u> |

Well information:

Depth of Well * 11.6 ft.Depth to Water * 6.26 ft.Length of Water Column 5.34 ft.

* Measurements taken from

X

Top of Well Casing

Top of Protective Casing

(Other, Specify)

* 4" Ø well

Water parameters: Lower submersible pump slowly through stagnant water column

15±0

Position pump in center of screened interval & maximum pumping rate of 0.5 liters/minute

Collect readings at every one minute intervals

| Elapsed Time | Depth To Water | Temperature | pH | Conductivity | Oxidation Reduction Potential | Dissolved Oxygen (mg/l) | Turbidity (NTU) | Flow Rate (ml/min). |
|--------------|----------------|-------------|------|--------------|-------------------------------|-------------------------|-----------------|---------------------|
| 5 | 6.90 | 12.97 | 4.63 | 3.04 | 52 | 0.53 | 9.89 | 250 slow |
| 10 | 6.96 | 13.16 | 4.54 | 3.01 | 38 | 0.24 | 8.78 | 210 |
| 15 | 7.00 | 13.16 | 4.59 | 2.99 | 13 | .10 | 4.33 | 210 slow |
| 20 | 7.00 | 12.74 | 4.71 | 3.02 | -13 | 0.0 | 4.22 | 200 |
| 25 | 7.00 | 12.40 | 4.89 | 3.05 | -39 | 0.0 | 3.34 | 180 |
| 30 | 7.01 | 12.31 | 5.00 | 3.10 | -52 | 0.0 | 11.2 | 200 |
| 35 | 7.00 | 12.30 | 5.05 | 3.10 | -56 | 0.0 | 17.6 | 180 |
| 40 | 6.99 | 12.45 | 5.10 | 3.07 | -62 | 0.0 | 18.5 | 180 |
| 45 | 6.99 | 12.10 | 5.10 | 3.01 | -67 | 0.0 | 18.3 | 180 |
| 50 | 6.99 | 12.92 | 5.07 | 2.95 | -71 | 0.0 | 16.1 | 180 |
| 55 | 6.99 | 12.89 | 5.03 | 2.91 | -72 | 0.0 | 15.7 | 180 |
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Water sample:

Time collected: 16:12

Total volume of purged water removed:

3 1/2 + -

Physical appearance at start

Color Milky white

Physical appearance at sampling

Color ClearOdor slightOdor strong solventSheen/Free Product NoneSheen/Free Product none

Samples collected:

| Container Size | Container Type | # Collected | Field Filtered | Preservative | Container pH |
|----------------|----------------|-------------|----------------|--------------|--------------|
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Notes: Collected Blind Dip. X-1

O'Brien & Gere Engineers, Inc.

Low Flow Ground Water Sampling Log

| | | | | | | |
|---------------|-----------------------|-------------------|------------------|-----------|--------------|----------------|
| Date | <u>6/7/01</u> | Personnel | <u>T. Prawel</u> | Weather | <u>Sunny</u> | <u>75° +/-</u> |
| Site Name | <u>Tonawanda Coke</u> | Evacuation Method | <u>Low Flow</u> | Well # | <u>MW-2</u> | |
| Site Location | <u>Tonawanda NY</u> | Sampling Method | <u>Low Flow</u> | Project # | <u>Z6080</u> | |

Well information:

Depth of Well * 11.8 ft.
 Depth to Water * 5.60 ft.
 Length of Water Column 6.2 ft.

* Measurements taken from

| | |
|----------|--------------------------|
| <u>X</u> | Top of Well Casing |
| | Top of Protective Casing |
| | (Other, Specify) |

Water parameters: Lower submersible pump slowly through stagnant water column

Position pump in center of screened interval & maximum pumping rate of 0.5 liters/minute

Collect readings at every three minute intervals

| Elapsed Time | Depth To Water | Temperature | pH | Conductivity | Oxidation Reduction Potential | Dissolved Oxygen (mg/l) | Turbidity (NTU) | Flow Rate (ml/min.) |
|--------------|----------------|-----------------------|------|------------------|-------------------------------|-------------------------|-----------------|---------------------|
| 1630 | 5.67 | | | | | | | |
| 5 | 6.46 | 12.80 | 7.01 | 312 | -156 | 1.80 | 1.80 | 420 |
| 10 | 6.81 | 12.62 | 7.10 | 309 | -161 | 0.48 | 1.07 | 300 |
| 15 | 7.06 | 12.58 | 7.13 | 307 | -155 | 0.05 | 1.01 | 260 |
| 20 | 7.30 | 12.78 | 7.14 | 3,05 | -162 | 0.00 | 0.66 | 230 |
| 25 | 7.47 | 12.93 | 7.16 | 3.03 | -161 | 0.00 | 0.95 | 170 |
| 30 | 7.67 | 12.89 | 7.17 | 3.02 | -167 | 0.00 | 0.83 | 170 |
| 35 | 7.84 | 12.85 | 7.18 | 3.00 | -183 | 0.00 | 0.67 | 150 |
| 40 | 7.94 | 12.96 | 7.18 | 2.94 | -178 | 0.00 | 1.19 | 130 |
| 45 | 8.07 | 13.05 | 7.19 | 2.88 | -168 | 0.00 | 1.11 | 130 |
| 50 | 8.17 | 13.30 | 7.18 | 2.83 | -153 | 0.00 | 1.26 | 150 |
| 55 | 8.39 | 13.46 | 7.18 | 2.78 | -144 | 0.00 | 1.57 | 150 |
| 60 | 8.58 | 13.57 | 7.20 | 2.79 | -133 | 0.00 | 1.12 | 150 |
| | Can't get | pump slower will turn | | up and pump down | | | | |
| 65 | 8.81 | 13.51 | 7.22 | 2.79 | -125 | 0.00 | 1.16 | 250 |
| 70 | 9.02 | 13.28 | 7.22 | 2.79 | -124 | 0.00 | 1.03 | 250 |
| 75 | 9.17 | 13.13 | 7.23 | 2.79 | -119 | 0.00 | 0.96 | 250 |
| 80 | 9.35 | 12.92 | 7.23 | 2.81 | -115 | 0.00 | 0.91 | 300 |
| 85 | 9.56 | 12.54 | 7.22 | 2.84 | -113 | 0.00 | 0.98 | 300 |
| 90 | 9.81 | 12.24 | 7.21 | 2.88 | -114 | 0.00 | 0.63 | 300 |

Water sample: 10.21 Dry ^AsunshineTime collected: 1045*

Total volume of purged water removed:

7

Physical appearance at start

6/9/01 - WL 10.85 @ 10¹³

Physical appearance at sampling

Color clear

Color

clear w/ AntsOdor Slight Sulfur

Odor

SlightSheen/Free Product None

Sheen/Free Product

None

Samples collected:

| Container Size | Container Type | # Collected | Field Filtered | Preservative | Container pH |
|----------------|----------------|-------------|----------------|--------------|--------------|
| | | | | | |
| | | | | | |
| | | | | | |
| | | | | | |

$$0.36 \times 653 = \\ .24 \text{ gal.}$$

Notes:

10am
12:55
* Sampled 6/9/01

O'Brien & Gere Engineers, Inc.

