

2014-2015 OU2 GROUNDWATER INVESTIGATION
RE122D1, RE122D2, RE122D3 (VPB156)
INSTALLATION REPORT

NAVAL WEAPONS INDUSTRIAL RESERVE PLANT
BETHPAGE, NY

Prepared for:



Department of the Navy
Naval Facilities Engineering Command, Atlantic
9324 Virginia Avenue
Building Z-144
Norfolk, Virginia 23511

January 2016

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NWIRP
SITE 1 OU2
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Department of the Navy
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9324 Virginia Avenue
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Prepared by:



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Contract Number: N62470-11-D-8013
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Brian Caldwell
Contract Task Order Manager

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List of Acronyms and Abbreviations

| | |
|----------|---|
| AOC | Area of Concern |
| bgs | below ground surface |
| COR | Continuously Operating Reference |
| EPA | Environmental Protection Agency, United States |
| ft | feet |
| GOCO | Government-Owned Contractor-Operated |
| GPS | Global Positioning System |
| IDW | Investigation Derived Waste |
| IR | Installation Restoration |
| Katahdin | Katahdin Analytical Services |
| NAD | North American Datum |
| NAVD | North American Vertical Datum |
| NAVFAC | Naval Facilities Engineering Command |
| NG | Northrop Grumman |
| NTU | nephelometric turbidity units |
| NWIRP | Naval Weapons Industrial Reserve Plant |
| NYSDEC | New York State Department of Environmental Conservation |
| ONCT | On-site Containment Treatment System |
| OU | Operable Unit |
| PCBs | Polychlorinated Biphenyls |
| POTW | Publicly Owned Treatment Works |
| ppb | parts per billion |
| PPE | Personal Protective Equipment |
| PVC | Polyvinylchloride |
| SAP | Sampling and Analysis Plan |
| SVOC | Semivolatile Organic Compounds |
| TCE | Trichloroethene |
| TCL | Target Compound List |
| TCLP | Toxicity Characteristic Leaching Procedure |
| TOC | Total Organic Carbon |
| UFP | United Federal Programs |
| US | United States |
| VOC | Volatile Organic Compounds |
| VPB | Vertical Profile Boring |

1.0 PROJECT BACKGROUND

Resolution Consultants has prepared this Data Summary Report for the Naval Facilities Engineering Command (NAVFAC), Mid-Atlantic under contract task order WE15 Contract N62470-11-D-8013. This report describes the installation of three monitoring wells and one quarterly groundwater monitoring event (specifically at the Vertical Profile Boring [VPB] 156 location) in 2014 and 2015 for the Naval Weapons Industrial Reserve Plant (NWIRP) Bethpage Operable Unit (OU) 2 Site 1 offsite plume. NWIRP Bethpage is located in east-central Nassau County, Long Island, New York, approximately 30 miles east of New York City (Figure 1).

1.1 Scope and Objectives

This report provides information on the installation of RE122D1, RE122D2 and RE122D3. The purpose of this investigation was to ascertain subsurface conditions and contaminant levels south of the On-site Containment Treatment system (ONCT) and to help ascertain the effectiveness of the ONCT. In addition, these wells help define the northwestern extent of the RE108 Hot Spot (defined as an area >1,000 parts per billion [ppb] of total volatile organic compounds [VOCs] north of Hempstead Turnpike). The locations of RE122D1, RE122D2 and RE122D3, VPBs and monitoring well locations are shown in Figure 2.

The field investigation included completing three monitoring wells, well development, soil/groundwater analysis, groundwater grab samples, and surveying. Field tasks were conducted in 2014 and 2015 in accordance with the *United Federal Programs Sampling and Analysis Plan (UFP SAP)*, Bethpage, New York (Resolution Consultants, 2013a). In addition, the work adhered to the following UFP SAP Addendums: *Groundwater Sampling Using Low Stress (Low Flow) Purging and Sampling Protocol* (Resolution Consultants, 2013b) and *Installation of Vertical Profile Borings and Monitoring Wells* (Resolution Consultants, 2013c).

Documentation of these activities is included in Appendix A of this report.

1.2 Site History

NWIRP Bethpage is in the Hamlet of Bethpage, Town of Oyster Bay, New York. Since its inception in 1941, the plant's primary mission was the research, prototyping, testing, design, engineering, fabrication, and primary assembly of military aircraft. The facilities at NWIRP included four plants used for assembly and prototype testing, a group of quality control laboratories, two warehouse complexes (north and south), a salvage storage area, water recharge basins, the Industrial Wastewater Treatment Plant, and several smaller support buildings.

The Navy's property originally totaled 109.5 acres and was formerly a Government-Owned Contractor-Operated (GOCO) facility that was operated by Northrop Grumman (NG) until September 1998. Prior to 2002, the NWIRP property was bordered on the north, west, and south by current or former NG facilities, and on the east by a residential neighborhood. By March 2008, approximately 100 acres of NWIRP property were transferred to Nassau County in three separate actions. The remaining 9 acres and access easements were retained by the Navy to continue remedial efforts at Installation Restoration (IR) Site 1 – Former Drum Marshalling Area and Site 4 – Former Underground Storage Tanks (Area of Concern [AOC] 22). A parcel of land connecting the two sites was also retained. Currently, the 9-acre parcel of NWIRP is bordered on the east by the residential neighborhood and on the north, south, and west by Steel Equities; however, a small portion is still owned by Nassau County. Access to the NWIRP is from South Oyster Bay Road.

1.3 Geology and Hydrogeology

Overburden at the site consists of well over 1,000 feet (ft) of unconsolidated deposits overlying crystalline bedrock of the Hartland Formation. Overburden is divided into four geologic units: the upper Pleistocene deposits, the Magothy Formation, the clay member of the Raritan Formation ("Raritan Clay") and the Lloyd Sand member of the Raritan Formation ("Lloyd Sand") (Geraghty and Miller, 1994).

The upper Pleistocene ranges in thickness from approximately 50 to 100 ft and consists of till and outwash deposits of medium to coarse sand and gravel with lenses of fine sand, silt and clay (Smolensky and Feldman, 1990); these deposits form the Upper Glacial Aquifer. Directly underlying this unit is the Magothy Formation with a thickness of 650 to 900 ft and lower extent of 700 to 1000 ft below ground surface (bgs) observed onsite. The Magothy is characterized by fine to medium sands and silts interbedded with zones of clays, silty sands and sandy clays. Sand and gravel lenses are found in some areas between depths of 600 and 880 ft bgs; these deposits form the main producing zones of the Magothy Aquifer.

Investigations performed by the Navy since 2012 indicate that the bottom of the Magothy (top of the Raritan Clay) can extend to depths of 700 to greater than 1,000 ft bgs. The top of the Raritan Clay deepens to the south-southeast, as evidenced by clay depths of 1,000 ft bgs (or more) in borings installed offsite. The Raritan Clay Unit is of continental origin and consists of clay, silty clay, clayey silt, and fine silty sand. This member acts as a confining layer over the Lloyd Sand Unit. The Lloyd Sand Unit is also of continental origin, having been deposited in a large fresh water lacustrine environment. The material consists of fine to coarse-grained sands, gravel, inter-bedded clay, and silty sand. These deposits form the Lloyd Aquifer.

The Upper Glacial Aquifer and the Magothy Aquifer comprise the aquifers of interest at the NWIRP. Regionally, these formations are generally considered to form a common, interconnected aquifer as the coarse nature of each unit near their contact and the lack of any regionally confining clay unit allows for the unrestricted flow of groundwater between the formations.

The Magothy Aquifer is the major source of public water in Nassau County. The most productive water bearing zones are the discontinuous lenses of sand and gravel that occur within the siltier matrix. The major water-bearing zones are coarse sand and gravel lenses located in the lower portion of the Magothy. The Magothy Aquifer is commonly regarded to function overall as an unconfined aquifer at shallow depths and a confined aquifer at deeper depths. The drilling program at the NWIRP has revealed that clay zones beneath the facility are common but laterally discontinuous. No confining clay units of facility-wide extent have been encountered. This is also the case for borings installed offsite.

Groundwater is encountered at a depth of approximately 50 ft bgs at the facility. Historically, because of pumping and recharge at the facility, groundwater depths have been measured to range from 40 to 60 ft bgs. The groundwater flow in the area is to the south-southeast.

2.0 FIELD PROGRAM

Three monitoring wells were installed in the vicinity of VPB156 between December 2014 and January 2015. Field investigation activities consisted of drilling, well installation, well development, sampling, soil/groundwater analysis, and surveying. Drilling during this investigation was performed by Delta Well and Pump Company of Ronkonkoma, New York. A description of these tasks is provided below.

2.1 Drilling and Well Construction

Monitoring wells RE122D1, RE122D2 and RE122D3 were installed using mud rotary drilling techniques (Figure 2). Depths of monitoring wells RE122D1, RE122D2 and RE122D3 were 545 ft, 615 ft and 740 ft, respectively. Well construction details are summarized in Table 1. Boring logs with lithologic descriptions of the well screen interval are included in the Appendix A. *Data Summary Report for VPB156* (Resolution Consultants, 2014) documents the installation of this VPB including detailed lithologic descriptions, continuous gamma plot and multiple VOC sample results over the entire boring length.

Prior to installing each monitoring well, the results of the groundwater samples, the geophysical logs, lithology and field data from the vertical profile borings were analyzed. Screen intervals were determined based on this analysis, intervals with the highest VOC concentrations as measured in the hydropunch samples, and coincident intervals with the highest apparent permeability based on the gamma logs. During the monitoring well installation, split spoon samples were collected every 5 ft in the screen interval. One soil sample per monitoring well was analyzed for Total Organic Carbon (TOC) via United States (US) Environmental Protection Agency (EPA) series SW-846 method 9060A by Katahdin Analytical Services (Katahdin). Data validation of TOC data was performed by Resolution Consultants. Data validation packages and analytical data tables are included in Appendix A.

Wells were constructed of 4-inch diameter, Schedule 80, National Sanitation Foundation-approved polyvinylchloride (PVC) riser pipe and .010-slot well screen. Wells were completed at the surface with a 12-inch diameter steel curb box. Well risers were set below grade and fit with lockable J plugs. Detailed monitoring well construction diagrams are included in Appendix A.

2.2 Well Development

Following installation, all monitoring wells were developed to evacuate silts and other fine-grained materials and to establish the filter pack to promote a hydraulic connection between the well and

the surrounding aquifer. Well development was not initiated until at least 24 hours after well installation.

Monitoring well screens were developed using a combination of air lifting, manual surging, and pumping with a submersible pump. Turbidity was monitored during development to determine stabilization. In compliance with New York State Department of Environmental Conservation (NYSDEC) policy, wells were developed until turbidity was less than 50 nephelometric turbidity units (NTUs) if possible. Table 2 summarizes total pumped volume from air and pump development and final turbidity. Well development logs are included in Appendix A.

2.3 Sampling

Following development, wells were allowed to stabilize for at least 2 weeks prior to groundwater sampling in accordance with low flow sampling procedures. Wells were purged using a bladder pump with a drop tube intake placed at the approximate midpoint of the screened interval. The following water quality parameters were continuously measured: water temperature, pH, conductivity, oxidation-reduction potential, dissolved oxygen and turbidity. Groundwater analytical samples were collected when water quality parameters stabilized. Samples were analyzed for VOCs via method 8260B and 1,4-dioxane via Method 8270C by Katahdin. All development and purge water was managed as investigation derived waste (IDW). Groundwater sample logs and data validation packages are included in Appendix A.

Monitoring wells RE122D1, RE122D2 and RE122D3 are sampled quarterly as part of the Navy's ongoing Environmental Restoration Program. Resolution Consultants sampled these three wells during the March 2015 quarterly monitoring event. Analytical results and stabilized field parameters for these data are summarized in Table 3 and 4, respectively. Data validation is documented in Appendix A.

2.4 Decontamination and Investigation Derived Waste (IDW)

Resolution Consultants utilized dedicated and disposable sampling equipment when possible to avoid the potential for cross-contamination of samples. The sampling equipment included dedicated plastic scoops, disposable Teflon or polyethylene tubing, disposable gloves, and laboratory supplied sample bottles. Hand held equipment and split spoons were decontaminated using Liquinox and water wash, a potable water rinse, followed by a distilled water rinse. Water was collected in 5-gallon pails or 55-gallon drums. Non dedicated sampling equipment was decontaminated as outlined in the *UFP SAP Addendum - Groundwater Sampling Using Low Stress (Low Flow) Purging and Sampling Protocol* (Resolution Consultants, 2013b).

As part of the IDW management practices and in accordance with the SAP, the investigation waste (consisting of soil cuttings, drilling muds, IDW fluids, and personal protective equipment [PPE]) generated during the groundwater monitoring well installation and sampling was containerized and staged at NWIRP Bethpage.

IDW solids were containerized in roll offs. Representative samples from each roll off were submitted to Katahdin for analysis of:

- Target Compound List (TCL) VOCs
- TCL Semi-volatile Organic Compounds (SVOCs)
- Toxicity Characteristic Leaching Procedure (TCLP) Metals
- Polychlorinated Biphenyls (PCBs)
- Total petroleum hydrocarbons
- Corrosivity
- Ignitability
- Reactive Cyanide
- Reactive Sulfide
- Paint Filter

IDW fluid generated during well development and purging was containerized in frac tanks and stored at NWIRP Bethpage for characterization and ultimate disposal to the Publicly Owned Treatment Works (POTW), in accordance with the facilities existing discharge permit. A representative water sample was collected from each frac tank and submitted to Katahdin for analysis of VOCs via Method SW 624, pH via Method SW 9040B, PCBs via Method 8082 and Total Metals via Method SW 846. All analytical criteria were met for disposal of water.

2.5 Surveying

A survey of the monitoring well locations was conducted at the end of fieldwork by C. T. Male, Inc., of Latham, NY, under the direct supervision of Resolution Consultants. The locations were tied into the existing base map developed for this investigation. The survey elevation is referenced to the North American Vertical Datum (NAVD) 1988 and has a vertical accuracy of 0.01 foot. Vertical control is based on observations of the Continuously Operating Reference (COR) Stations Queens and Central Islip. The horizontal location is referenced to the North American Datum (NAD) 1983

(2011) N.Y. Long Island Zone 3104 and has an accuracy of 0.1 foot. Local horizontal and vertical control is based on Global Positioning System (GPS) observations using the NYSNet Real Time Network.

A table of survey data (latitude/longitude, northing/easting, elevations of ground, rim and PVC) and a survey map is included in Appendix A.

3.0 REFERENCES

Geraghty and Miller, Inc., 1994. *Remedial Investigation Report, Grumman Aerospace Corporation, Bethpage, New York*. Revised September 1994.

Naval Facilities Engineering Command (NAVFAC), 2003. *Record of Decision Naval Weapons Industrial Reserve Plant Bethpage, New York, Operable Unit 2 – Groundwater*, NYS Registry: 1-30-003B. April 2003.

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Tables

TABLE 1
MONITORING WELL CONSTRUCTION SUMMARY
2014-2015 OU2 GROUNDWATER INVESTIGATION
NWIRP BETHPAGE, NY

| MONITORING WELL | WELL COMPLETION DATE | GROUND ELEVATION (MSL) | PVC ELEVATION (INNER CASING) (MSL) | WELL DEPTH (ft bgs) | CASING DEPTH (ft bgs) | SCREEN INTERVAL (ft bgs) | SUMP DEPTH INTERVAL (ft bgs) | BORING DEPTH (ft bgs) |
|-----------------|----------------------|------------------------|------------------------------------|---------------------|-----------------------|--------------------------|------------------------------|-----------------------|
| RE122D1 | 1/29/2015 | 97.74 | 97.42 | 545 | 53 | 520 - 540 | 540 - 545 | 557 |
| RE122D2 | 1/14/2015 | 97.70 | 97.35 | 615 | 53 | 590 - 610 | 610 - 615 | 627 |
| RE122D3 | 12/11/2014 | 97.62 | 97.27 | 740 | 53 | 715 - 735 | 735 - 740 | 752 |

MSL - mean sea level
ft bgs - feet below ground surface

TABLE 2
 MONITORING WELL DEVELOPMENT SUMMARY
 2014-2015 GROUNDWATER INVESTIGATION
 NWIRP BETHPAGE, NY

| MONITORING WELL | AIR DEVELOPMENT | | PUMP DEVELOPMENT | | | APPROX. TOTAL DEVELOPMENT VOLUME (GAL) | FINAL TURBIDITY (NTUs) |
|-----------------|-------------------|----------------------|-------------------|---------------------------|----------------------|--|------------------------|
| | DATE | APPROX. VOLUME (GAL) | DATE | FINAL PUMP DEPTH (FT BGS) | APPROX. VOLUME (GAL) | | |
| RE122D1 | 2/9/2015 | 4,500 | 2/13/2015 | 520-540 | 3,500 | 8,000 | 0.02 |
| RE122D2 | 2/10/2015 | 5,500 | 2/18/2015 | 590-610 | 4000 | 9,500 | 0.02 |
| RE122D3 | 2/10/15 & 2/11/15 | 7,000 | 2/19/15 & 2/23/15 | 715-735 | 7000 | 14,000 | 20.15 |

GAL - gallon

FT BGS - feet below ground surface

NTUs - Nephelometric Turbidity Units

Table 3. Analytical Data Summary

| Location | NYSDEC Groundwater Guidance or Standard Value (Note 1) | RE122D1 | RE122D2 | RE122D3 |
|---------------------------------------|--|--------------------|--------------------|--------------------|
| Sample Date | | 3/24/2015 | 3/24/2015 | 3/24/2015 |
| Sample ID | | RE122D1-GW-032415 | RE122D2-GW-032415 | RE122D3-GW-032415 |
| Sample type code | | N | N | N |
| VOC 8260C (ug/L) | | | | |
| 1,1,1-TRICHLOROETHANE | 5 | < 0.50 U | 0.71 J | < 0.50 U |
| 1,1,2,2-TETRACHLOROETHANE | 5 | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 5 | 7.2 | 31 | < 0.50 U |
| 1,1,2-TRICHLOROETHANE | 1 | 0.40 J | 2.8 | < 0.50 U |
| 1,1-DICHLOROETHANE | 5 | < 0.50 U | 1.7 | < 0.50 U |
| 1,1-DICHLOROETHENE | 5 | < 0.50 U | 8.7 | < 0.50 U |
| 1,2,4-TRICHLOROBENZENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,2-DIBROMO-3-CHLOROPROPANE | 0.04 | < 0.75 U | < 0.75 U | < 0.75 U |
| 1,2-DIBROMOETHANE | NL | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,2-DICHLOROBENZENE | 3 | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,2-DICHLOROETHANE | 5 | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,2-DICHLOROETHENE, TOTAL | 5 | 2.0 | 6.0 | < 1.0 U |
| 1,2-DICHLOROPROPANE | 1 | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,3-DICHLOROBENZENE | 3 | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,4-DICHLOROBENZENE | 3 | < 0.50 U | < 0.50 U | < 0.50 U |
| 1,4-DIOXANE (Method 8270D_SIM) | NL | 8.1 | 14 | < 0.19 U |
| 2-BUTANONE | 50 | < 2.5 UJ | < 2.5 UJ | < 2.5 UJ |
| 2-HEXANONE | 50 | < 2.5 UJ | < 2.5 UJ | < 2.5 UJ |
| 4-METHYL-2-PENTANONE | NL | < 2.5 U | < 2.5 U | < 2.5 U |
| ACETONE | 50 | < 2.5 UJ | < 2.5 UJ | < 2.5 UJ |
| BENZENE | 1 | < 0.50 U | < 0.50 U | < 0.50 U |
| BROMODICHLOROMETHANE | 50 | < 0.50 U | < 0.50 U | < 0.50 U |
| BROMOFORM | 50 | < 0.50 U | < 0.50 U | < 0.50 U |
| BROMOMETHANE | 5 | < 1.0 UJ | < 1.0 UJ | < 1.0 UJ |
| CARBON DISULFIDE | 60 | < 0.50 UJ | < 0.50 UJ | < 0.50 UJ |
| CARBON TETRACHLORIDE | 5 | < 0.50 U | 1.5 | < 0.50 U |
| CHLOROBENZENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U |
| CHLOROETHANE | 5 | < 1.0 UJ | < 1.0 UJ | < 1.0 U |
| CHLOROFORM | 7 | 0.62 J | 2.4 | < 0.50 U |
| CHLOROMETHANE | 5 | < 1.0 UJ | < 1.0 UJ | < 1.0 UJ |
| CIS-1,2-DICHLOROETHENE | 5 | 2.0 | 6.0 | < 0.50 U |
| CIS-1,3-DICHLOROPROPENE | 0.4 | < 0.50 U | < 0.50 U | < 0.50 U |
| CYCLOHEXANE | NL | < 0.50 UJ | < 0.50 UJ | < 0.50 UJ |
| DIBROMOCHLOROMETHANE | 5 | < 0.50 U | < 0.50 U | < 0.50 U |
| DICHLORODIFLUOROMETHANE | 5 | < 1.0 UJ | < 1.0 UJ | < 1.0 UJ |
| ETHYLBENZENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U |
| ISOPROPYLBENZENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U |
| M- AND P-XYLENE | NL | < 1.0 U | < 1.0 U | < 1.0 U |
| METHYL ACETATE | NL | < 0.75 U | < 0.75 U | < 0.75 U |
| METHYL CYCLOHEXANE | NL | < 0.50 U | < 0.50 U | < 0.50 UJ |
| METHYL TERT-BUTYL ETHER | 10 | < 0.50 U | < 0.50 U | < 0.50 U |
| METHYLENE CHLORIDE | 5 | < 2.5 U | < 2.5 U | < 2.5 U |
| O-XYLENE | NL | < 0.50 U | < 0.50 U | < 0.50 U |
| STYRENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U |
| TETRACHLOROETHENE | 5 | 1.3 J | 2.7 J | < 0.50 UJ |
| TOLUENE | 5 | < 0.50 U | 0.63 J | 0.37 J |
| TRANS-1,2-DICHLOROETHENE | 5 | < 0.50 U | < 0.50 U | < 0.50 U |
| TRANS-1,3-DICHLOROPROPENE | 0.4 | < 0.50 U | < 0.50 U | < 0.50 U |
| TRICHLOROETHENE | 5 | 570 J | 4600 J | 6.8 |
| TRICHLOROFLUOROMETHANE | 5 | < 1.0 U | < 1.0 U | < 1.0 U |
| VINYL CHLORIDE | 2 | < 1.0 U | < 1.0 U | < 1.0 U |
| XYLENES, TOTAL | 5 | < 1.5 U | < 1.5 U | < 1.5 U |

Notes:

1 New York State Department of Environmental Conservation Division of Water Technical and Operation Guidance series
(6 NYCRR 700-706, Part 703.5 summarized in TOGS 1.1.1)

Ambient water quality standards and groundwater effluent limitations, class GA; NL = Not Listed

Bold = Detected; **Bold and Italics** = Not detected exceeds NYS Groundwater Standards or guidance value

Yellow highlighted values exceed Groundwater Standards or guidance value

Sample type codes: N - normal environmental sample, FD - field duplicate

U = Nondetected result. The analyte was analyzed for, but was not detected above the reported sample quantitation limit.

UJ = The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte.

J = The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.

M = the matrix spike or matrix spike duplicate did not meet recovery or precision requirements.

Table 4
Stabilized Field Parameters

| Well | Date | Temperature (°C) | pH | Specific Conductance (µS/cm) | DO (mg/L) | ORP (mV) | Turbidity (NTU) | Depth to water (ft bgs) | Flow rate (ml/min) |
|---------|-----------|------------------|------|------------------------------|-----------|----------|-----------------|-------------------------|--------------------|
| RE122D1 | 3/24/2015 | 12.90 | 6.5 | 0.153 | 6.35 | 144.7 | 25.8 | 39.27 | 500 |
| RE122D2 | 3/24/2015 | 12.81 | 5.65 | 0.121 | 5.09 | -15.3 | 3.86 | 39.58 | 500 |
| RE122D3 | 3/24/2015 | 12.52 | 5.55 | 0.051 | 2.82 | 200.7 | 95.4 | 40.7 | 500 |

°C - degrees Celsius

µS/cm - Microsiemens per Centimeter

mg/L - milligrams per liter

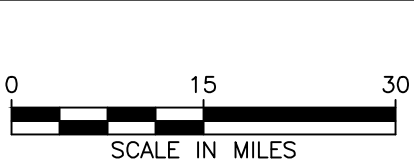
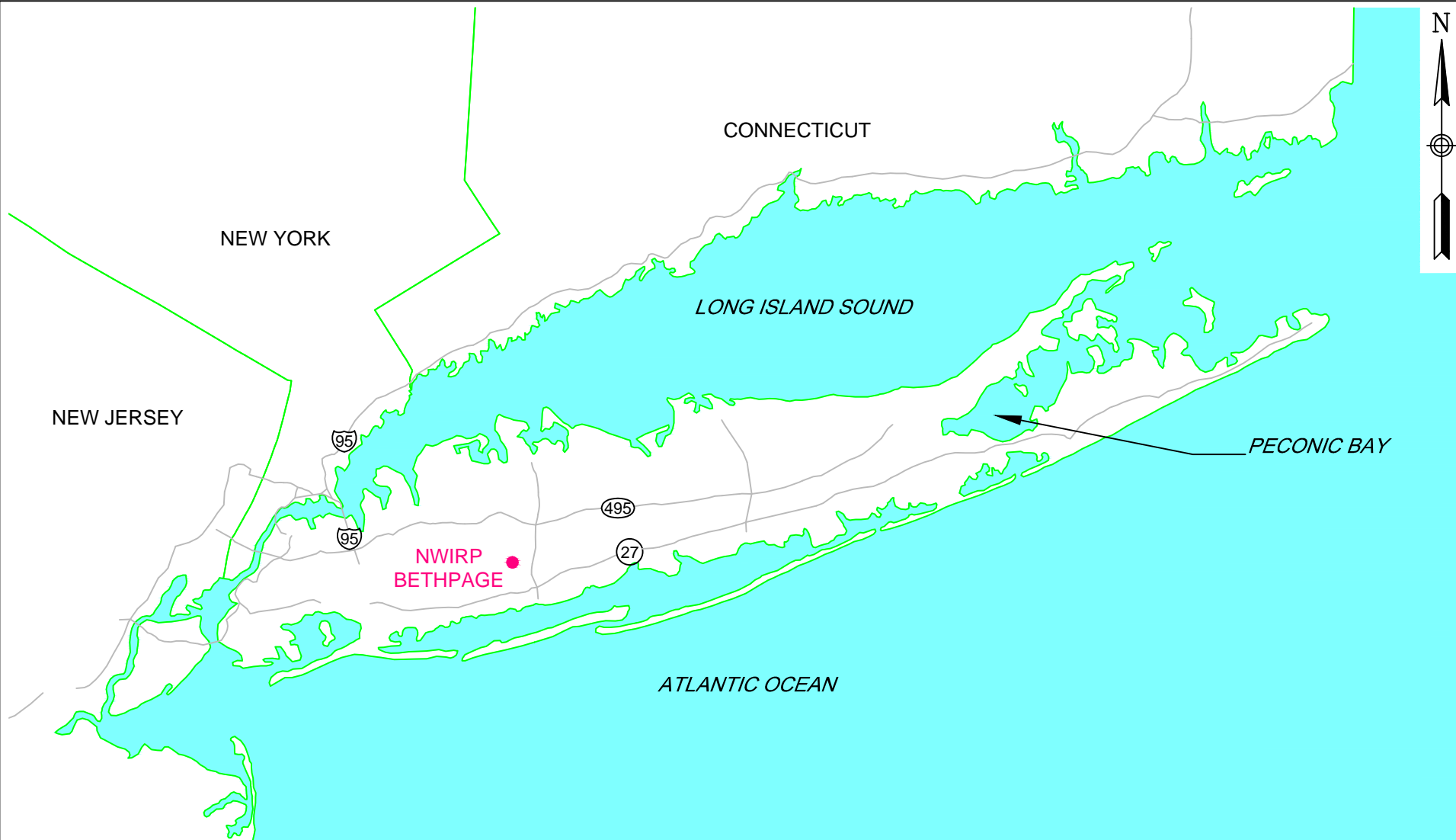
mV - Millivolts

NTU - Nephelometric Turbidity Unit

ft bgs - feet below ground surface

ml/min - milliliters per minute

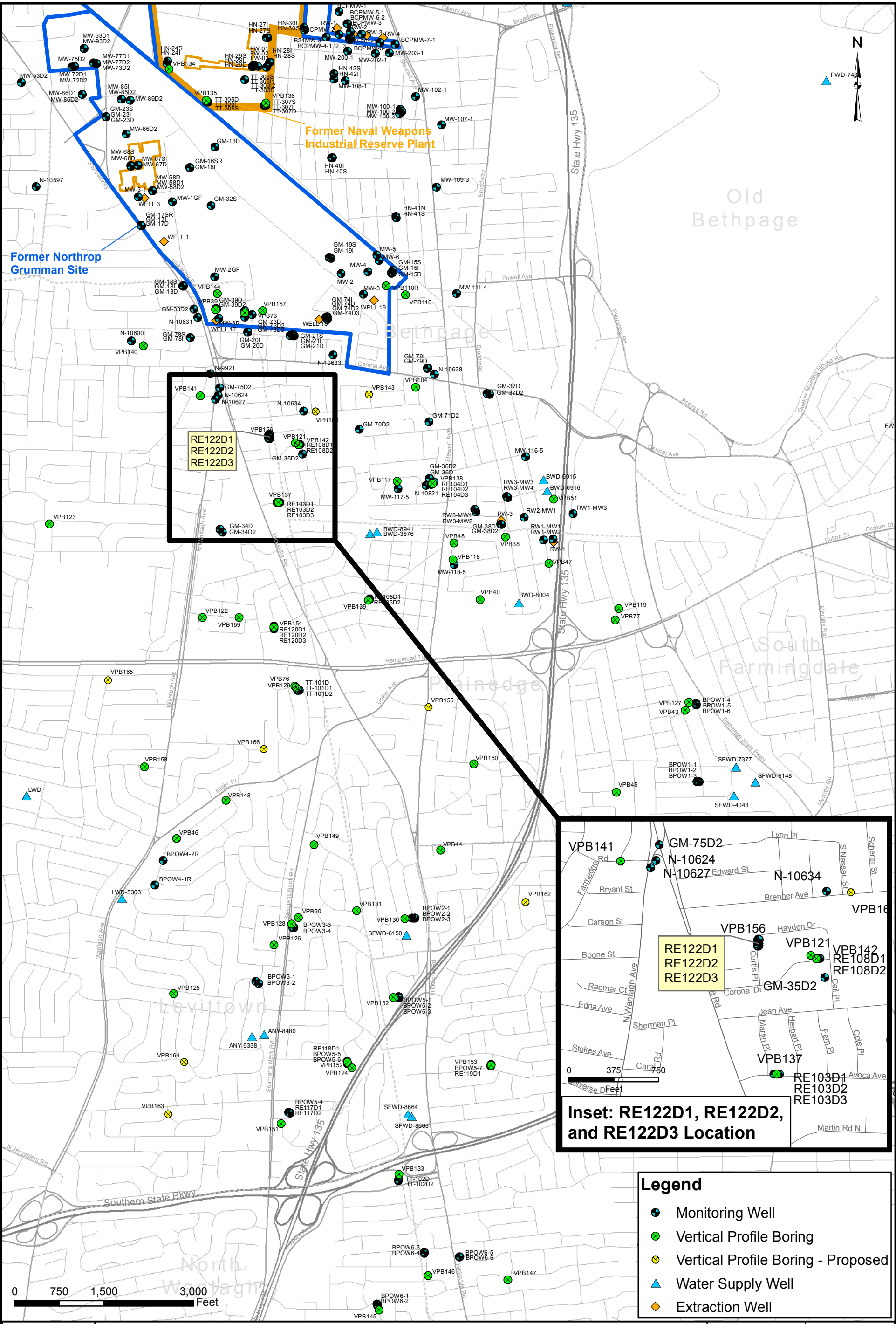
Figures



GENERAL LOCATION MAP
NWIRP BETHPAGE
BETHPAGE, NEW YORK

| | | | |
|-------------------------------------|--|--------------------|----------|
| CONTRACT NUMBER N62470-11-D-8013 | | CTO NUMBER WE15 | |
| APPROVED BY --- | | DATE --- | |
| APPROVED BY --- | | DATE --- | |
| FIGURE NO. 1 | | | REV 0 |