Low Flow Ground Water Sampling Log

| Date | 6/8/01 | Personnel | T. Prawel | Weather | Sunny 75° +/- | | | |
|--|----------------------|--|--------------------|---------------------------------|-------------------------------|---------------------------------------|---------------------------------|---------------------|
| Site Name | Tonawanda Coke | Evacuation Method | Low Flow | Well # | MW - 3 | | | |
| Site Location | Tonawanda NY | Sampling Method | Low Flow | Project # | 26080 | | | |
| Well information: | | | | | | | | |
| Depth of Well * | 11.60 ft. | * Measurements taken from | | | | | | |
| Depth to Water * | 5.94 ft. | Top of Well Casing | | | | | | |
| Length of Water Column | 5.66 ft. | X Top of Protective Casing (Other, Specify) | | | | | | |
| Water parameters: Lower submersible pump slowly through stagnant water column Position pump in center of screened interval & maximum pumping rate of 0.5 liters/minute Collect readings at every three minute intervals | | | | | | | | |
| Elapsed Time | Depth To Water | Temperature | pH | Conductivity | Oxidation Reduction Potential | Dissolved Oxygen (mg/l) | Turbidity (NTU) | Flow Rate (ml/min.) |
| 13:15 | 6.10 = 6.50 TOC | | | | | | | |
| 5 | 6.92 | 13.90 | 7.25 | 1.80 | -141 | 6.67 | 4.20 | 320 |
| 10 | 7.18 | 14.09 | 7.27 | 1.69 | -144 | 9.16 | 2.22 | 230 |
| 15 | 7.32 | 14.07 | 7.29 | 1.65 | -142 | 7.17 | 2.31 | 150 |
| 20 | 7.46 | 14.20 | 7.30 | 1.51 | -139 | 6.12 | 2.07 | 140 |
| 25 | 7.60 | 14.47 | 7.30 | 1.54 | -136 | 1.33 | 2.70 | 140 |
| 30 | 7.75 | 14.45 | 7.31 | 1.50 | -133 | 1.29 | 2.10 | 140 |
| 35 | 7.90 | 14.45 | 7.32 | 1.47 | -131 | 0.98 | 2.89 | 140 |
| 40 | 8.04 | 14.66 | 7.33 | 1.44 | -128 | 0.73 | 2.42 | 140 |
| 45 | 8.15 | 14.72 | 7.33 | 1.42 | -126 | 0.67 | 2.53 | 140 |
| 50 | 8.43 will not static | 14.46 | 7.35 | F.R. and pump down 1.38 -124 | | 0.62 | 2.89 | 300 |
| 55 | 8.69 | 14.05 | 7.38 | 1.26 | -118 | 0.52 | 2.67 | 300 |
| 60 | 8.97 | 13.90 | 7.38 | 1.21 | -110 | 0.42 | 1.54 | 340 |
| 65 | 9.23 | 13.80 | 7.35 | 1.23 | -110 | 0.21 | 1.47 | 340 |
| 70 | 9.58 | 13.63 | 7.34 | 1.28 | -113 | 0.00 | 1.26 | 380 |
| 75 | 9.95 | 13.43 | 7.32 | 1.37 | -116 | 0.00 | 1.08 | 400 |
| 80 | 10.30 | 13.20 | 7.31 | 1.47 | -119 | 0.00 | 1.29 | 400 |
| 85 | 10.72 | 13.12 | 7.31 | 1.53 | -121 | 0.00 | 1.09 | 400 |
| 90 | 11.15 Dry | 11.55 TOC | | | | | | |
| Water sample: 6/9/01 WL 10.30 @ 10:10 | | | | | | Total volume of purged water removed: | 6 | |
| Time collected: 11:40 * | | | | | | Physical appearance at start | Physical appearance at sampling | |
| Physical appearance at start | Color | clear w/ black part's | Color | | | | | |
| Odor | Odor | slight solvent | Odor | | | | | |
| Sheen/Free Product | Sheen/Free Product | None | Sheen/Free Product | | | | | |
| Samples collected: | | | | | | | | |
| Container Size | Container Type | # Collected | Field Filtered | Preservative | Container pH | | | |
| | | | | | | | | |
| | | | | | | | | |
| | | | | | | | | |
| Notes: | | | | | | | | |

* Sampled 6/9/01

O'Brien & Gere Engineers, Inc.