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RE122D1
RE122D2
RE122D3

Inset: RE122D1, RE122D2, and RE122D3 Location

VPB141, GM-75D2, N-10624, N-10627, N-10634, VPB162, VPB156, VPB121, VPB142, RE108D1, RE108D2, GM-35D2, VPB137, RE103D1, RE103D2, RE103D3

| Legend | |
|--------|------------------------------------|
| | Monitoring Well |
| | Vertical Profile Boring |
| | Vertical Profile Boring - Proposed |
| | Water Supply Well |
| | Extraction Well |



RE122D1, RE122D2, AND RE122D3 LOCATION MAP
NAVAL WEAPONS INDUSTRIAL RESERVE PLANT
BETHPAGE, NEW YORK

| | |
|------------------------------------|--------------------|
| CONTRACT NUMBER N62470-11-D8013 | CTO NUMBER WE15 |
| APPROVED BY PS | DATE 8/18/2015 |
| APPROVED BY | DATE |
| FIGURE NO. 2 | REV 0 |

Appendix A
RE122D1, RE122D2, RE122D3

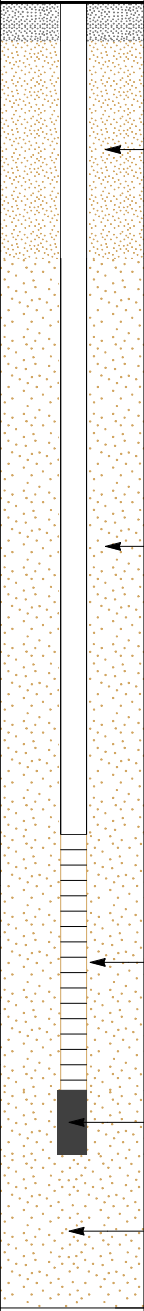

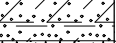

Section 1

Boring Logs

| | | |
|---|---|--|
| Client: Department of the Navy, Naval Facilities Engineering Command, Mid-Atlantic | | Logged By: V. Varricchio |
| Location: Curtis Pl and Hayden Dr, Bethpage, NY | | Drilling Company: DELTA WELL AND PUMP COMPANY |
| Project #: 60266526 | Ground Elevation (msl): 97.74 | Well Screen Interval (ft): 520-540 |
| Start Date: 1/21/2015 | Drilling Method: Auger (0-50' bgs) Mud Rotary (>50' bgs) | Water Level (ft): |
| Finish Date: 1/29/2015 | Northing: 207818.19 Easting: 1124982.02 | Total Depth (ft): 557.0 |

| DEPTH (ft) | PID (ppm) | Formation | USCS | GRAPHIC LOG | MATERIAL DESCRIPTION | Well Completion | Well Construction |
|------------|-----------|-----------|------|-------------|---|-----------------|-----------------------------------|
| 0 | | | | | 0-523 ft bgs: See VPB156 for Descriptions | | |
| 50 | | | | | | | 10" Diameter Steel Casing |
| 100 | | | | | | | |
| 150 | | | | | | | |
| 200 | | | | | | | |
| 250 | | | | | | | Bentonite Grout |
| 300 | | | | | | | |
| 350 | | | | | | | |
| 400 | | | | | | | |
| 450 | | | | | | | 4" Diameter Schedule 80 PVC Riser |

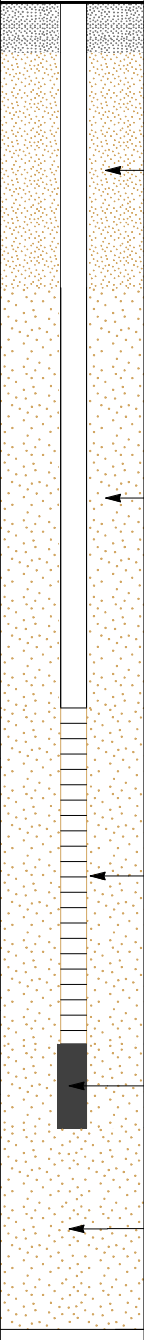


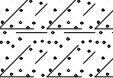
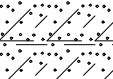
| | | |
|---|---|--|
| Client: Department of the Navy, Naval Facilities Engineering Command, Mid-Atlantic | | Logged By: V. Varricchio |
| Location: Curtis Pl and Hayden Dr, Bethpage, NY | | Drilling Company: DELTA WELL AND PUMP COMPANY |
| Project #: 60266526 | Ground Elevation (msl): 97.74 | Well Screen Interval (ft): 520-540 |
| Start Date: 1/21/2015 | Drilling Method: Auger (0-50' bgs) Mud Rotary (>50' bgs) | Water Level (ft): |
| Finish Date: 1/29/2015 | Northing: 207818.19 Easting: 1124982.02 | Total Depth (ft): 557.0 |

| DEPTH (ft) | PID (ppm) | Formation | USCS | GRAPHIC LOG | MATERIAL DESCRIPTION | Well Completion | Well Construction | | |
|---------------------------------|-----------|-----------|------|---|---|--|--|----------------|--|
| 456 | | | | | 0-523 ft bgs: See VPB156 for Descriptions (<i>continued</i>) |  | 4" Diameter Schedule 80 PVC Riser (<i>continued</i>) | | |
| 458 | | | | | | | | #0 Filter Sand | |
| 460 | | | | | | | | | |
| 462 | | | | | | | | | |
| 464 | | | | | | | | | |
| 466 | | | | | | | | | |
| 468 | | | | | | | | | |
| 470 | | | | | | | | | |
| 472 | | | | | | | | | |
| 474 | | | | | | | | | |
| 476 | | | | | | | | | |
| 478 | | | | | | | | | |
| 480 | | | | | | | | | |
| 482 | | | | | | | | | |
| 484 | | | | | | | | | |
| 486 | | | | | | | | | |
| 488 | | | | | | | | | |
| 490 | | | | | | | | | |
| 492 | | | | | | | | | |
| 494 | | | | | | | | | |
| 496 | | | | | | | | | |
| 498 | | | | | | | | | |
| 500 | | | | | | | | | |
| 502 | | | | | | | | | |
| 504 | | | | | | | | | |
| 506 | | | | | | | | | |
| 508 | | | | | | | | | |
| 510 | | | | | | | | | |
| 512 | | | | | | | | | |
| 514 | | | | | | | | | |
| 516 | | | | | | | | | |
| 518 | | | | | | | | | |
| 520 | | | | | | | | | |
| 522 | | | | | | | | | |
| 524 | 0.0 | | CH |  | Light grey (10 YR 7/2) fat CLAY, trace Silt | | | | |
| 526 | | | SW |  | Light brown (10 YR 6/2) fine to medium well-graded SAND, trace Silt | | | | |
| 528 | | | | | White (10 YR 8/1) poorly graded fine SAND | | | | |
| 530 | 0.0 | | SP |  | | | | | |
| 532 | | | | | Yellowish brown (10 YR 5/8) poorly graded fine SAND | | | | |
| 534 | 0.0 | | SP | | | | | | |
| 536 | | | | | White (10 YR 8/1) poorly graded fine SAND | | | | |
| 538 | | | | | White (10 YR 8/1) poorly graded fine SAND, iron banding | | | | |
| 540 | 0.0 | | SP | | | | | | |
| 542 | | | | | | | | | |
| 544 | | | | | | | Sump | | |
| 546 | | | | | | | | | |
| 548 | | | | | | | | | |
| 550 | | | | | | | | | |
| 552 | | | | | | | | | |
| 554 | | | | | | | | | |
| 556 | | | | | | | #1 Sand to bottom | | |
| End of boring at 557.0 ft. bgs. | | | | | | | | | |

| | | | | | |
|---|--|---|--|---|--|
| Client: Department of the Navy, Naval Facilities Engineering Command, Mid-Atlantic | | | Logged By: V. Varricchio | | |
| Location: Curtis Pl and Hayden Dr, Bethpage, NY | | | Drilling Company: DELTA WELL AND PUMP COMPANY | | |
| Project #: 60266526 | | Ground Elevation (msl): 97.7 | | Well Screen Interval (ft): 590-610 | |
| Start Date: 1/5/2015 | | Drilling Method: Auger (0-50' bgs) Mud Rotary (>50' bgs) | | Water Level (ft): | |
| Finish Date: 1/14/2015 | | Northing: 207789.07 Easting: 1124979.19 | | Total Depth (ft): 627.0 | |

| DEPTH (ft) | PID (ppm) | Formation | USCS | GRAPHIC LOG | MATERIAL DESCRIPTION | Well Completion | Well Construction |
|------------|-----------|-----------|------|-------------|--|-----------------|-----------------------------------|
| 0 | | | | | 0-593 ft bgs; See VPB 156 for Descriptions | | |
| 50 | | | | | | | 10" Diameter Steel Casing |
| 100 | | | | | | | |
| 150 | | | | | | | |
| 200 | | | | | | | |
| 250 | | | | | | | Bentonite Grout |
| 300 | | | | | | | |
| 350 | | | | | | | |
| 400 | | | | | | | |
| 450 | | | | | | | |
| 500 | | | | | | | 4" Diameter Schedule 80 PVC Riser |

| | | |
|---|---|--|
| Client: Department of the Navy, Naval Facilities Engineering Command, Mid-Atlantic | | Logged By: V. Varricchio |
| Location: Curtis Pl and Hayden Dr, Bethpage, NY | | Drilling Company: DELTA WELL AND PUMP COMPANY |
| Project #: 60266526 | Ground Elevation (msl): 97.7 | Well Screen Interval (ft): 590-610 |
| Start Date: 1/5/2015 | Drilling Method: Auger (0-50' bgs) Mud Rotary (>50' bgs) | Water Level (ft): |
| Finish Date: 1/14/2015 | Northing: 207789.07 Easting: 1124979.19 | Total Depth (ft): 627.0 |

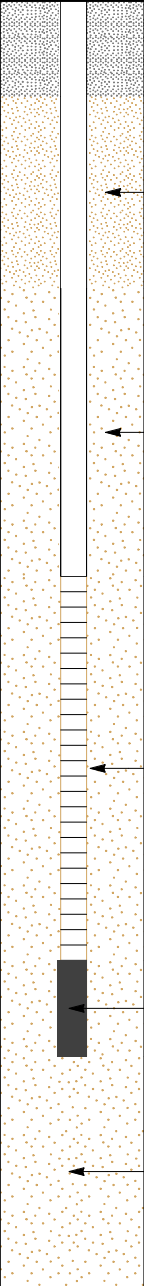
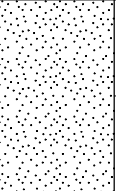
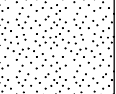
| DEPTH (ft) | PID (ppm) | Formation | USCS | GRAPHIC LOG | MATERIAL DESCRIPTION | Well Completion | Well Construction |
|------------|-----------|-----------|------|---|--|--|--|
| 548 | | | | | 0-593 ft bgs; See VPB 156 for Descriptions (<i>continued</i>) |  | 4" Diameter Schedule 80 PVC Riser (<i>continued</i>) |
| 550 | | | | | | | |
| 552 | | | | | | | |
| 554 | | | | | | | |
| 556 | | | | | | | |
| 558 | | | | | | | |
| 560 | | | | | | | |
| 562 | | | | | | | |
| 564 | | | | | | | |
| 566 | | | | | | | |
| 568 | | | | | | | |
| 570 | | | | | | | |
| 572 | | | | | | | |
| 574 | | | | | | | |
| 576 | | | | | | | |
| 578 | | | | | | | |
| 580 | | | | | | | |
| 582 | | | | | | | |
| 584 | | | | | | | |
| 586 | | | | | | | |
| 588 | | | | | | | |
| 590 | | | | | | | |
| 592 | | | | | | | |
| 594 | 0 | | SM |  | Light brownish grey (10 YR 6/2) silty fine SAND, lean Clay, trace iron/lignite banding | | #0 Filter Sand |
| 596 | | | | | | | |
| 598 | 0 | | SM |  | Light brownish grey (10 YR 6/2) silty fine SAND, trace lean Clay, iron/lignite banding | | #0 Filter Sand |
| 600 | | | | | | | |
| 602 | 0 | | SW |  | Grey (10 YR 6/1) fine to medium SAND, trace Silt and subrounded coarse sand | | #1 Filter Sand |
| 604 | | | | | | | |
| 606 | 0 | | SW |  | Grey (10 YR 6/1) fine to medium SAND, trace Silt | | #1 Filter Sand |
| 608 | | | | | | | |
| 610 | | | | | | | |
| 612 | | | | | | | |
| 614 | | | | | | | Sump |
| 616 | | | | | | | |
| 618 | | | | | | | |
| 620 | | | | | | | |
| 622 | | | | | | | #1 Sand to bottom |
| 624 | | | | | | | |
| 626 | | | | | | | |

End of boring at 627.0 ft. bgs.

| | | | | | |
|---|--|---|--|---|--|
| Client: Department of the Navy, Naval Facilities Engineering Command, Mid-Atlantic | | | Logged By: V. Varricchio | | |
| Location: Curtis Pl and Hayden Dr, Bethpage, NY | | | Drilling Company: DELTA WELL AND PUMP COMPANY | | |
| Project #: 60266526 | | Ground Elevation (msl): 97.62 | | Well Screen Interval (ft): 715-735 | |
| Start Date: 12/2/2014 | | Drilling Method: Auger (0-50' bgs) Mud Rotary (>50' bgs) | | Water Level (ft): | |
| Finish Date: 12/11/2014 | | Northing: 207774.53 Easting: 1124980.64 | | Total Depth (ft): 752.0 | |

| DEPTH (ft) | PID (ppm) | Formation | USCS | GRAPHIC LOG | MATERIAL DESCRIPTION | Well Completion | Well Construction |
|------------|-----------|-----------|------|-------------|---|-----------------|-----------------------------------|
| 0 | | | | | 0-718 ft bgs: See VPB156 for Descriptions | | |
| 50 | | | | | | | 10" Diameter Steel Casing |
| 100 | | | | | | | |
| 150 | | | | | | | |
| 200 | | | | | | | |
| 250 | | | | | | | Bentonite Grout |
| 300 | | | | | | | |
| 350 | | | | | | | |
| 400 | | | | | | | |
| 450 | | | | | | | |
| 500 | | | | | | | |
| 550 | | | | | | | 4" Diameter Schedule 80 PVC Riser |
| 600 | | | | | | | |
| 650 | | | | | | | |

| | | |
|---|---|--|
| Client: Department of the Navy, Naval Facilities Engineering Command, Mid-Atlantic | | Logged By: V. Varricchio |
| Location: Curtis Pl and Hayden Dr, Bethpage, NY | | Drilling Company: DELTA WELL AND PUMP COMPANY |
| Project #: 60266526 | Ground Elevation (msl): 97.62 | Well Screen Interval (ft): 715-735 |
| Start Date: 12/2/2014 | Drilling Method: Auger (0-50' bgs) Mud Rotary (>50' bgs) | Water Level (ft): |
| Finish Date: 12/11/2014 | Northing: 207774.53 Easting: 1124980.64 | Total Depth (ft): 752.0 |

| DEPTH (ft) | PID (ppm) | Formation | USCS | GRAPHIC LOG | MATERIAL DESCRIPTION | Well Completion | Well Construction |
|------------|-----------|-----------|------|---|--|--|--|
| 686 | | | | | 0-718 ft bgs: See VPB156 for Descriptions (<i>continued</i>) |  | 4" Diameter Schedule 80 PVC Riser (<i>continued</i>) |
| 688 | | | | | | | |
| 690 | | | | | | | |
| 692 | | | | | | | |
| 694 | | | | | | | |
| 696 | | | | | | | |
| 698 | | | | | | | |
| 700 | | | | | | | |
| 702 | | | | | | | |
| 704 | | | | | | | |
| 706 | | | | | | | |
| 708 | | | | | | | |
| 710 | | | | | | | #0 Filter Sand |
| 712 | | | | | | | |
| 714 | | | | | | | |
| 716 | | | | | | | |
| 718 | 0 | | |  | Poorly graded medium SAND | | |
| 720 | | | SP | | | | |
| 722 | | | | | | | |
| 724 | 0 | | |  | Poorly graded medium SAND, Iron banding ~6-7" | | |
| 726 | | | SP | | | | |
| 728 | | | | | | | |
| 730 | | | | | | | |
| 732 | | | | | | | |
| 734 | | | | | | | |
| 736 | | | | | | | |
| 738 | | | | | | | Sump |
| 740 | | | | | | | |
| 742 | | | | | | | |
| 744 | | | | | | | |
| 746 | | | | | | | |
| 748 | | | | | | | |
| 750 | | | | | | | #1 Sand to bottom |
| 752 | | | | | End of boring at 752.0 ft. bgs. | | |