Low Flow Ground Water Sampling Log

| Date | 6/8/01 | Personnel | T. Pravet | Weather | Sunny 70's | | | |
|--|----------------------------|--|---------------------------------|--------------|-------------------------------|---|-----------------|---------------------|
| Site Name | Tonawanda Coke | Evacuation Method | Low Flow | Well # | mw-4 | | | |
| Site Location | Tonawanda NY | Sampling Method | Low Flow | Project # | 26080 | | | |
| Well information: | | | | | | | | |
| Depth of Well * | 11.9 ft. | * Measurements taken from | | | | | | |
| Depth to Water * | 5.52 ft. | Top of Well Casing | | | | | | |
| Length of Water Column | 6.08 ft. | X Top of Protective Casing (Other, Specify) | | | | | | |
| Water parameters: Lower submersible pump slowly through stagnant water column Position pump in center of screened interval & maximum pumping rate of 0.5 liters/minute Collect readings at every three minute intervals | | | | | | | | |
| Elapsed Time | Depth To Water | Temperature | pH | Conductivity | Oxidation Reduction Potential | Dissolved Oxygen (mg/l) | Turbidity (NTU) | Flow Rate (ml/min.) |
| 115 | 5.82 WL = 6.12 Tot | | | | | | | |
| 5 | 6.61 | 14.50 | 7.12 | 1.79 | 70 | 0.82 | 1.81 | 380 |
| 10 | 6.77 | 14.32 | 7.20 | 1.71 | 32 | 0.22 | 0.98 | 240 |
| 15 | 6.88 | 14.72 | 7.27 | 1.68 | -4 | 0.09 | 0.71 | 180 |
| 20 | 6.99 | 14.74 | 7.29 | 1.65 | -18 | 0.15 | 0.54 | 180 |
| 25 | 7.09 | 14.88 | 7.30 | 1.62 | -30 | 0.25 | 0.69 | 140 |
| 30 | 7.18 | 15.07 | 7.32 | 1.60 | -41 | 0.41 | 1.15 | 140 |
| 35 | 7.33 | 15.22 | 7.34 | 1.57 | -46 | 0.60 | 0.60 | 130 |
| 40 | 7.44 | 15.14 | 7.36 | 1.55 | -45 | 1.00 | 0.79 | 140 |
| 45 | 7.55 | 14.94 | 7.28 | 1.63 | -39 | 1.07 | 0.90 | 130 |
| 50 | 7.65 | 15.20 | 7.30 | 1.58 | -38 | 1.40 | 0.60 | 140 |
| 55 | 7.79 | 15.07 | 7.33 | 1.56 | -37 | 1.73 | 0.91 | 150 |
| 60 | 7.90 | 15.31 | 7.37 | 1.51 | -37 | 2.37 | 1.34 | 130 |
| 65 | 8.00 | 15.34 | 7.40 | 1.49 | -36 | 2.76 | 0.94 | 120 |
| 70 | 8.10 | 15.30 | 7.40 | 1.49 | -36 | 2.90 | 1.18 | 120 |
| 75 | 8.19 | 15.26 | 7.38 | 1.48 | -40 | 2.74 | 1.24 | 120 |
| 80 | 8.27 | 15.29 | 7.34 | 1.48 | -48 | 2.49 | 1.21 | 120 |
| 85 | 8.35 | 15.37 | 7.33 | 1.48 | -52 | 2.29 | 1.47 | 120 |
| 90 | Turn up Flow Rate 149.8 | will not 7.30 | 90 static 1.69 | | -57 | 1.86 | 1.03 | 400 |
| 95 | 8.69 | | | | | | | |
| Water sample: Dry @ 12.55 | | | | | | Total volume of purged water removed: 4.5 gal's | | |
| Time collected: 10:15 AM 6/9/01 - WL 5.85 | | | | | | | | |
| Physical appearance at start | | | Physical appearance at sampling | | | | | |
| Color | clear | = 6.7 x 6.3 | Color | clear | | | | |
| Odor | None | = 4.37 gal.) | Odor | slight | | | | |
| Sheen/Free Product | None | | Sheen/Free Product | None | | | | |
| Samples collected: | | | | | | | | |
| Container Size | Container Type | # Collected | Field Filtered | Preservative | Container pH | | | |
| | | | | | | | | |
| | | | | | | | | |
| | | | | | | | | |
| | | | | | | | | |
| Notes: NL7.85 @ 1500 | | | | | | | | |
| * Sampled 6/9/01 | | | | | | | | |

O'Brien & Gere Engineers, Inc.

Low Flow Ground Water Sampling Log

| | | | | | |
|---------------|-----------------------|-------------------|-----------------|-----------|-------------------|
| Date | <u>6/5/01</u> | Personnel | <u>TPP</u> | Weather | <u>Sunny 65°+</u> |
| Site Name | <u>Tonawanda Coke</u> | Evacuation Method | <u>Low Flow</u> | Well # | <u>MW-5</u> |
| Site Location | <u>Tonawanda NY</u> | Sampling Method | <u>Low Flow</u> | Project # | <u>26080</u> |

Well information:

Depth of Well * 11.1 ft.
 Depth to Water * 6.67 ft.
 Length of Water Column 4.43 ft.

* Measurements taken from

| |
|--------------------------|
| X |
| Top of Well Casing |
| Top of Protective Casing |
| (Other, Specify) |

Top of Well Casing
 Top of Protective Casing
 (Other, Specify)

Water parameters: Lower submersible pump slowly through stagnant water column

Position pump in center of screened interval & maximum pumping rate of 0.5 liters/minute

Collect readings at every three minute intervals

| Elapsed Time | Depth To Water | Temperature | pH | Conductivity | Oxidation Reduction Potential | Dissolved Oxygen (mg/l) | Turbidity (NTU) | Flow Rate (ml/min.) |
|---------------------------|----------------|---------------|------|--------------|-------------------------------|-------------------------|-----------------|---------------------|
| 0 | 6.85 = 7.42 Td | 15.79 | 6.13 | 382 | -44 | 7.48 | 41.2 | 310 slow |
| 5 | 8.50 | 13.53 | 6.97 | 2.53 | -122 | 7.91 | 3.45 | 210 slow |
| 10 | 8.91 | 12.91 | 7.06 | 2.35 | -121 | 8.16 | 13.2 | 180 |
| 15 | 9.22 | 12.85 | 7.09 | 2.32 | -108 | 3.24 | 34.8 | 190 slow |
| 20 | 9.60 | 11.83 | 7.10 | 2.30 | -103 | 1.16 | 6.68 | 165 |
| 25 | 10.11 | 11.96 | 7.10 | 2.28 | -99 | 0.31 | 1.21 | 135 |
| 30 | 10.40 | 11.90 | 7.10 | 2.29 | -92 | 0.15 | 2.63 | 110 |
| 35 | 10.85 | 11.94 | 7.09 | 2.27 | -96 | 0.24 | 2.33 | 100 |
| | 10.47 = 11.00 | | | | | | | |
| 11:35 | Dry - C | 2.5 gal's +/- | | | | | | |
| | | | | | | | | |
| Per T. Alexander | | | | | | | | |
| 2 Jars | | | | | | | | |
| 2 - 1/2 full Liters Glass | | | | | | | | |
| '1/2 full Metals | | | | | | | | |
| '1/2 full CN | | | | | | | | |
| | | | | | | | | |
| Sampled 16:00 6/7/01 | | | | | | | | |
| | | | | | | | | |
| | | | | | | | | |

Water sample:

Time collected: 16:00 6/7/01

Total volume of purged water removed:

Physical appearance at start

Physical appearance at sampling

Color ClearColor Odor Slight SolventOdor Sheen/Free Product NoneSheen/Free Product

Samples collected:

| Container Size | Container Type | # Collected | Field Filtered | Preservative | Container pH |
|----------------|----------------|-------------|----------------|--------------|--------------|
| | | | | | |
| | | | | | |
| | | | | | |
| | | | | | |

Notes: 16:50 WL 9.91 Bmp8.85 @ 13:30 = 0.35 gal13:00 WL 9.36 Bmp

O'Brien & Gere Engineers, Inc.