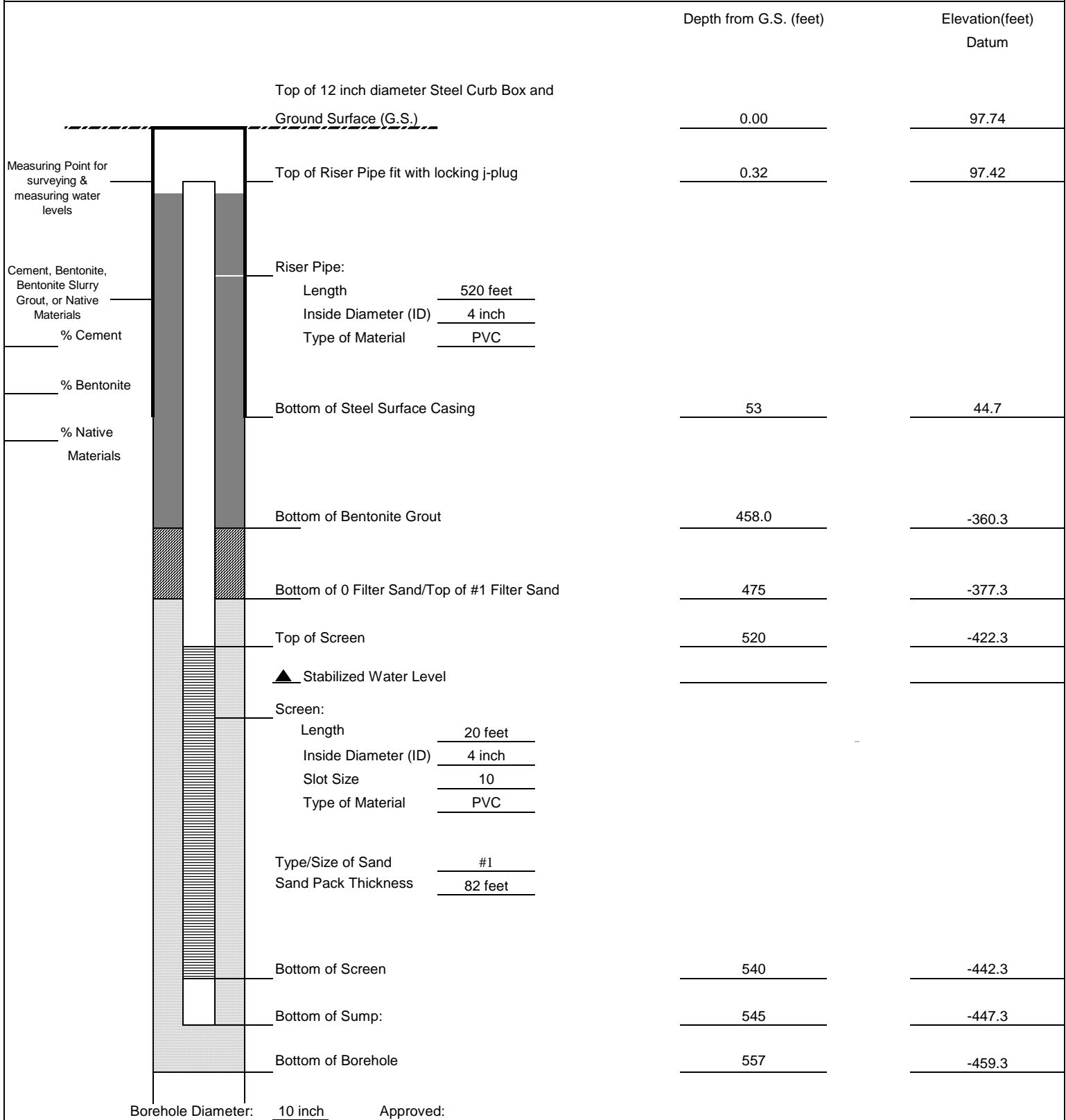
Section 2

Monitoring Well Construction Logs



| | | |
|--|--------------------------|---------------------------------------|
| Client: NAVFAC | Project Number: 60266526 | WELL ID: RE122D1 |
| Site Location: NWIRP BETHPAGE, NY | | |
| Well Location: Curtis Pl. and Hayden Dr., Bethpage | | Date Installed: 1/21/2015 - 1/29/2015 |
| Method: MUD ROTARY | | Inspector: V. VARRICCHIO |
| Coords: Northing: 207818.1929 Easting: 1124982.01 | | Contractor: DELTA WELL & PUMP |

MONITORING WELL CONSTRUCTION DETAIL



Borehole Diameter: 10 inch Approved: _____

Describe Measuring Point:

Signature _____

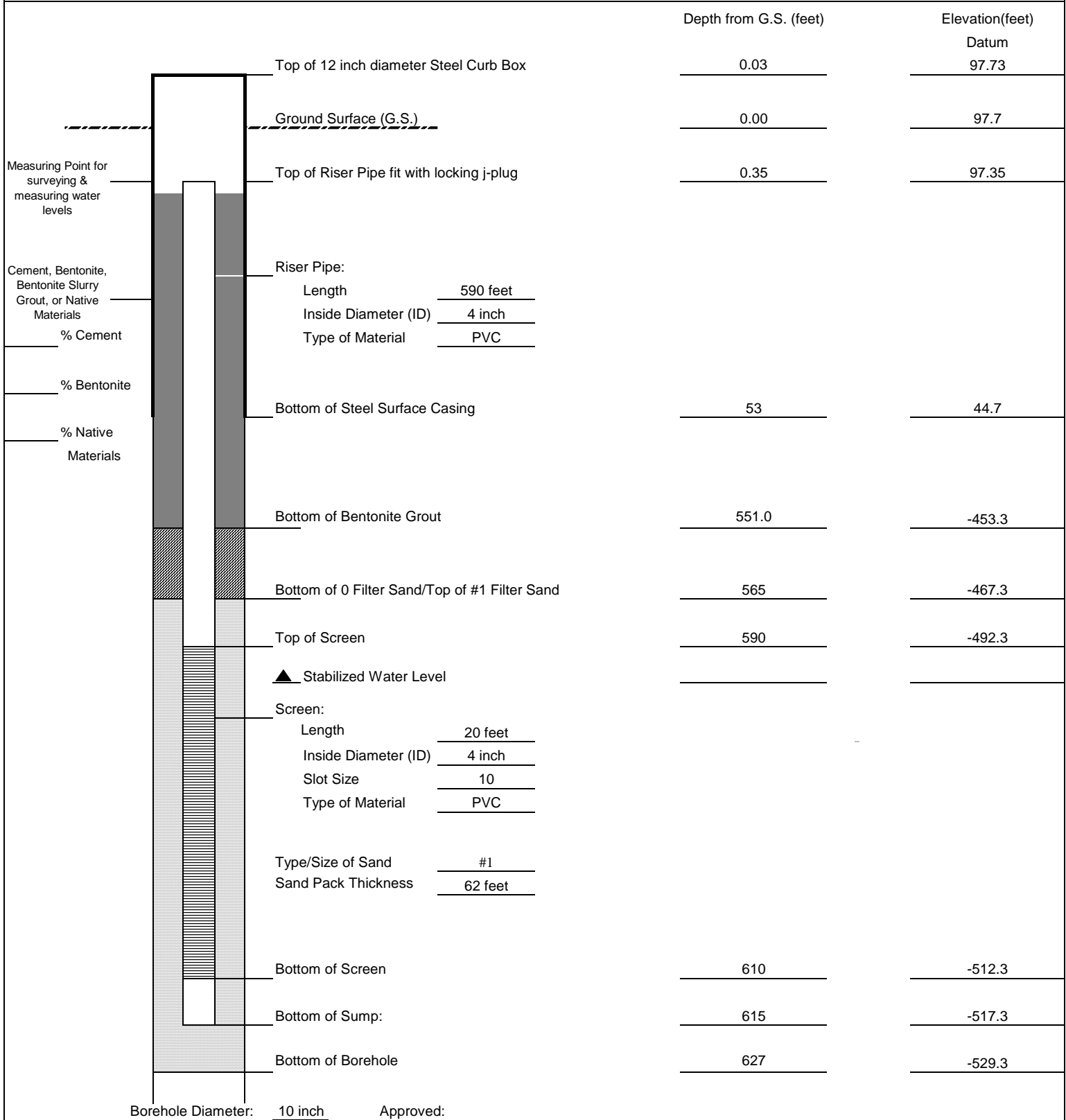
Date _____

Ground Surface _____



| | | |
|--|--------------------------|--------------------------------------|
| Client: NAVFAC | Project Number: 60266526 | WELL ID: RE122D2 |
| Site Location: NWIRP BETHPAGE, NY | | |
| Well Location: Curtis Pl. and Hayden Dr., Bethpage | | Date Installed: 1/5/2015 - 1/14/2015 |
| Method: MUD ROTARY | | Inspector: V. VARRICCHIO |
| Coords: Northing: 207789.07 Easting: 1124979.19 | | Contractor: DELTA WELL & PUMP |

MONITORING WELL CONSTRUCTION DETAIL



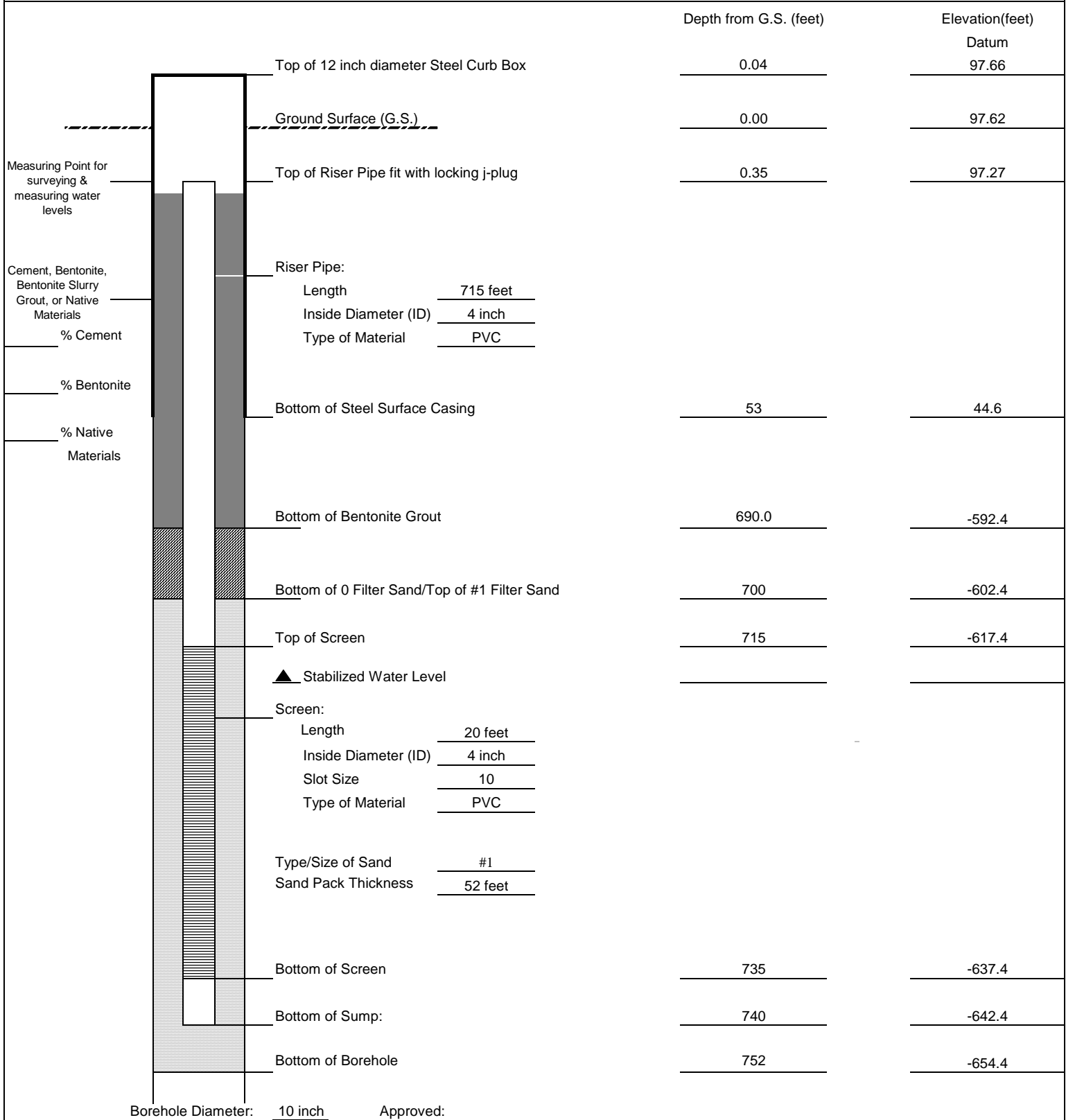
Describe Measuring Point: _____
 Ground Surface _____

Approved: _____
 Signature _____ Date _____



| | | |
|--|--------------------------|--|
| Client: NAVFAC | Project Number: 60266526 | WELL ID: RE122D3 |
| Site Location: NWIRP BETHPAGE, NY | | |
| Well Location: Curtis Pl. and Hayden Dr., Bethpage | | Date Installed: 12/2/2014 - 12/11/2014 |
| Method: MUD ROTARY | | Inspector: V. VARRICCHIO |
| Coords: Northing: 207774.52 Easting: 1124980.63 | | Contractor: DELTA WELL & PUMP |

MONITORING WELL CONSTRUCTION DETAIL



Describe Measuring Point: _____
 Ground Surface _____
 Signature _____ Date _____

Section 3

Groundwater Sample Log Sheets



RESOLUTION CONSULTANTS

Well ID: RE12201

Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 3/24/15 Time: Start 8:15 am/pm
 Project No: 60266526 Finish 12:15 am/pm
 Site Location: Curtis & Hayden
 Weather Conds: Sunny 25° Collector(s): JC

1. WATER LEVEL DATA: (measured from Top of Casing)

a. Total Well Length 575 c. Length of Water Column _____ (a-b) Casing Diameter/Material 4-inch PVC
 b. Water Table Depth 39.26 d. Calculated System Volume (see back) 20ft screen 13.1gal