Low Flow Ground Water Sampling Log

| | | | | | |
|---------------|----------------|-------------------|-----------|-----------|--------------------|
| Date | 6/6/01 | Personnel | T. Prawel | Weather | overcast/sunny 65° |
| Site Name | Tonawanda Coke | Evacuation Method | Low Flow | Well # | MW-6 |
| Site Location | Tonawanda NY | Sampling Method | Low Flow | Project # | 26080 |

Well information:

Depth of Well * 11.4 ft.
 Depth to Water * 5.55 ft.
 Length of Water Column 5.85 ft.

* Measurements taken from

| |
|---|
| X |
| |
| |

Top of Well Casing
 Top of Protective Casing
 (Other, Specify)

Water parameters: Lower submersible pump slowly through stagnant water column

Position pump in center of screened interval & maximum pumping rate of 0.5 liters/minute

Collect readings at every three minute intervals

| Elapsed Time | Depth To Water | Temperature | pH | Conductivity | Oxidation Reduction Potential | Dissolved Oxygen (mg/l) | Turbidity (NTU) | Flow Rate (ml/min.) |
|------------------|----------------|-------------|------|--------------|-------------------------------|-------------------------|-----------------|---------------------|
| 10 ¹⁰ | 5.65 | | | | | | | |
| 5 | 5.70 | 12.22 | 3.73 | 2.65 | 255 | 7.60 | 11.5 | 320 |
| 10 | 5.70 | 12.09 | 4.23 | 2.56 | 163 | 6.88 | 3.11 | 320 |
| 15 | 5.70 | 12.07 | 4.32 | 2.56 | 135 | 6.36 | 1.32 | 330 |
| 20 | 5.70 | 12.01 | 4.36 | 2.58 | 116 | 5.97 | 0.78 | 330 |
| 25 | 5.70 | 12.00 | 4.38 | 2.59 | 107 | 4.95 | 0.54 | 320 |
| 30 | 5.70 | 11.99 | 4.39 | 2.61 | 98 | 4.15 | 0.48 | 320 |
| 35 | 5.70 | 11.98 | 4.40 | 2.61 | 93 | 3.89 | 0.50 | 320 |
| 40 | 5.70 | 12.00 | 4.41 | 2.61 | 88 | 3.68 | 0.57 | 320 |
| 45 | 5.71 | 11.99 | 4.42 | 2.61 | 85 | 3.31 | 0.39 | 350 8000 |
| 50 | 5.71 | 11.96 | 4.42 | 2.61 | 81 | 3.17 | 0.18 | 350 |
| 55 | 5.71 | 11.95 | 4.42 | 2.60 | 78 | 2.93 | 0.19 | 350 |
| 60 | 5.71 | 11.93 | 4.42 | 2.60 | 74 | 2.80 | 0.17 | 330 |
| 65 | 5.71 | 11.87 | 4.43 | 2.60 | 72 | 2.36 | 0.50 | 330 |
| 70 | 5.71 | 11.83 | 4.43 | 2.60 | 68 | 1.91 | 0.14 | 330 |
| 75 | 5.71 | 11.75 | 4.44 | 2.63 | 70 | 0.0 | 0.13 | 330 |
| 80 | 5.71 | 11.77 | 4.45 | 2.64 | 69 | 0.0 | 0.14 | 330 |
| 85 | 5.71 | 11.89 | 4.46 | 2.63 | 70 | 0.0 | 0.20 | 330 |

Water sample:

Time collected: 1245Total volume of purged water removed: 7

Physical appearance at start

Color clear
 Odor Slight Solvent
 Sheen/Free Product None

Physical appearance at sampling

Color clear
 Odor Heavy Solvent
 Sheen/Free Product None

Samples collected:

| Container Size | Container Type | # Collected | Field Filtered | Preservative | Container pH |
|----------------|----------------|-------------|----------------|--------------|--------------|
| | | | | | |
| | | | | | |
| | | | | | |
| | | | | | |

Notes:

| Date | <u>6/6/01</u> | Personnel | <u>T. Prawel</u> | Weather | <u>Sunny 56° +/-</u> | | | |
|---|-----------------------|---------------------------------------|-------------------------------------|--------------------------|-------------------------------|-------------------------|-----------------|---------------------|
| Site Name | <u>Tonawanda Coke</u> | Evacuation Method | <u>Low Flow</u> | Well # | <u>MW-7</u> | | | |
| Site Location | <u>Tonawanda NY</u> | Sampling Method | <u>Low Flow</u> | Project # | <u>Z6080</u> | | | |
| Well information: | | | | | | | | |
| Depth of Well * | <u>11.8</u> | ft. | * Measurements taken from | | | | | |
| Depth to Water * | <u>6.22</u> | ft. | <input type="checkbox"/> | Top of Well Casing | | | | |
| Length of Water Column | <u>5.58</u> | ft. | <input checked="" type="checkbox"/> | Top of Protective Casing | | | | |
| | | | <input type="checkbox"/> | (Other, Specify) | | | | |
| Water parameters: Lower submersible pump slowly through stagnant water column Position pump in center of screened interval & maximum pumping rate of 0.5 liters/minute Collect readings at every three minute intervals <i>9:00</i> | | | | | | | | |
| Elapsed Time | Depth To Water | Temperature | pH | Conductivity | Oxidation Reduction Potential | Dissolved Oxygen (mg/l) | Turbidity (NTU) | Flow Rate (ml/min.) |
| 0 | <u>6.28 = 6.61</u> | <u>70C</u> | | | | | | <u>450</u> |
| 5 | <u>6.61</u> | <u>13.84</u> | <u>2.24</u> | <u>5.54</u> | <u>388</u> | <u>.50</u> | <u>1.33</u> | <u>450</u> |
| 10 | <u>6.61</u> | <u>13.56</u> | <u>2.20</u> | <u>5.68</u> | <u>386</u> | <u>4.72</u> | <u>0.88</u> | <u>420</u> |
| 15 | <u>6.61</u> | <u>13.53</u> | <u>2.25</u> | <u>5.71</u> | <u>382</u> | <u>5.50</u> | <u>0.90</u> | <u>410</u> |
| 20 | <u>6.61</u> | <u>13.56</u> | <u>2.26</u> | <u>5.66</u> | <u>381</u> | <u>0.05</u> | <u>0.71</u> | <u>410</u> |
| 25 | <u>6.61</u> | <u>13.45</u> | <u>2.27</u> | <u>5.63</u> | <u>380</u> | <u>10.50</u> | <u>0.87</u> | <u>410</u> |
| 30 | <u>6.61</u> | <u>13.37</u> | <u>2.27</u> | <u>5.61</u> | <u>379</u> | <u>5.72</u> | <u>0.98</u> | <u>400</u> |
| 35 | <u>6.61</u> | <u>13.40</u> | <u>2.28</u> | <u>5.58</u> | <u>378</u> | <u>5.12</u> | <u>0.66</u> | <u>400</u> |
| 40 | <u>6.61</u> | <u>13.27</u> | <u>2.28</u> | <u>5.56</u> | <u>377</u> | <u>4.20</u> | <u>0.46</u> | <u>400</u> |
| 45 | <u>6.61</u> | <u>13.23</u> | <u>2.27</u> | <u>5.52</u> | <u>376</u> | <u>0.23</u> | <u>0.42</u> | <u>400</u> |
| 50 | <u>6.61</u> | <u>13.22</u> | <u>2.25</u> | <u>5.51</u> | <u>375</u> | <u>0.0</u> | <u>0.58</u> | <u>400</u> |
| 55 | <u>6.61</u> | <u>13.33</u> | <u>2.26</u> | <u>5.49</u> | <u>375</u> | <u>0.0</u> | <u>0.33</u> | <u>400</u> |
| 60 | <u>6.61</u> | <u>13.30</u> | <u>2.28</u> | <u>5.49</u> | <u>376</u> | <u>0.0</u> | <u>0.45</u> | <u>400</u> |
| Water sample: | | | | | | | | |
| Time collected: | <u>10:05</u> | Total volume of purged water removed: | | | | | | <u>7</u> |
| Physical appearance at start | | | Physical appearance at sampling | | | | | |
| Color | <u>clear</u> | Color | <u>clear</u> | | | | | |
| Odor | <u>none</u> | Odor | <u>slight solvent</u> | | | | | |
| Sheen/Free Product | <u>none</u> | Sheen/Free Product | <u>none</u> | | | | | |
| Samples collected: | | | | | | | | |
| Container Size | Container Type | # Collected | Field Filtered | Preservative | Container pH | | | |
| | | | | | | | | |
| | | | | | | | | |
| | | | | | | | | |
| | | | | | | | | |
| Notes: | | | | | | | | |

O'Brien & Gere Engineers, Inc.

Low Flow Ground Water Sampling Log

| | | | | | |
|---------------|----------------|-------------------|----------|-----------|--------------|
| Date | 4/5/01 | Personnel | TPP | Weather | Sunny 65°+/- |
| Site Name | Tonawanda Coke | Evacuation Method | Low Flow | Well # | MW-8 |
| Site Location | Tonawanda NY | Sampling Method | Low Flow | Project # | 26080 |

Well information:

Depth of Well * 15.2 ft.
 Depth to Water * 3.81 ft.
 Length of Water Column 11.39 ft.

* Measurements taken from

| |
|--------------------------|
| X |
| Top of Well Casing |
| Top of Protective Casing |
| (Other, Specify) |

Top of Well Casing
 Top of Protective Casing
 (Other, Specify)

Water parameters: Lower submersible pump slowly through stagnant water column

Position pump in center of screened interval & maximum pumping rate of 0.5 liters/minute

13' 0"

Collect readings at every three minute intervals

| Elapsed Time | Depth To Water | Temperature | pH | Conductivity | Oxidation Reduction Potential | Dissolved Oxygen (mg/l) | Turbidity (NTU) | Flow Rate (ml/min.) |
|--------------|----------------|-------------|------|--------------|-------------------------------|-------------------------|-----------------|---------------------|
| 0 | 4.91 | 14.85 | 7.08 | 2.21 | -93 | 0.93 | 3.99 | 230 |
| 5 | 5.20 | 11.57 | 7.33 | 2.15 | -76 | 1.37 | 2.20 | 200 |
| 10 | 6.61 | 11.20 | 7.38 | 2.10 | -55 | 1.41 | 1.52 | 200 |
| 15 | 7.74 | 11.20 | 7.38 | 2.08 | -37 | 1.35 | 1.62 | 190 |
| 20 | 8.92 | 11.21 | 7.39 | 2.08 | -25 | 1.26 | 1.57 | 190 |
| 25 | 10.21 | 11.11 | 7.37 | 2.06 | -17 | 1.10 | 0.99 | 200 |
| 30 | 10.51 | 11.00 | 7.37 | 2.06 | -10 | 1.07 | 2.05 | 200 |
| 35 | 10.87 | 11.02 | 7.40 | 2.05 | -3 | 1.33 | 1.09 | 200 slow |
| 40 | 11.17 | 11.02 | 7.38 | 2.03 | 3 | 1.36 | 1.18 | 170 |
| 45 | 11.45 | 11.06 | 7.39 | 2.03 | 8 | 0.98 | 0.68 | 170 |
| 50 | 11.72 | 11.04 | 7.46 | 2.03 | 11 | 1.01 | 0.68 | 170 |
| 55 | 12.02 | 10.96 | 7.49 | 2.03 | 14 | 1.00 | 0.45 | 170 |
| 60 | 12.28 | 10.95 | 7.51 | 2.03 | 17 | 1.00 | 0.62 | 170 slow |
| 65 | 12.49 | 11.13 | 7.53 | 2.04 | 20 | 1.00 | 1.01 | 140 |
| 70 | 12.72 | 11.14 | 7.46 | 2.04 | 23 | 1.19 | 0.68 | 140 |
| 75 | 12.74 | 10.50 | 7.47 | 2.07 | 29 | 1.21 | 0.62 | 140 |
| 80 | 12.75 | 10.51 | 7.47 | 2.07 | 29 | 1.21 | 0.62 | 140 |

Water sample:

Time collected: 14:20Total volume of purged water removed: 4

Physical appearance at start

Color Clear
 Odor None
 Sheen/Free Product None

Physical appearance at sampling

Color Clear
 Odor Slight
 Sheen/Free Product None

Samples collected:

| Container Size | Container Type | # Collected | Field Filtered | Preservative | Container pH |
|----------------|----------------|-------------|----------------|--------------|--------------|
| | | | | | |
| | | | | | |
| | | | | | |
| | | | | | |

Notes:

O'Brien & Gere Engineers, Inc.**Low Flow Ground Water Sampling Log**

| | | | | | |
|---------------|-----------------------|-------------------|------------------|-----------|-------------------------|
| Date | <u>6/6/01</u> | Personnel | <u>T. Prawel</u> | Weather | <u>overcast 68° + L</u> |
| Site Name | <u>Tonawanda Coke</u> | Evacuation Method | <u>Low Flow</u> | Well # | <u>MW-9</u> |
| Site Location | <u>Tonawanda NY</u> | Sampling Method | <u>Low Flow</u> | Project # | <u>Z6080</u> |

Well information:

Depth of Well * 12.90 ft.
 Depth to Water * 6.70 ft.
 Length of Water Column 6.20 ft.

* Measurements taken from

| |
|-------------------------------------|
| <input checked="" type="checkbox"/> |
| |
| |

Top of Well Casing

Top of Protective Casing

(Other, Specify)

Water parameters: Lower submersible pump slowly through stagnant water column

Position pump in center of screened interval & maximum pumping rate of 0.5 liters/minute

Collect readings at every three minute intervals

| Elapsed Time | Depth To Water | Temperature | pH | Conductivity | Oxidation Reduction Potential | Dissolved Oxygen (mg/l) | Turbidity (NTU) | Flow Rate (ml/min.) |
|--------------|----------------|-------------|------|--------------|-------------------------------|-------------------------|-----------------|---------------------|
| 1307 | 6.77 | | | | | | | |
| 5 | 7.30 | 12.98 | 6.99 | 2.02 | -6 | 7.39 | 6.21 | 390 |
| 10 | 7.25 | 12.96 | 7.07 | 2.01 | -36 | 7.23 | 2.48 | 330 |
| 15 | 7.26 | 12.90 | 7.12 | 1.93 | -58 | 6.76 | 1.75 | 330 |
| 20 | 7.27 | 13.00 | 7.16 | 1.86 | -74 | 5.88 | 1.14 | 330 |
| 25 | 7.28 | 13.01 | 7.19 | 1.82 | -86 | 5.52 | 0.98 | 330 |
| 30 | 7.28 | 13.00 | 7.21 | 1.79 | -94 | 5.26 | 0.73 | 320 |
| 35 | 7.29 | 13.00 | 7.24 | 1.77 | -103 | 4.95 | 0.60 | 330 |
| 40 | 7.24 | 13.06 | 7.25 | 1.73 | -108 | 4.67 | 0.55 | 300 |
| 45 | 7.25 | 13.04 | 7.26 | 1.74 | -110 | 0.54 | 0.54 | 300 |
| 50 | 7.25 | 13.01 | 7.27 | 1.72 | -114 | 0.0 | 0.52 | 300 |
| 55 | 7.25 | 13.09 | 7.28 | 1.71 | -117 | 0.00 | 0.57 | 300 |
| 60 | 7.25 | 13.18 | 7.29 | 1.69 | -119 | 0.0 | 0.52 | 300 |
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Water sample:Time collected: 1415

Total volume of purged water removed:

5

Physical appearance at start

Color Clear
 Odor Slight Solvent
 Sheen/Free Product None

Physical appearance at sampling

Color clear
 Odor solvent
 Sheen/Free Product none

Samples collected:

| Container Size | Container Type | # Collected | Field Filtered | Preservative | Container pH |
|----------------|----------------|-------------|----------------|--------------|--------------|
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Notes:

O'Brien & Gere Engineers, Inc.

Low Flow Ground Water Sampling Log

| Date | 6/6/01 | Personnel | T. Pravell | Weather | overcast 68°+/- | | | |
|--|----------------|--|--------------------|--------------|-------------------------------|-------------------------|-----------------|---------------------|
| Site Name | Tonawanda Coke | Evacuation Method | Low Flow | Well # | MW - 11 | | | |
| Site Location | Tonawanda NY | Sampling Method | Low Flow | Project # | 26080 | | | |
| Well information: | | | | | | | | |
| Depth of Well * | 11.40 ft. | * Measurements taken from | | | | | | |
| Depth to Water * | 5.94 ft. | <input checked="" type="checkbox"/> Top of Well Casing | | | | | | |
| Length of Water Column | 5.46 ft. | <input type="checkbox"/> Top of Protective Casing | | | | | | |
| (Other, Specify) | | | | | | | | |
| Water parameters: Lower submersible pump slowly through stagnant water column Position pump in center of screened interval & maximum pumping rate of 0.5 liters/minute Collect readings at every three minute intervals | | | | | | | | |
| Elapsed Time | Depth To Water | Temperature | pH | Conductivity | Oxidation Reduction Potential | Dissolved Oxygen (mg/l) | Turbidity (NTU) | Flow Rate (ml/min.) |
| 1440 | 5.92 | | | | | | | |
| 5 | 7.09 | 1324 | 7.52 | 1.81 | +4 | 0.51 | 3.65 | 290 |
| 10 | 7.45 | 1346 | 7.52 | 1.79 | -4 | 0.41 | 2.68 | 210 |
| 15 | 7.76 | 1339 | 7.49 | 1.79 | -16 | 0.04 | 2.18 | 180 |
| 20 | 8.23 | 1344 | 7.46 | 1.81 | -32 | 0.02 | 2.80 | 150 |
| 25 | 8.70 | 1349 | 7.45 | 1.80 | -42 | 0.00 | 5.47 | 150 |
| 30 | 9.15 | 1341 | 7.46 | 1.80 | -50 | 0.00 | 16.4 | 150 |
| 35 | 9.50 | 1306 | 7.48 | 1.81 | -46 | 0.00 | 15.5 | 150 |
| 40 | 9.76 | 1275 | 7.41 | 1.79 | -25 | 0.50 | 7.49 | 150 |
| 45 | | | | | | | | |
| 50 | | | | | | | | |
| 1423 | Dry C | 9.86 BMLP | w/ 2 gal +/- Purge | | | | | |
| 1410 | 6/7/01 | Sampled | R recharge | | | | | |
| Water sample: | | | | | | | | |
| Time collected: | 1410 | Total volume of purged water removed: | 2 | | | | | |
| Physical appearance at start | | Physical appearance at sampling | | | | | | |
| Color | Clear | Color | Clear | | | | | |
| Odor | Slight | Odor | Slight | | | | | |
| Sheen/Free Product | None | Sheen/Free Product | None | | | | | |
| Samples collected: | | | | | | | | |
| Container Size | Container Type | # Collected | Field Filtered | Preservative | Container pH | | | |
| | | | | | | | | |
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| | | | | | | | | |
| Notes: NL. 6.15 - 6/7/01 C 13 ³⁹ | | | | | | | | |