2. WELL PURGE DATA

a. Purge Method: Geotech bladder pump with drop tube assembly
 b. Acceptance Criteria defined (see workplan)
 - Temperature ± 3% - D.O. ± 10% (values >0.5 mg/L) Turbidity ± 10%
 - pH ± 0.1 unit - ORP ± 10mV
 - Sp. Cond. ± 3% - Drawdown < 0.3' Remove a minimum 1 screen volume
 c. Field Testing Equipment used:

| Make | Model | Serial Number |
|--------------|-----------------|---------------|
| <u>VSI</u> | <u>556 MP3</u> | <u>55474</u> |
| <u>Hanna</u> | <u>HI 98703</u> | <u>69177</u> |

| Time (24hr) | Volume Removed (Liters) | Temp. (°C) | pH | Spec. Cond. (mS/cm) | DO (mg/L) | ORP (mV) | Turbidity (NTU) | Flow Rate (ml/min) | Depth to water (ft) | Color/Odor |
|-------------|-------------------------|--------------|-------------|---------------------|--------------|--------------|-----------------|--------------------|---------------------|---------------|
| <u>0900</u> | | <u>10.33</u> | <u>7.76</u> | <u>0.158</u> | <u>11.10</u> | <u>190.4</u> | <u>2.00</u> | <u>400</u> | <u>39.35</u> | <u>clear</u> |
| <u>0910</u> | | <u>11.42</u> | <u>6.08</u> | <u>0.153</u> | <u>10.54</u> | <u>172.1</u> | | <u>400</u> | <u>39.45</u> | |
| <u>0915</u> | | <u>12.20</u> | <u>5.79</u> | <u>0.153</u> | <u>9.78</u> | <u>168.2</u> | <u>9.83</u> | <u>500</u> | <u>39.45</u> | <u>cloudy</u> |
| <u>0920</u> | | <u>12.26</u> | <u>5.79</u> | <u>0.153</u> | <u>9.80</u> | <u>168.1</u> | | <u>500</u> | <u>39.45</u> | |
| <u>0925</u> | | <u>12.18</u> | <u>6.13</u> | <u>0.235</u> | <u>9.16</u> | <u>165.8</u> | <u>>50</u> | <u>500</u> | <u>39.25</u> | |
| <u>0930</u> | | <u>12.34</u> | <u>6.76</u> | <u>0.242</u> | <u>8.69</u> | <u>168.3</u> | | <u>500</u> | <u>39.25</u> | |
| <u>0935</u> | | <u>12.37</u> | <u>6.75</u> | <u>0.241</u> | <u>8.76</u> | <u>167.2</u> | <u>>50</u> | <u>500</u> | <u>39.25</u> | |

d. Acceptance criteria pass/fail

| | | | |
|-------------------------------------|-------------------------------------|--------------------------|--------------------------|
| | Yes | No | N/A |
| Has required volume been removed | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Has required turbidity been reached | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Have parameters stabilized | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |

If no or N/A - Explain below.

(continued on back)

3. SAMPLE COLLECTION:

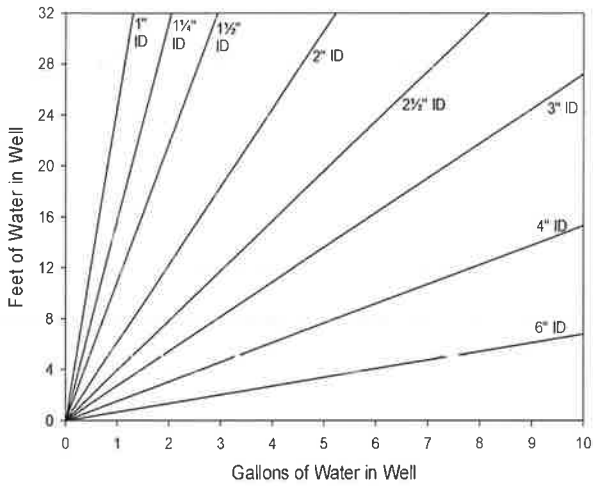
Method: Geotech bladder pump with drop tube assembly

| Sample ID | Container Type | No. of Containers | Preservation | Analysis Req. | Time |
|--------------------------|-------------------|-------------------|--------------|--------------------|-------------|
| <u>RE12201-GW-032415</u> | <u>40-mL vial</u> | <u>3</u> | <u>HCl</u> | <u>VOCs</u> | <u>1105</u> |
| <u>RE12201-GW-032415</u> | <u>1-L amber</u> | <u>2</u> | <u>none</u> | <u>1,4-Dioxane</u> | <u>1105</u> |

Comments hit bottom ~15ft to long ~535 of tubing

Signature _____ Date 3/24/15

Purge Volume Calculation



| Volume / Linear Ft. of Pipe | | |
|-----------------------------|--------|--------|
| ID (in) | Gallon | Liter |
| 0.25 | 0.0025 | 0.0097 |
| 0.375 | 0.0057 | 0.0217 |
| 0.5 | 0.0102 | 0.0386 |
| 0.75 | 0.0229 | 0.0869 |
| 1 | 0.0408 | 0.1544 |
| 1.25 | 0.0637 | 0.2413 |
| 1.5 | 0.0918 | 0.3475 |
| 2 | 0.1632 | 0.6178 |
| 2.5 | 0.2550 | 0.9653 |
| 3 | 0.3672 | 1.3900 |
| 4 | 0.6528 | 2.4711 |
| 6 | 1.4688 | 5.5600 |

1 screen volume
 15 ft = ~~5.56~~ / 9.8 G
 20 ft = ~~7.41~~ / 13.1 G
 25 ft = ~~9.27~~ / 16.3 G

Well ID:

(continued from front)

| Time (24 hr) | Volume | Temp (°C) | pH | Spec. Cond. (mS/cm) | DO (mg/L) | ORP (mV) | Turbidity (NTU) | Flow Rate (ml/min) | Depth to water (ft) | Color/Odor |
|-----------------|---------------------|--------------|------|------------------------|--------------|-------------|--------------------|-----------------------|------------------------|----------------|
| | Removed (Liters) | | | | | | | | | |
| 0740 | 56 | 12.14 | 6.74 | 0.239 | 8.73 | 163.2 | | 500 | 39.27 | cloudy |
| 1000 | | 12.24 | 6.61 | 0.157 | 7.84 | 170.4 | | 500 | 39.27 | |
| 1010 | | 12.74 | 6.53 | 0.157 | 6.91 | 173.7 | | 500 | 39.27 | dusty clearing |
| 1015 | | 12.73 | 6.51 | 0.155 | 6.54 | 172.5 | 234 | 500 | 39.27 | " |
| 1020 | | 12.71 | 6.51 | 0.153 | 6.57 | 161.4 | 225 | 520 | 39.27 | " |
| 1025 | | 12.69 | 6.51 | 0.155 | 6.49 | 152.8 | 223 | 500 | 39.27 | " |
| 1030 | 104 | 12.84 | 6.49 | 0.154 | 6.54 | 148.1 | 44.4 | 500 | 39.27 | " |
| 1035 | | 12.78 | 6.50 | 0.154 | 6.47 | 145.8 | 63.0 | 500 | 39.27 | " |
| 1040 | | 12.42 | 6.50 | 0.154 | 6.43 | 145.2 | 64.6 | 500 | 39.27 | " |
| 1045 | | 12.66 | 6.49 | 0.153 | 6.36 | 145.6 | 32.5 | 500 | 39.27 | " |
| 1050 | | 12.84 | 6.50 | 0.154 | 6.34 | 145.1 | 29.0 | 500 | 39.27 | " |
| 1055 | | 12.90 | 6.50 | 0.153 | 6.35 | 144.7 | 25.8 | 500 | 39.27 | " |
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RESOLUTION CONSULTANTS

Well ID: RE12202

Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 3/24/15 Time: Start 8:15 am/pm
 Project No: 60266526 Finish 12:00 am/pm
 Site Location: Curtis & Hayden
 Weather Conds: Sunny 25° Collector(s): Paul Kareth

1. WATER LEVEL DATA: (measured from Top of Casing)

a. Total Well Length 615 c. Length of Water Column _____ (a-b) Casing Diameter/Material
 b. Water Table Depth 39.53 d. Calculated System Volume (see back) 20ft screen, 13.1 gal
 4-inch PVC

2. WELL PURGE DATA

a. Purge Method: Geotech bladder pump with drop tube assembly
 b. Acceptance Criteria defined (see workplan)
 - Temperature ± 3% - D.O. ± 10% (values >0.5 mg/L) Turbidity ± 10%
 - pH ± 0.1 unit - ORP ± 10mV
 - Sp. Cond. ± 3% - Drawdown < 0.3' Remove a minimum 1 screen volume

c. Field Testing Equipment used:

| Make | Model | Serial Number |
|--------------|-----------------|---------------|
| <u>YSI</u> | <u>556MPS</u> | <u>71124</u> |
| <u>Hanna</u> | <u>HI 98703</u> | <u>69177</u> |

| Time (24hr) | Volume Removed (Liters) | Temp. (°C) | pH | Spec. Cond. (mS/cm) | DO (mg/L) | ORP (mV) | Turbidity (NTU) | Flow Rate (ml/min) | Depth to water (ft) | Color/Odor |
|-------------|-------------------------|--------------|-------------|---------------------|--------------|--------------|-----------------|--------------------|---------------------|-------------------------------------|
| <u>925</u> | <u>ON</u> | | | | | | | | | |
| <u>0935</u> | | <u>9.37</u> | <u>7.07</u> | <u>0.158</u> | <u>15.82</u> | <u>-54.6</u> | <u>3.86</u> | <u>480</u> | <u>39.78</u> | <u>clear</u> |
| <u>0940</u> | | <u>11.64</u> | <u>6.17</u> | <u>0.153</u> | <u>10.47</u> | <u>-43</u> | <u>2.75</u> | | <u>39.45</u> | <u> </u> |
| <u>0945</u> | | <u>10.38</u> | <u>5.59</u> | <u>0.150</u> | <u>7.01</u> | <u>-32.3</u> | | <u>350</u> | | |
| <u>0950</u> | | <u>10.41</u> | <u>5.60</u> | <u>0.150</u> | <u>7.20</u> | <u>-31.4</u> | | | <u>39.44</u> | <u>irritate adjust hose fitting</u> |
| <u>1000</u> | | | | | | | | <u>450</u> | | |
| <u>1010</u> | | <u>11.12</u> | <u>5.85</u> | <u>0.120</u> | <u>5.70</u> | <u>-36.9</u> | | | | |

d. Acceptance criteria pass/fail

| | Yes | No | N/A |
|-------------------------------------|-------------------------------------|--------------------------|--------------------------|
| Has required volume been removed | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Has required turbidity been reached | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Have parameters stabilized | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |

If no or N/A - Explain below.

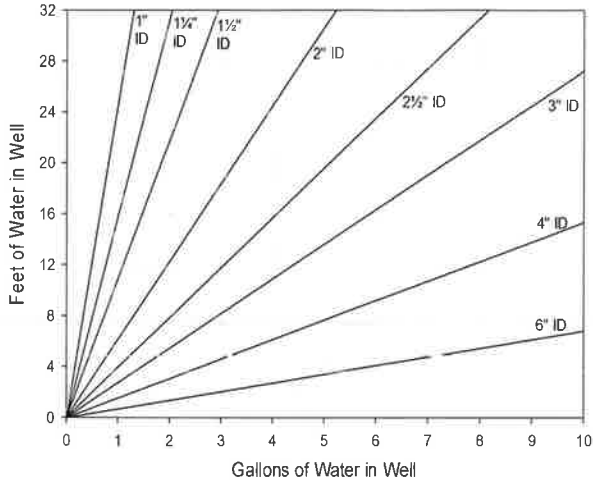
3. SAMPLE COLLECTION: Method: Geotech bladder pump with drop tube assembly

| Sample ID | Container Type | No. of Containers | Preservation | Analysis Req. | Time |
|--------------------------|-------------------|-------------------|--------------|--------------------|---------------------|
| <u>RE12202-GW-052415</u> | <u>40-mL vial</u> | <u>9/8</u> | <u>HCl</u> | <u>VOCs</u> | <u>1130 MS, MSD</u> |
| <u>RE12202-GW-032415</u> | <u>1-L amber</u> | <u>6/2</u> | <u>none</u> | <u>1,4-Dioxane</u> | <u>1130 MS, MSD</u> |

Comments: hit bottom, tubing was not cut

Signature: Paul Kareth Date: 3/24/15

Purge Volume Calculation



| ID (in) | Gallon | Liter |
|---------|--------|--------|
| 0.25 | 0.0025 | 0.0097 |
| 0.375 | 0.0057 | 0.0217 |
| 0.5 | 0.0102 | 0.0386 |
| 0.75 | 0.0229 | 0.0869 |
| 1 | 0.0408 | 0.1544 |
| 1.25 | 0.0637 | 0.2413 |
| 1.5 | 0.0918 | 0.3475 |
| 2 | 0.1632 | 0.6178 |
| 2.5 | 0.2550 | 0.9653 |
| 3 | 0.3672 | 1.3900 |
| 4 | 0.6528 | 2.4711 |
| 6 | 1.4688 | 5.5600 |

1 screen volume
 15 ft = ~~56.2~~ L / 9.8 G
 20 ft = ~~74.9~~ L / 13.1 G
 25 ft = ~~93.6~~ L / 16.3 G

Well ID: RE-122020 925

(continued from front)

| Time (24 hr) | Volume Removed (Liters) | Temp (°C) | pH | Spec. Cond. (mS/cm) | DO (mg/L) | ORP (mV) | Turbidity (NTU) | Flow Rate (ml/min) | Depth to water (ft) | Color/Odor |
|-----------------|-------------------------------|--------------|------|------------------------|--------------|-------------|--------------------|-----------------------|------------------------|----------------|
| | | | | | | | | | | |
| 1015 | | 11.31 | 5.74 | 0.117 | 5.59 | -33.3 | 15.1 | | 39.51 | |
| 1020 | 5 gal | 12.01 | 5.71 | 0.119 | 5.32 | -30.9 | | 500 | | |
| 1020 | | 11.99 | 5.68 | 0.119 | 5.57 | -29.0 | 14.6 | | | |
| 1025 | | 12.59 | 5.68 | 0.121 | 5.41 | -28.4 | | | | |
| 1030 | | 12.70 | 5.67 | 0.121 | 5.32 | -27.6 | | | | |
| 1035 | | 12.51 | 5.67 | 0.121 | 5.31 | -25.2 | | | 39.52 | |
| 1040 | | 12.77 | 5.66 | 0.121 | 5.27 | -23.5 | 5.62 | | | |
| 1045 | | 12.41 | 5.65 | 0.120 | 5.25 | -21.9 | | 500 | | |
| 1050 | | 12.70 | 5.65 | 0.120 | 5.22 | -20.9 | | | | |
| 1055 | 10 gal | 12.79 | 5.65 | 0.120 | 5.19 | -19.2 | | | | |
| 1100 | | 12.83 | 5.65 | 0.121 | 5.17 | -18.4 | 3.86 | | | |
| 1105 | | 12.62 | 5.65 | 0.121 | 5.10 | -16.6 | | | | |
| 1110 | | 12.91 | 5.65 | 0.121 | 5.09 | -15.3 | | 500 | 39.58 | |
| 1115 | | | | | | | | | | |
| 1130 | | | | | | | | | | Sample MS, MSD |
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Well ID: RE122-03

Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 3/24/15 Time: Start 9:15 am/pm
 Project No: 60266526 Finish 12:15 am/pm
 Site Location: Curtis & Hayden
 Weather Conds: Sunny 25°C Collector(s): SC

1. WATER LEVEL DATA: (measured from Top of Casing)
 a. Total Well Length 740 c. Length of Water Column _____ (a-b) Casing Diameter/Material 4-inch PVC
 b. Water Table Depth 40.19 d. Calculated System Volume (see back) soft screen 13.1 gal