O'Brien & Gere Engineers, Inc.

Low Flow Ground Water Sampling Log

Date 6/8/01
Site Name Tonawanda
Site Location Cohr

Personnel T. Prawel
Evacuation Method Low Flow
Sampling Method Low Flow

Weather Sunny 75°+
Well # MW-12
Project # 26080

Well information:

Depth of Well * 12.30 ft.
Depth to Water * 12.00 ft.
Length of Water Column 0.30 ft.

* Measurements taken from

| |
|---|
| X |
| |
| |

Top of Well Casing
Top of Protective Casing
(Other, Specify)

Water parameters: Lower submersible pump slowly through stagnant water column

Position pump in center of screened interval & maximum pumping rate of 0.5 liters/minute

Collect readings at every three minute intervals

| Elapsed Time | Depth To Water | Temperature | pH | Conductivity | Oxidation Reduction Potential | Dissolved Oxygen (mg/l) | Turbidity (NTU) | Flow Rate (ml/min.) |
|--------------|----------------------------|-------------|----|--------------|-------------------------------|-------------------------|-----------------|---------------------|
| | | | | | | | | |
| | No Water in well to sample | | | | | | | |
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Water sample:

Time collected: _____

Total volume of purged water removed: 0

Physical appearance at start

Physical appearance at sampling

Color _____

Color _____

Odor _____

Odor _____

Sheen/Free Product _____

Sheen/Free Product _____

Samples collected:

| Container Size | Container type | # Collected | Field Filtered | Preservative | Container pH |
|----------------|----------------|-------------|----------------|--------------|--------------|
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Notes:

Brown & Quert Laboratories, Inc.

5000 Brittonfield Parkway
East Syracuse, New York 13057
(315) 437-0200

Chain of Custody

| Client: | Aliced / Harragwell | | Phone # | 26080 | Analysis/Method | |
|---------------------------|---------------------|----------------|---------------|---------------|-------------------|----------------|
| Project: | Tsunananda Colle | | | | | |
| Sampled by: | T. Powell | | | | | |
| Client Contact: | | | | | | |
| Sample Description | | | | | | |
| Sample Location | Date Collected | Time Collected | Sample Matrix | Comp. or Grab | No. of Containers | Comments |
| MW-2 | 6/8/01 | 1045 | Water | Grab | 7 | 3 1 1 - |
| MW-4 | 6/8/01 | 1115 | Water | Grab | 8 | 3 2 1 - |
| MW-3 | 6/8/01 | 1140 | Water | Grab | 8 | 3 2 1 - |
| Trip Blank | | | | | | |
| | 6/8/01 | | | | | |
| Requisitioned by: | <u>Tom Powell</u> | | Date: 6/8/01 | Time: 1300 | Received by: | Date: _____ |
| Reinquired by: | | | Date: _____ | Time: _____ | Received by: | Date: _____ |
| Reinquired by: | | | Date: _____ | Time: _____ | Received by Lab: | Date: _____ |
| Shipment Method: | FED Ex | | | | A/R Bill Number: | 8271-1284-1739 |
| Comments: | | | | | | |
| Turnaround Time Required: | | | | | | |
| Routine | | | | | | |
| Rush (Specify) | | | | | | |
| Cooler Temperature: | | | | | | |

Brien & Gere Laboratories, Inc.

5000 Brittonfield Parkway
East Syracuse, New York 13057
(315) 437-0200

Chain of Custody

| Client: | Project: | Sampled by: | Client Contact: | Analysis/Method | | | | | | | | | |
|---------------------------|----------------|----------------|-----------------|-----------------|-------------------|---------------|---------------|-------------------|-----------|---|---|---|--|
| | | | | Date Collected | Time Collected | Sample Matrix | Comp. or Grab | No. of Containers | Comments | | | | |
| Raw - 11 | 6/7/01 | 14:10 | water | Grab | 8 | 3 | 2 | 1 | 1 | 1 | 1 | | |
| MW-5 | 6/7/01 | 11:00 | water | Grab | 6 | 2 | 1 | 1 | 1 | 1 | 1 | | |
| | | | | | | | | | | | | | |
| Sample Description | | | | | | | | | | | | | |
| Sample Location | Date Collected | Time Collected | Sample Matrix | Comp. or Grab | No. of Containers | Comments | | | | | | | |
| Raw - 11 | 6/7/01 | 14:10 | water | Grab | 8 | 3 | 2 | 1 | 1 | 1 | 1 | 1 | |
| MW-5 | 6/7/01 | 11:00 | water | Grab | 6 | 2 | 1 | 1 | 1 | 1 | 1 | 1 | |
| | | | | | | | | | | | | | |
| Retrnished by: | T. P. Blank | 6/7/01 | 1:00 | Date: 6/8/01 | Time: 17:00 | Received by: | Date: | Time: | Comments: | | | | |
| Retrnished by: | T. P. Blank | 6/7/01 | 1:00 | Date: 6/8/01 | Time: 17:00 | Received by: | Date: | Time: | Comments: | | | | |
| Retrnished by: | T. P. Blank | 6/7/01 | 1:00 | Date: 6/8/01 | Time: 17:00 | Received by: | Date: | Time: | Comments: | | | | |
| Shipment Method: | FED EX | | | | | | | | Comments: | | | | |
| ArbID Number: | 827 | 12 | 84 | 1740 | | | | | Comments: | | | | |
| Turnaround Time Required: | Routine | | | | | | | | Comments: | | | | |
| | Rush (Specify) | | | | | | | | Comments: | | | | |
| Cooler Temperature: | | | | | | | | | Comments: | | | | |