2. WELL PURGE DATA
 a. Purge Method: Geotech bladder pump with drop tube assembly
 b. Acceptance Criteria defined (see workplan)
 - Temperature ± 3% - D.O. ± 10% (values >0.5 mg/L) Turbidity ± 10%
 - pH ± 0.1 unit - ORP ± 10mV
 - Sp. Cond. ± 3% - Drawdown < 0.3' Remove a minimum 1 screen volume
 c. Field Testing Equipment used:

| Volume | Make | Model | Serial Number |
|--------|-----------------|----------------|------------------------|
| | <u>YSI 556</u> | <u>U28070X</u> | <u>602336494 71977</u> |
| | <u>Hann MPS</u> | <u>U54589X</u> | <u>14J21 69177</u> |

| Time (24hr) | Volume Removed (Liters) | Temp. (°C) | pH | Spec. Cond. (mS/cm) | DO (mg/L) | ORP (mV) | Turbidity (NTU) | Flow Rate (ml/min) | Depth to water (ft) | Color/Odor |
|--------------|-------------------------|--------------|-------------|---------------------|-------------|--------------|-----------------|--------------------|---------------------|---------------|
| <u>9:50</u> | | | | | | | | | | |
| <u>10:00</u> | | <u>11.63</u> | <u>5.99</u> | <u>0.055</u> | <u>8.55</u> | <u>149.1</u> | <u>250</u> | <u>500</u> | <u>40.26</u> | <u>cloudy</u> |
| <u>10:05</u> | | <u>11.43</u> | <u>5.82</u> | <u>0.054</u> | <u>8.72</u> | <u>155.6</u> | <u>30.5</u> | | | |
| <u>10:10</u> | | <u>11.72</u> | <u>5.47</u> | <u>0.050</u> | <u>5.86</u> | <u>185.7</u> | <u>291</u> | <u>500</u> | <u>40.70</u> | <u>cloudy</u> |
| <u>10:20</u> | | <u>11.42</u> | <u>5.56</u> | <u>0.051</u> | <u>5.62</u> | <u>187.8</u> | <u>299</u> | <u>500</u> | <u>40.70</u> | <u>"</u> |
| <u>10:25</u> | | <u>11.71</u> | <u>6.13</u> | <u>0.087</u> | <u>5.32</u> | <u>167.9</u> | | | <u>40.70</u> | <u>"</u> |
| <u>10:30</u> | <u>56</u> | <u>12.22</u> | <u>5.89</u> | <u>0.063</u> | <u>4.92</u> | <u>177.2</u> | <u>141</u> | <u>500</u> | <u>40.70</u> | |

d. Acceptance criteria pass/fail

| | Yes | No | N/A |
|-------------------------------------|-------------------------------------|--------------------------|--------------------------|
| Has required volume been removed | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Has required turbidity been reached | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Have parameters stabilized | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |

If no or N/A - Explain below.

3. SAMPLE COLLECTION: Method: Geotech bladder pump with drop tube assembly

| Sample ID | Container Type | No. of Containers | Preservation | Analysis Req. | Time |
|--------------------------|-------------------|-------------------|--------------|--------------------|--------------|
| <u>RE12203-GW-032415</u> | <u>40-mL vial</u> | <u>3</u> | <u>HCl</u> | <u>VOCs</u> | <u>11:50</u> |
| <u>RE12203-GW-032415</u> | <u>1-L amber</u> | <u>2</u> | <u>none</u> | <u>1,4-Dioxane</u> | <u>11:50</u> |

Comments: tubing hits bottom! ~34ft too long!

Signature: [Signature] Date: 03/24/15

Section 4

Analytical Data Validation



DATA VALIDATION REPORT

| | | |
|------------------------|--|-------------------------------|
| Project: | Regional Groundwater Investigation — NWIRP Bethpage | |
| Laboratory: | Katahdin Analytical | |
| Sample Delivery Group: | SI1843 | |
| Analyses/Method: | Volatile Organic Compounds by U.S. EPA SW-846 Method 8260C 1,4-Dioxane by U.S. EPA SW-846 Method 8270D via Selective Ion Monitoring (SIM) | |
| Validation Level: | 3 | |
| Project Number: | 0888812477.SA.DV | |
| Prepared by: | Dana Miller/Resolution Consultants | Completed on: 06/08/2015 |
| Reviewed by: | Tina Cantwell/Resolution Consultants | File Name: SI1843_8260C_8270D |

SUMMARY

This report summarizes data review findings for samples listed below, collected by Resolution Consultants from the Regional Groundwater Investigation — NWIRP Bethpage site on 24 March 2015 in accordance with the following Sampling and Analysis Plans:

- *Sampling and Analysis Plan, Bethpage, New York.* (Resolution Consultants April 2013).
- *UFP SAP Addendum, Installation of Vertical Profile Borings and Monitoring Wells, Operable Unit 2, NWIRP Bethpage, New York.* (Resolution Consultants November 2013).
- *UFP SAP Addendum, Inclusion of Additional Target Analytes for Volatile Organics Analyses, NWIRP Bethpage OU2, Bethpage, New York.* (Resolution Consultants August 2014).

| Sample ID | Matrix/Sample Type | Analysis |
|---------------------|--------------------|-------------------|
| DUPLICATE-GW-032315 | Field Duplicate | 8260C / 8270D_SIM |
| RE103D1-GW-032315 | Groundwater | 8260C / 8270D_SIM |
| RE103D2-GW-032315 | Groundwater | 8260C / 8270D_SIM |
| RE103D3-GW-032315 | Groundwater | 8260C / 8270D_SIM |
| RE104D1-GW-032315 | Groundwater | 8260C / 8270D_SIM |
| RE104D2-GW-032315 | Groundwater | 8260C / 8270D_SIM |
| RE104D3-GW-032315 | Groundwater | 8260C / 8270D_SIM |
| RE122D1-GW-032415 | Groundwater | 8260C / 8270D_SIM |
| RE122D2-GW-032415 | Groundwater | 8260C / 8270D_SIM |
| RE122D3-GW-032415 | Groundwater | 8260C / 8270D_SIM |
| TT101D1-GW-032415 | Groundwater | 8260C / 8270D_SIM |

| Sample ID | Matrix/Sample Type | Analysis |
|-------------------|--------------------|-------------------|
| TT101D2-GW-032415 | Groundwater | 8260C / 8270D_SIM |
| TT101D-GW-032415 | Groundwater | 8260C / 8270D_SIM |
| TRIPBLANK031615 | Trip Blank | 8260C |

Data validation activities were conducted using the following guidance documents: *Test Methods for Evaluating Solid Waste, Physical/Chemical Methods SW-846, specifically Method 8260C, Volatile Organic Compounds by Gas Chromatography/Mass Spectrometry* (U.S. EPA, 2006), *SW-846 Method 8270D, Semivolatile Organic Compounds by Gas Chromatograph/Mass Spectrometry* (U.S. EPA, 2007), *U.S. Environmental Protection Agency (U.S. EPA) Contract Laboratory Program National Functional Guidelines for Superfund Organic Methods Data Review* (June 2008), and *Department of Defense Quality Systems Manual for Environmental Laboratories, Version 4.2* (October 2010). In the absence of method-specific information, laboratory quality control (QC) limits, project-specific requirements and/or professional judgment were used as appropriate.

REVIEW ELEMENTS

The data were evaluated based on the following parameters (where applicable to the method):

- ✓ Data completeness (chain-of-custody)/sample integrity
- ✓ Holding times and sample preservation
- ✓ Gas chromatography/Mass spectrometer performance checks
- ✗ Initial calibration/continuing calibration verification
- ✓ Laboratory blanks/trip blanks
- ✗ Surrogate spike recoveries
- ✓ Matrix spike and/or matrix spike duplicate results
- ✗ Laboratory control sample (LCS)/laboratory control sample duplicate (LCSD) results
- ✓ Field duplicates
- ✓ Internal standards
- ✓ Sample results/reporting issues

The symbol (✓) indicates that no validation qualifiers were applied based on this parameter. NA indicates that the parameter was not included as part of this data set or was not applicable to this validation and therefore not reviewed. Acceptable data parameters for which all criteria were met and no qualification was performed and non-conformance or other issues that were noted during validation, but did not result in qualification of data are not discussed further.

The symbol (X) indicates that a QC non-conformance resulted in the qualification of data. Any QC non-conformance that resulted in the qualification of data is discussed below.

RESULTS

Initial Calibration/Continuing Calibration Verification

Calibration data were reviewed for conformance with the QC acceptance criteria to ensure that:

- the initial calibration percent relative standard deviation, correlation coefficient/coefficient of determination, and/or response factor method acceptance criteria were met;
- the initial calibration verification (ICV) standard percent recovery acceptance criteria were met;
- the continuing calibration verification standard (CCV) method percent difference or percent drift (%Ds) and response factor acceptance criteria were met; and
- the retention time method acceptance criteria were met.

Data qualification to the analytes associated with the specific calibration verification was as follows:

ICV Recovery Non-conformance:

| Criteria | Actions | |
|-----------------|------------------|----------------------|
| | Detected Results | Non-detected Results |
| Recovery > 120% | J | UJ |
| Recovery < 80% | J | UJ |

Notes:

J = Estimated
UJ = Undetected and estimated

CCV Linearity Non-conformance:

| Criteria | Actions | |
|-----------------------------|------------------|----------------------|
| | Detected Results | Non-detected Results |
| %Difference or %Drift > 20% | J | UJ |

Notes:

J = Estimated
UJ = Undetected and estimated

ICV and CCV non-conformances are summarized in Attachment A in Table's A-1 and A-2.

Surrogate Spike Recoveries

Surrogates provide information needed to assess the accuracy of analyses. Known amounts of surrogate compounds, or compounds which are not likely to be found in the actual samples, are added to each organic sample to check for accuracy. If surrogate percent recoveries (%Rs) are close to the known concentrations, the reported target compound concentrations are assumed to be accurate. Data qualification on the basis of surrogate recovery was as follows:

Surrogate Recovery Non-conformance Chart:

| Criteria | Action | |
|---------------|----------|------------------|
| | Detected | Non-detected |
| % R > UL | J | No qualification |
| 20% ≤ %R < LL | J | UJ |
| %R < 20% | J | Rejected |

Notes:

- %R = Percent recovery
- UL = Upper limit
- LL = Lower limit
- J = Estimated
- UJ = Undetected and estimated

Surrogate recovery non-conformance is summarized in Attachment A in Table A-3.

Laboratory Control Samples / Laboratory Control Sample Duplicate

LCS %Rs is used to monitor the overall accuracy and performance of each step during analysis, including sample preparation. The laboratory analyzed LCSs in duplicate when matrix spike/matrix spike duplicates were not reported. In these instances, the laboratory determined precision between the duplicated values. Non-conformance is summarized in Attachment A in Table A-4. Data qualification to the analytes associated with the specific LCS / LCS duplicate was as follows:

Laboratory Control Sample / Laboratory Control Sample Duplicate Non-conformance Chart:

| Criteria | Action | |
|----------------|----------|------------------|
| | Detected | Non-detected |
| %R or RPD > UL | J | No qualification |
| %R < LL | J | UJ |
| %R < 20% | J | Rejected |

Notes:

| | | |
|-----|---|-----------------------------|
| %R | = | Percent recovery |
| RPD | = | Relative percent difference |
| UL | = | Upper limit |
| LL | = | Lower limit |
| J | = | Estimated |
| UJ | = | Undetected and estimated |

Qualifications Actions

The data was reviewed independently from the laboratory to assess data quality. All compounds detected at concentrations less than the limit of quantitation but greater than the method detection limit were qualified by the laboratory as estimated (J). This "J" qualifier was retained during data validation. Any sample that was analyzed at a dilution because of high concentrations of target or non-target analytes was checked to confirm that the results and/or sample-specific limit of quantitation and limit of detections were adjusted accordingly by the laboratory.

No results were rejected; therefore, analytical completeness was calculated to be 100 percent. Data not qualified during data review are considered usable by the project. The remaining results qualified as estimated may be high or low, but the data are usable for their intended purpose, according to U.S. Environmental Protection Agency and Department of Defense guidelines. Final data review qualifiers used to describe results and how they should be interpreted by the end data user are provided in Attachment B and Attachment C. Attachment D provides final results after data review.

ATTACHMENTS

- Attachment A: Non-Conformance Summary Tables
- Attachment B: Qualifier Codes and Explanations
- Attachment C: Reason Codes and Explanations
- Attachment D: Final Results after Data Review

Attachment A
Non-Conformance Summary Table

| Table A-1 Initial Calibration Verification Non-Conformance | | | | | | |
|---|-------------------------|------------|--------|--------|---------------------|-----------|
| Method | Analyte | ICV ID | %R | Limit | Associated Samples | Qualifier |
| 8260C | Dichlorodifluoromethane | WG160458-7 | 132.53 | 80-120 | DUPLICATE-GW-032315 | UJ |
| 8260C | Dichlorodifluoromethane | WG160458-7 | 132.53 | 80-120 | RE103D1-GW-032315 | UJ |
| 8260C | Dichlorodifluoromethane | WG160458-7 | 132.53 | 80-120 | RE103D2-GW-032315 | UJ |
| 8260C | Dichlorodifluoromethane | WG160458-7 | 132.53 | 80-120 | RE103D3-GW-032315 | UJ |
| 8260C | Dichlorodifluoromethane | WG160458-7 | 132.53 | 80-120 | RE104D1-GW-032315 | UJ |
| 8260C | Dichlorodifluoromethane | WG160458-7 | 132.53 | 80-120 | RE104D2-GW-032315 | UJ |
| 8260C | Dichlorodifluoromethane | WG160458-7 | 132.53 | 80-120 | RE104D3-GW-032315 | UJ |
| 8260C | Dichlorodifluoromethane | WG160458-7 | 132.53 | 80-120 | RE122D1-GW-032415 | UJ |
| 8260C | Dichlorodifluoromethane | WG160458-7 | 132.53 | 80-120 | RE122D2-GW-032415 | UJ |
| 8260C | Dichlorodifluoromethane | WG160458-7 | 132.53 | 80-120 | RE122D3-GW-032415 | UJ |
| 8260C | Dichlorodifluoromethane | WG160458-7 | 132.53 | 80-120 | TT101D1-GW-032415 | J |
| 8260C | Dichlorodifluoromethane | WG160458-7 | 132.53 | 80-120 | TT101D2-GW-032415 | J |
| 8260C | Dichlorodifluoromethane | WG160458-7 | 132.53 | 80-120 | TT101D-GW-032415 | UJ |
| 8260C | Dichlorodifluoromethane | WG160458-7 | 132.53 | 80-120 | TRIPBLANK031615 | UJ |
| 8260C | Chloromethane | WG160458-7 | 121.77 | 80-120 | DUPLICATE-GW-032315 | UJ |
| 8260C | Chloromethane | WG160458-7 | 121.77 | 80-120 | RE103D1-GW-032315 | UJ |
| 8260C | Chloromethane | WG160458-7 | 121.77 | 80-120 | RE103D2-GW-032315 | UJ |
| 8260C | Chloromethane | WG160458-7 | 121.77 | 80-120 | RE103D3-GW-032315 | UJ |
| 8260C | Chloromethane | WG160458-7 | 121.77 | 80-120 | RE104D1-GW-032315 | UJ |
| 8260C | Chloromethane | WG160458-7 | 121.77 | 80-120 | RE104D2-GW-032315 | UJ |
| 8260C | Chloromethane | WG160458-7 | 121.77 | 80-120 | RE104D3-GW-032315 | UJ |
| 8260C | Chloromethane | WG160458-7 | 121.77 | 80-120 | RE122D1-GW-032415 | UJ |
| 8260C | Chloromethane | WG160458-7 | 121.77 | 80-120 | RE122D2-GW-032415 | UJ |
| 8260C | Chloromethane | WG160458-7 | 121.77 | 80-120 | RE122D3-GW-032415 | UJ |
| 8260C | Chloromethane | WG160458-7 | 121.77 | 80-120 | TT101D1-GW-032415 | UJ |
| 8260C | Chloromethane | WG160458-7 | 121.77 | 80-120 | TT101D2-GW-032415 | UJ |
| 8260C | Chloromethane | WG160458-7 | 121.77 | 80-120 | TT101D-GW-032415 | UJ |
| 8260C | Chloromethane | WG160458-7 | 121.77 | 80-120 | TRIPBLANK031615 | UJ |
| 8260C | Bromomethane | WG160458-7 | 130.01 | 80-120 | DUPLICATE-GW-032315 | UJ |
| 8260C | Bromomethane | WG160458-7 | 130.01 | 80-120 | RE103D1-GW-032315 | UJ |
| 8260C | Bromomethane | WG160458-7 | 130.01 | 80-120 | RE103D2-GW-032315 | UJ |
| 8260C | Bromomethane | WG160458-7 | 130.01 | 80-120 | RE103D3-GW-032315 | UJ |
| 8260C | Bromomethane | WG160458-7 | 130.01 | 80-120 | RE104D1-GW-032315 | UJ |
| 8260C | Bromomethane | WG160458-7 | 130.01 | 80-120 | RE104D2-GW-032315 | UJ |
| 8260C | Bromomethane | WG160458-7 | 130.01 | 80-120 | RE104D3-GW-032315 | UJ |
| 8260C | Bromomethane | WG160458-7 | 130.01 | 80-120 | RE122D1-GW-032415 | UJ |
| 8260C | Bromomethane | WG160458-7 | 130.01 | 80-120 | RE122D2-GW-032415 | UJ |
| 8260C | Bromomethane | WG160458-7 | 130.01 | 80-120 | RE122D3-GW-032415 | UJ |
| 8260C | Bromomethane | WG160458-7 | 130.01 | 80-120 | TT101D1-GW-032415 | UJ |
| 8260C | Bromomethane | WG160458-7 | 130.01 | 80-120 | TT101D2-GW-032415 | UJ |
| 8260C | Bromomethane | WG160458-7 | 130.01 | 80-120 | TT101D-GW-032415 | UJ |
| 8260C | Bromomethane | WG160458-7 | 130.01 | 80-120 | TRIPBLANK031615 | UJ |
| 8260C | Carbon Disulfide | WG160458-7 | 544.89 | 80-120 | DUPLICATE-GW-032315 | UJ |
| 8260C | Carbon Disulfide | WG160458-7 | 544.89 | 80-120 | RE103D1-GW-032315 | J |
| 8260C | Carbon Disulfide | WG160458-7 | 544.89 | 80-120 | RE103D2-GW-032315 | J |
| 8260C | Carbon Disulfide | WG160458-7 | 544.89 | 80-120 | RE103D3-GW-032315 | J |
| 8260C | Carbon Disulfide | WG160458-7 | 544.89 | 80-120 | RE104D1-GW-032315 | UJ |

Table A-1
Initial Calibration Verification Non-Conformance

| Method | Analyte | ICV ID | %R | Limit | Associated Samples | Qualifier |
|--------|------------------|------------|--------|--------|---------------------|-----------|
| 8260C | Carbon Disulfide | WG160458-7 | 544.89 | 80-120 | RE104D2-GW-032315 | UJ |
| 8260C | Carbon Disulfide | WG160458-7 | 544.89 | 80-120 | RE104D3-GW-032315 | UJ |
| 8260C | Carbon Disulfide | WG160458-7 | 544.89 | 80-120 | RE122D1-GW-032415 | UJ |
| 8260C | Carbon Disulfide | WG160458-7 | 544.89 | 80-120 | RE122D2-GW-032415 | UJ |
| 8260C | Carbon Disulfide | WG160458-7 | 544.89 | 80-120 | RE122D3-GW-032415 | UJ |
| 8260C | Carbon Disulfide | WG160458-7 | 544.89 | 80-120 | TT101D1-GW-032415 | UJ |
| 8260C | Carbon Disulfide | WG160458-7 | 544.89 | 80-120 | TT101D2-GW-032415 | UJ |
| 8260C | Carbon Disulfide | WG160458-7 | 544.89 | 80-120 | TT101D-GW-032415 | UJ |
| 8260C | Carbon Disulfide | WG160458-7 | 544.89 | 80-120 | TRIPBLANK031615 | UJ |
| 8260C | Acetone | WG160458-7 | 137.34 | 80-120 | DUPLICATE-GW-032315 | UJ |
| 8260C | Acetone | WG160458-7 | 137.34 | 80-120 | RE103D1-GW-032315 | UJ |
| 8260C | Acetone | WG160458-7 | 137.34 | 80-120 | RE103D2-GW-032315 | UJ |
| 8260C | Acetone | WG160458-7 | 137.34 | 80-120 | RE103D3-GW-032315 | UJ |
| 8260C | Acetone | WG160458-7 | 137.34 | 80-120 | RE104D1-GW-032315 | UJ |
| 8260C | Acetone | WG160458-7 | 137.34 | 80-120 | RE104D2-GW-032315 | UJ |
| 8260C | Acetone | WG160458-7 | 137.34 | 80-120 | RE104D3-GW-032315 | UJ |
| 8260C | Acetone | WG160458-7 | 137.34 | 80-120 | RE122D1-GW-032415 | UJ |
| 8260C | Acetone | WG160458-7 | 137.34 | 80-120 | RE122D2-GW-032415 | UJ |
| 8260C | Acetone | WG160458-7 | 137.34 | 80-120 | RE122D3-GW-032415 | UJ |
| 8260C | Acetone | WG160458-7 | 137.34 | 80-120 | TT101D1-GW-032415 | UJ |
| 8260C | Acetone | WG160458-7 | 137.34 | 80-120 | TT101D2-GW-032415 | UJ |
| 8260C | Acetone | WG160458-7 | 137.34 | 80-120 | TT101D-GW-032415 | UJ |
| 8260C | Acetone | WG160458-7 | 137.34 | 80-120 | TRIPBLANK031615 | UJ |
| 8260C | 2-Butanone | WG160458-7 | 134.65 | 80-120 | DUPLICATE-GW-032315 | UJ |
| 8260C | 2-Butanone | WG160458-7 | 134.65 | 80-120 | RE103D1-GW-032315 | UJ |
| 8260C | 2-Butanone | WG160458-7 | 134.65 | 80-120 | RE103D2-GW-032315 | UJ |
| 8260C | 2-Butanone | WG160458-7 | 134.65 | 80-120 | RE103D3-GW-032315 | UJ |
| 8260C | 2-Butanone | WG160458-7 | 134.65 | 80-120 | RE104D1-GW-032315 | UJ |
| 8260C | 2-Butanone | WG160458-7 | 134.65 | 80-120 | RE104D2-GW-032315 | UJ |
| 8260C | 2-Butanone | WG160458-7 | 134.65 | 80-120 | RE104D3-GW-032315 | UJ |
| 8260C | 2-Butanone | WG160458-7 | 134.65 | 80-120 | RE122D1-GW-032415 | UJ |
| 8260C | 2-Butanone | WG160458-7 | 134.65 | 80-120 | RE122D2-GW-032415 | UJ |
| 8260C | 2-Butanone | WG160458-7 | 134.65 | 80-120 | RE122D3-GW-032415 | UJ |
| 8260C | 2-Butanone | WG160458-7 | 134.65 | 80-120 | TT101D1-GW-032415 | UJ |
| 8260C | 2-Butanone | WG160458-7 | 134.65 | 80-120 | TT101D2-GW-032415 | UJ |
| 8260C | 2-Butanone | WG160458-7 | 134.65 | 80-120 | TT101D-GW-032415 | UJ |
| 8260C | 2-Butanone | WG160458-7 | 134.65 | 80-120 | TRIPBLANK031615 | UJ |
| 8260C | Cyclohexane | WG160458-7 | 193.37 | 80-120 | DUPLICATE-GW-032315 | UJ |
| 8260C | Cyclohexane | WG160458-7 | 193.37 | 80-120 | RE103D1-GW-032315 | UJ |
| 8260C | Cyclohexane | WG160458-7 | 193.37 | 80-120 | RE103D2-GW-032315 | UJ |
| 8260C | Cyclohexane | WG160458-7 | 193.37 | 80-120 | RE103D3-GW-032315 | UJ |
| 8260C | Cyclohexane | WG160458-7 | 193.37 | 80-120 | RE104D1-GW-032315 | UJ |
| 8260C | Cyclohexane | WG160458-7 | 193.37 | 80-120 | RE104D2-GW-032315 | UJ |
| 8260C | Cyclohexane | WG160458-7 | 193.37 | 80-120 | RE104D3-GW-032315 | UJ |
| 8260C | Cyclohexane | WG160458-7 | 193.37 | 80-120 | RE122D1-GW-032415 | UJ |
| 8260C | Cyclohexane | WG160458-7 | 193.37 | 80-120 | RE122D2-GW-032415 | UJ |
| 8260C | Cyclohexane | WG160458-7 | 193.37 | 80-120 | RE122D3-GW-032415 | UJ |
| 8260C | Cyclohexane | WG160458-7 | 193.37 | 80-120 | TT101D1-GW-032415 | UJ |

| Table A-1 Initial Calibration Verification Non-Conformance | | | | | | |
|---|-------------------|------------|--------|--------|---------------------|-----------|
| Method | Analyte | ICV ID | %R | Limit | Associated Samples | Qualifier |
| 8260C | Cyclohexane | WG160458-7 | 193.37 | 80-120 | TT101D2-GW-032415 | UJ |
| 8260C | Cyclohexane | WG160458-7 | 193.37 | 80-120 | TT101D-GW-032415 | UJ |
| 8260C | Cyclohexane | WG160458-7 | 193.37 | 80-120 | TRIPBLANK031615 | UJ |
| 8260C | Tetrachloroethene | WG160458-7 | 124.99 | 80-120 | DUPLICATE-GW-032315 | UJ |
| 8260C | Tetrachloroethene | WG160458-7 | 124.99 | 80-120 | RE103D1-GW-032315 | J |
| 8260C | Tetrachloroethene | WG160458-7 | 124.99 | 80-120 | RE103D2-GW-032315 | J |
| 8260C | Tetrachloroethene | WG160458-7 | 124.99 | 80-120 | RE103D3-GW-032315 | J |
| 8260C | Tetrachloroethene | WG160458-7 | 124.99 | 80-120 | RE104D1-GW-032315 | J |
| 8260C | Tetrachloroethene | WG160458-7 | 124.99 | 80-120 | RE104D2-GW-032315 | UJ |
| 8260C | Tetrachloroethene | WG160458-7 | 124.99 | 80-120 | RE104D3-GW-032315 | UJ |
| 8260C | Tetrachloroethene | WG160458-7 | 124.99 | 80-120 | RE122D1-GW-032415 | J |
| 8260C | Tetrachloroethene | WG160458-7 | 124.99 | 80-120 | RE122D2-GW-032415 | J |
| 8260C | Tetrachloroethene | WG160458-7 | 124.99 | 80-120 | RE122D3-GW-032415 | UJ |
| 8260C | Tetrachloroethene | WG160458-7 | 124.99 | 80-120 | TT101D1-GW-032415 | UJ |
| 8260C | Tetrachloroethene | WG160458-7 | 124.99 | 80-120 | TT101D2-GW-032415 | UJ |
| 8260C | Tetrachloroethene | WG160458-7 | 124.99 | 80-120 | TT101D-GW-032415 | J |
| 8260C | Tetrachloroethene | WG160458-7 | 124.99 | 80-120 | TRIPBLANK031615 | UJ |
| 8260C | 2-Hexanone | WG160458-7 | 130.94 | 80-120 | DUPLICATE-GW-032315 | UJ |
| 8260C | 2-Hexanone | WG160458-7 | 130.94 | 80-120 | RE103D1-GW-032315 | UJ |
| 8260C | 2-Hexanone | WG160458-7 | 130.94 | 80-120 | RE103D2-GW-032315 | UJ |
| 8260C | 2-Hexanone | WG160458-7 | 130.94 | 80-120 | RE103D3-GW-032315 | UJ |
| 8260C | 2-Hexanone | WG160458-7 | 130.94 | 80-120 | RE104D1-GW-032315 | UJ |
| 8260C | 2-Hexanone | WG160458-7 | 130.94 | 80-120 | RE104D2-GW-032315 | UJ |
| 8260C | 2-Hexanone | WG160458-7 | 130.94 | 80-120 | RE104D3-GW-032315 | UJ |
| 8260C | 2-Hexanone | WG160458-7 | 130.94 | 80-120 | RE122D1-GW-032415 | UJ |
| 8260C | 2-Hexanone | WG160458-7 | 130.94 | 80-120 | RE122D2-GW-032415 | UJ |
| 8260C | 2-Hexanone | WG160458-7 | 130.94 | 80-120 | RE122D3-GW-032415 | UJ |
| 8260C | 2-Hexanone | WG160458-7 | 130.94 | 80-120 | TT101D1-GW-032415 | UJ |
| 8260C | 2-Hexanone | WG160458-7 | 130.94 | 80-120 | TT101D2-GW-032415 | UJ |
| 8260C | 2-Hexanone | WG160458-7 | 130.94 | 80-120 | TT101D-GW-032415 | UJ |
| 8260C | 2-Hexanone | WG160458-7 | 130.94 | 80-120 | TRIPBLANK031615 | UJ |

Notes:

- ICV = Initial calibration verification
- %R = Percent recovery
- J = Detected analyte in associate sample qualified estimated "J" due to potential bias.
- UJ = Non-detected analyte in associated sample qualified estimated "UJ" due to potential bias.

| Table A-2 Continuing Calibration Verification Non-Conformance | | | | | | |
|--|--------------------|---------|-------|-------|---------------------|-----------|
| Method | Analyte | CCV ID | %D | Limit | Associated Samples | Qualifier |
| 8260C | Methyl cyclohexane | C2012.D | 49.68 | 20 | RE104D2-GW-032315 | UJ |
| 8260C | Methyl cyclohexane | C2012.D | 49.68 | 20 | RE122D3-GW-032415 | UJ |
| 8260C | Methyl cyclohexane | C2012.D | 49.68 | 20 | TRIPBLANK031615 | UJ |
| 8260C | Methyl cyclohexane | C2012.D | 49.68 | 20 | TT101D1-GW-032415 | UJ |
| 8260C | Methyl cyclohexane | C2012.D | 49.68 | 20 | TT101D2-GW-032415 | UJ |
| 8260C | Chloroethane | C1970.D | 40.73 | 20 | RE103D1-GW-032315 | UJ |
| 8260C | Chloroethane | C1970.D | 40.73 | 20 | RE103D2-GW-032315 | UJ |
| 8260C | Chloroethane | C1970.D | 4.073 | 20 | RE103D3-GW-032315 | UJ |
| 8260C | Chloroethane | C1986.D | 36.53 | 20 | DUPLICATE-GW-032315 | UJ |

| Table A-2 Continuing Calibration Verification Non-Conformance | | | | | | |
|--|--------------|---------|-------|-------|--------------------|-----------|
| Method | Analyte | CCV ID | %D | Limit | Associated Samples | Qualifier |
| 8260C | Chloroethane | C1986.D | 36.53 | 20 | RE104D1-GW-032315 | UJ |
| 8260C | Chloroethane | C1986.D | 36.53 | 20 | RE104D3-GW-032315 | UJ |
| 8260C | Chloroethane | C1986.D | 36.53 | 20 | RE122D1-GW-032415 | UJ |
| 8260C | Chloroethane | C1986.D | 36.53 | 20 | RE122D2-GW-032415 | UJ |
| 8260C | Chloroethane | C1986.D | 36.53 | 20 | TT101D-GW-032415 | UJ |

Notes:

CCV = Continuing calibration verification
 %D = Percent difference
 UJ = Non-detected analyte in associated sample qualified estimated "UJ" due to potential bias.