Original Laboratory Copy Client

Brien & Gere Laboratories, Inc.

Chain of Custody

5000 Brittonfield Parkway
East Syracuse, New York 13057
(315) 437-0200

| Client: | Project: | Sampled by: | Client Contact: | Analysis/Method | | | | | | Comments |
|---------------------------|------------------|--------------|-----------------|------------------|----------------|---------------|---------------|-------------------|---|----------|
| | | | | Date Collected | Time Collected | Sample Matrix | Comp. or Grab | No. of Containers | | |
| Sample Description | | | | | | | | | | |
| MW-10 | 6/5/01 | 1045 | Wat | Grab | 8 | 3 | 2 | 1 | - | |
| MW-8 | 6/5/01 | 1420 | Wat | Grab | 8 | 3 | 2 | 1 | - | |
| MW-1 | 6/5/01 | 1610 | Wat | Grab | 8 | 3 | 2 | 1 | - | |
| MW-7 | 6/6/01 | 1005 | water | Grab | 24 | 9 | 4 | 3 | 3 | |
| MW-6 | 6/6/01 | 1245 | water | Grab | 8 | 3 | 2 | 1 | - | |
| MW-9 | 6/6/01 | 1415 | water | Grab | 8 | 3 | 2 | 1 | - | |
| X-1 | - | - | Wat | Grab | 8 | 3 | 2 | 1 | - | |
| Tri-Blank | - | - | Water | - | - | - | - | - | - | |
| Requisitioned by: | Treasor Off Hand | Date: 6/6/01 | Time: 1700 | Received by: | | | | | | |
| Requisitioned by: | | Date: | Time: | Received by: | | | | | | |
| Requisitioned by: | | Date: | Time: | Received by Lab: | | | | | | |
| Shipment Method: | Fed Ex | | | AIRBN Number: | 82711284 | 1680 | | | | |
| Turnaround Time Required: | | | | | | | | | | |
| Routine _____ | | | | | | | | | | |
| Rush (Specify) _____ | | | | | | | | | | |
| Cooler Temperature: _____ | | | | | | | | | | |
| Comments: | | | | | | | | | | |
| 82711284 1691 | | | | | | | | | | |
| 82711284 1706 | | | | | | | | | | |

Ground Water Elevation Summary

Attachment 1

**Former Allied Specialty Chemical Site
Tonawanda, New York**

Ground Water Elevations

| Well I.D. | Total Depth (ft bTOC) | Top of Casing Elevation (ft) | 9/27/91 | | 12/5/91 | | 3/5/92 | | 6/15/92 | | 5/12/99 | | 6/5/01 | |
|-----------|-----------------------|------------------------------|-------------|-----------|-------------|-----------|-------------|-----------|-------------|-----------|-------------|-----------|-------------|-----------|
| | | | Depth to GW | Elevation |
| MW-1 | 11.6 | 585.39 | 8.20 | 577.19 | 6.49 | 578.90 | 6.88 | 578.51 | 7.00 | 578.39 | 6.59 | 578.80 | 6.26 | 579.13 |
| MW-2 | 11.8 | 583.78 | 0.10 | 574.68 | 9.22 | 574.56 | 8.67 | 575.11 | 7.68 | 576.10 | 6.08 | 577.70 | 5.60 | 578.18 |
| MW-3 | 11.6 | 582.51 | 8.12 | 574.39 | 6.96 | 575.55 | 5.77 | 576.74 | 6.48 | 576.03 | 6.20 | 576.31 | 5.94 | 576.57 |
| MW-4 | 11.9 | 583.87 | 6.01 | 577.86 | 5.84 | 578.03 | 5.92 | 577.95 | 5.98 | 577.89 | 5.83 | 578.04 | 5.82 | 578.35 |
| MW-5 | 11.1 | 586.47 | 8.54 | 577.93 | 6.10 | 580.37 | 6.53 | 579.94 | 6.76 | 579.71 | 6.83 | 579.64 | 6.67 | 579.80 |
| MW-6 | 11.4 | 584.62 | 7.52 | 577.10 | 5.90 | 578.72 | 6.22 | 578.40 | 6.30 | 578.32 | 5.93 | 578.69 | 5.55 | 579.07 |
| MW-7 | 11.8 | 585.24 | 8.22 | 577.02 | 6.57 | 578.67 | 6.88 | 578.36 | 6.97 | 578.27 | 6.60 | 578.64 | 6.22 | 579.02 |
| MW-8 | 15.2 | 586.08 | 7.50 | 578.58 | 5.70 | 580.38 | 9.50 | 576.58 | 7.42 | 578.66 | 4.18 | 581.90 | 3.81 | 582.27 |
| MW-9 | 12.9 | 585.13 | N | N | N | N | N | N | N | N | N | N | N | 578.43 |
| MW-10 | 12.6 | 586.28 | N | N | N | N | N | N | N | N | N | N | N | 578.54 |
| MW-11 | 11.4 | 584.80 | N | N | N | N | N | N | N | N | N | N | N | 578.86 |
| MW-12 | 12.2 | 581.48 | N | N | N | N | N | N | N | N | N | N | N | DRY |

36-inch invert - Inlet A 574.36

36-inch invert - Outfall #1 570.89

48-inch invert - Outfall #2 571.26

NOTES: NI - Not Installed

Elevations recorded by ERM-Northeast during 1991 and 1992 and by O'Brien & Gere during 1999 and 2001