| Table A-3 Surrogate Non-Conformance | | | | | |
|--|-----------------------|-----|--------|---------------------------------|-----------------------------|
| Method | Analyte | %R | Limits | Associated Sample | Qualifier |
| 8260C | 1,2-Dichloroethane-d4 | 123 | 70-120 | RE103D1-GW-032315 (diluted run) | Trichloroethene qualified J |
| 8260C | 1,2-Dichloroethane-d4 | 126 | 70-120 | RE103D2-GW-032315 (diluted run) | Trichloroethene qualified J |
| 8260C | 1,2-Dichloroethane-d4 | 127 | 70-120 | RE103D3-GW-032315 (diluted run) | Trichloroethene qualified J |
| 8260C | 1,2-Dichloroethane-d4 | 124 | 70-120 | RE122D1-GW-032415 (diluted run) | Trichloroethene qualified J |
| 8260C | 1,2-Dichloroethane-d4 | 122 | 70-120 | RE122D2-GW-032415 (diluted run) | Trichloroethene qualified J |
| 8260C | 1,2-Dichloroethane-d4 | 129 | 70-120 | TT101D1-GW-032415 (diluted run) | Trichloroethene qualified J |
| 8260C | Dibromofluoromethane | 117 | 85-115 | RE103D2-GW-032315 (diluted run) | Trichloroethene qualified J |
| 8260C | Dibromofluoromethane | 116 | 85-115 | RE122D1-GW-032415 (diluted run) | Trichloroethene qualified J |
| 8260C | Dibromofluoromethane | 118 | 85-115 | TT101D1-GW-032415 (diluted run) | Trichloroethene qualified J |

Notes:

%R = Percent recovery
 J = Detected analyte qualified estimated "J" because %R is greater than the upper control limit in associated sample.

| Table A-4 Laboratory Control Sample Non-Conformance | | | | | | |
|--|----------|--------------------|------|--------|-------------------|-----------|
| LCS | Batch | Analyte | %R | Limits | Associated Sample | Qualifier |
| WG160459-1 | WG160459 | Carbon Disulfide | 534 | 35-160 | RE103D1-GW-032315 | J |
| WG160459-1 | WG160459 | Carbon Disulfide | 534 | 35-160 | RE103D2-GW-032315 | J |
| WG160459-1 | WG160459 | Carbon Disulfide | 534 | 35-160 | RE103D3-GW-032315 | J |
| WG160576-1 | WG160576 | Methyl cyclohexane | 50.2 | 73-125 | RE104D2-GW-032315 | UJ |
| WG160576-1 | WG160576 | Methyl cyclohexane | 50.2 | 73-125 | RE122D3-GW-032415 | UJ |
| WG160576-1 | WG160576 | Methyl cyclohexane | 50.2 | 73-125 | TT101D1-GW-032415 | UJ |
| WG160576-1 | WG160576 | Methyl cyclohexane | 50.2 | 73-125 | TT101D2-GW-032415 | UJ |

Notes:

LCS = Laboratory control sample
 %R = Percent recovery
 J = Detected analyte qualified estimated "J" because %R is greater than the upper control limit in associated sample.
 UJ = Non-detected analyte in associated sample qualified estimated "UJ" because %R is lower than lower control limit.

Attachment B
Qualifier Codes and Explanations

| Qualifier | Explanation |
|-----------|--|
| J | The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample. |
| UJ | The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual quantitation limit necessary to accurately and precisely measure the analyte in the sample. |
| U | The analyte was analyzed for, but was not detected above the reported sample quantitation limit. |

Attachment C
Reason Codes and Explanations

| Reason Code | Explanation |
|-------------|---|
| be | Equipment blank contamination |
| bf | Field blank contamination |
| bl | Laboratory blank contamination |
| bt | Trip blank contamination |
| c | Calibration issue |
| d | Reporting limit raised due to chromatographic interference |
| fd | Field duplicate relative percent difference |
| h | Holding times |
| i | Internal standard areas |
| k | Estimated Maximum Possible Concentration |
| l | Laboratory control sample |
| lc | Labeled compound recovery |
| ld | Laboratory duplicate relative percent difference |
| lp | Laboratory control sample/laboratory control sample duplicate relative percent difference |
| m | Matrix spike recovery |
| mc | Method compliance non-conformance |
| md | Matrix spike/matrix spike duplicate relative percent difference |
| nb | Negative laboratory blank contamination |
| p | Chemical preservation issue |
| r | Dual column relative percent difference |
| q | Quantitation issue |
| s | Surrogate recovery |
| su | Ion suppression |
| t | Temperature preservation issue |
| x | Percent solids |
| y | Serial dilution results |
| z | Interference check sample results (metals) |

Attachment D
Final Results after Data Review

| Sample Delivery Group | | | | SI1843 | | |
|-----------------------|---------------------------------------|-----------------|-------|-------------------|------|-----|
| Lab ID | | | | SI1843-1 | | |
| Sample ID | | | | RE103D3-GW-032315 | | |
| Sample Date | | | | 3/23/2015 | | |
| Sample Type | | | | Groundwater | | |
| Method | Analyte | CAS No | Units | Result | Qual | RC |
| 8260C | 1,1,1-TRICHLOROETHANE | 71-55-6 | UG_L | 0.5 | U | |
| 8260C | 1,1,2,2-TETRACHLOROETHANE | 79-34-5 | UG_L | 0.5 | U | |
| 8260C | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 76-13-1 | UG_L | 3.5 | | |
| 8260C | 1,1,2-TRICHLOROETHANE | 79-00-5 | UG_L | 0.5 | U | |
| 8260C | 1,1-DICHLOROETHANE | 75-34-3 | UG_L | 0.5 | J | |
| 8260C | 1,1-DICHLOROETHENE | 75-35-4 | UG_L | 0.69 | J | |
| 8260C | 1,2,4-TRICHLOROBENZENE | 120-82-1 | UG_L | 0.5 | U | |
| 8260C | 1,2-DIBROMO-3-CHLOROPROPANE | 96-12-8 | UG_L | 0.75 | U | |
| 8260C | 1,2-DIBROMOETHANE | 106-93-4 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROBENZENE | 95-50-1 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROETHANE | 107-06-2 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROETHENE, TOTAL | 540-59-0 | UG_L | 1.2 | J | |
| 8260C | 1,2-DICHLOROPROPANE | 78-87-5 | UG_L | 0.5 | U | |
| 8260C | 1,3-DICHLOROBENZENE | 541-73-1 | UG_L | 0.5 | U | |
| 8260C | 1,4-DICHLOROBENZENE | 106-46-7 | UG_L | 0.5 | U | |
| 8260C | 2-BUTANONE | 78-93-3 | UG_L | 2.5 | UJ | c |
| 8260C | 2-HEXANONE | 591-78-6 | UG_L | 2.5 | UJ | c |
| 8260C | 4-METHYL-2-PENTANONE | 108-10-1 | UG_L | 2.5 | U | |
| 8260C | ACETONE | 67-64-1 | UG_L | 2.5 | UJ | c |
| 8260C | BENZENE | 71-43-2 | UG_L | 0.5 | U | |
| 8260C | BROMODICHLOROMETHANE | 75-27-4 | UG_L | 0.5 | U | |
| 8260C | BROMOFORM | 75-25-2 | UG_L | 0.5 | U | |
| 8260C | BROMOMETHANE | 74-83-9 | UG_L | 1 | UJ | c |
| 8260C | CARBON DISULFIDE | 75-15-0 | UG_L | 0.31 | J | l,c |
| 8260C | CARBON TETRACHLORIDE | 56-23-5 | UG_L | 0.5 | U | |
| 8260C | CHLOROBENZENE | 108-90-7 | UG_L | 0.5 | U | |
| 8260C | CHLOROETHANE | 75-00-3 | UG_L | 1 | UJ | c |
| 8260C | CHLOROFORM | 67-66-3 | UG_L | 0.9 | J | |
| 8260C | CHLOROMETHANE | 74-87-3 | UG_L | 1 | UJ | c |
| 8260C | CIS-1,2-DICHLOROETHENE | 156-59-2 | UG_L | 1.2 | | |
| 8260C | CIS-1,3-DICHLOROPROPENE | 10061-01-5 | UG_L | 0.5 | U | |
| 8260C | CYCLOHEXANE | 110-82-7 | UG_L | 0.5 | UJ | c |
| 8260C | DIBROMOCHLOROMETHANE | 124-48-1 | UG_L | 0.5 | U | |
| 8260C | DICHLORODIFLUOROMETHANE | 75-71-8 | UG_L | 1 | UJ | c |
| 8260C | ETHYLBENZENE | 100-41-4 | UG_L | 0.5 | U | |
| 8260C | ISOPROPYLBENZENE | 98-82-8 | UG_L | 0.5 | U | |
| 8260C | M- AND P-XYLENE | 108-38-3/106-42 | UG_L | 1 | U | |
| 8260C | METHYL ACETATE | 79-20-9 | UG_L | 0.75 | U | |
| 8260C | METHYL CYCLOHEXANE | 108-87-2 | UG_L | 0.5 | U | |
| 8260C | METHYL TERT-BUTYL ETHER | 1634-04-4 | UG_L | 0.5 | U | |
| 8260C | METHYLENE CHLORIDE | 75-09-2 | UG_L | 2.5 | U | |
| 8260C | O-XYLENE | 95-47-6 | UG_L | 0.5 | U | |
| 8260C | STYRENE | 100-42-5 | UG_L | 0.5 | U | |
| 8260C | TETRACHLOROETHENE | 127-18-4 | UG_L | 0.47 | J | c |
| 8260C | TOLUENE | 108-88-3 | UG_L | 0.5 | U | |
| 8260C | TRANS-1,2-DICHLOROETHENE | 156-60-5 | UG_L | 0.5 | U | |
| 8260C | TRANS-1,3-DICHLOROPROPENE | 10061-02-6 | UG_L | 0.5 | U | |
| 8260C | TRICHLOROETHENE | 79-01-6 | UG_L | 570 | J | s |
| 8260C | TRICHLOROFLUOROMETHANE | 75-69-4 | UG_L | 1 | U | |
| 8260C | VINYL CHLORIDE | 75-01-4 | UG_L | 1 | U | |
| 8260C | XYLENES, TOTAL | 1330-20-7 | UG_L | 1.5 | U | |
| 8270D_SIM | 1,4-DIOXANE | 123-91-1 | UG_L | 1.3 | | |

| Sample Delivery Group | | | | SI1843 | | |
|-----------------------|---------------------------------------|-----------------|-------|-------------------|------|-----|
| Lab ID | | | | SI1843-2 | | |
| Sample ID | | | | RE103D2-GW-032315 | | |
| Sample Date | | | | 3/23/2015 | | |
| Sample Type | | | | Groundwater | | |
| Method | Analyte | CAS No | Units | Result | Qual | RC |
| 8260C | 1,1,1-TRICHLOROETHANE | 71-55-6 | UG_L | 0.5 | U | |
| 8260C | 1,1,2,2-TETRACHLOROETHANE | 79-34-5 | UG_L | 0.5 | U | |
| 8260C | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 76-13-1 | UG_L | 8.1 | | |
| 8260C | 1,1,2-TRICHLOROETHANE | 79-00-5 | UG_L | 0.54 | J | |
| 8260C | 1,1-DICHLOROETHANE | 75-34-3 | UG_L | 0.92 | J | |
| 8260C | 1,1-DICHLOROETHENE | 75-35-4 | UG_L | 1.6 | | |
| 8260C | 1,2,4-TRICHLOROBENZENE | 120-82-1 | UG_L | 0.5 | U | |
| 8260C | 1,2-DIBROMO-3-CHLOROPROPANE | 96-12-8 | UG_L | 0.75 | U | |
| 8260C | 1,2-DIBROMOETHANE | 106-93-4 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROBENZENE | 95-50-1 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROETHANE | 107-06-2 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROETHENE, TOTAL | 540-59-0 | UG_L | 1.9 | J | |
| 8260C | 1,2-DICHLOROPROPANE | 78-87-5 | UG_L | 0.5 | U | |
| 8260C | 1,3-DICHLOROBENZENE | 541-73-1 | UG_L | 0.5 | U | |
| 8260C | 1,4-DICHLOROBENZENE | 106-46-7 | UG_L | 0.5 | U | |
| 8260C | 2-BUTANONE | 78-93-3 | UG_L | 2.5 | UJ | c |
| 8260C | 2-HEXANONE | 591-78-6 | UG_L | 2.5 | UJ | c |
| 8260C | 4-METHYL-2-PENTANONE | 108-10-1 | UG_L | 2.5 | U | |
| 8260C | ACETONE | 67-64-1 | UG_L | 2.5 | UJ | c |
| 8260C | BENZENE | 71-43-2 | UG_L | 0.5 | U | |
| 8260C | BROMODICHLOROMETHANE | 75-27-4 | UG_L | 0.5 | U | |
| 8260C | BROMOFORM | 75-25-2 | UG_L | 0.5 | U | |
| 8260C | BROMOMETHANE | 74-83-9 | UG_L | 1 | UJ | c |
| 8260C | CARBON DISULFIDE | 75-15-0 | UG_L | 0.47 | J | l,c |
| 8260C | CARBON TETRACHLORIDE | 56-23-5 | UG_L | 0.5 | U | |
| 8260C | CHLOROBENZENE | 108-90-7 | UG_L | 0.5 | U | |
| 8260C | CHLOROETHANE | 75-00-3 | UG_L | 1 | UJ | c |
| 8260C | CHLOROFORM | 67-66-3 | UG_L | 1.2 | | |
| 8260C | CHLOROMETHANE | 74-87-3 | UG_L | 1 | UJ | c |
| 8260C | CIS-1,2-DICHLOROETHENE | 156-59-2 | UG_L | 1.9 | | |
| 8260C | CIS-1,3-DICHLOROPROPENE | 10061-01-5 | UG_L | 0.5 | U | |
| 8260C | CYCLOHEXANE | 110-82-7 | UG_L | 0.5 | UJ | c |
| 8260C | DIBROMOCHLOROMETHANE | 124-48-1 | UG_L | 0.5 | U | |
| 8260C | DICHLORODIFLUOROMETHANE | 75-71-8 | UG_L | 1 | UJ | c |
| 8260C | ETHYLBENZENE | 100-41-4 | UG_L | 0.5 | U | |
| 8260C | ISOPROPYLBENZENE | 98-82-8 | UG_L | 0.5 | U | |
| 8260C | M- AND P-XYLENE | 108-38-3/106-42 | UG_L | 1 | U | |
| 8260C | METHYL ACETATE | 79-20-9 | UG_L | 0.75 | U | |
| 8260C | METHYL CYCLOHEXANE | 108-87-2 | UG_L | 0.5 | U | |
| 8260C | METHYL TERT-BUTYL ETHER | 1634-04-4 | UG_L | 0.5 | U | |
| 8260C | METHYLENE CHLORIDE | 75-09-2 | UG_L | 2.5 | U | |
| 8260C | O-XYLENE | 95-47-6 | UG_L | 0.5 | U | |
| 8260C | STYRENE | 100-42-5 | UG_L | 0.5 | U | |
| 8260C | TETRACHLOROETHENE | 127-18-4 | UG_L | 1.1 | J | c |
| 8260C | TOLUENE | 108-88-3 | UG_L | 0.5 | U | |
| 8260C | TRANS-1,2-DICHLOROETHENE | 156-60-5 | UG_L | 0.5 | U | |
| 8260C | TRANS-1,3-DICHLOROPROPENE | 10061-02-6 | UG_L | 0.5 | U | |
| 8260C | TRICHLOROETHENE | 79-01-6 | UG_L | 940 | J | s |
| 8260C | TRICHLOROFLUOROMETHANE | 75-69-4 | UG_L | 1 | U | |
| 8260C | VINYL CHLORIDE | 75-01-4 | UG_L | 1 | U | |
| 8260C | XYLENES, TOTAL | 1330-20-7 | UG_L | 1.5 | U | |
| 8270D_SIM | 1,4-DIOXANE | 123-91-1 | UG_L | 3 | | |

| Sample Delivery Group | | | | SI1843 | | |
|-----------------------|---------------------------------------|-----------------|-------|-------------------|------|-----|
| Lab ID | | | | SI1843-3 | | |
| Sample ID | | | | RE103D1-GW-032315 | | |
| Sample Date | | | | 3/23/2015 | | |
| Sample Type | | | | Groundwater | | |
| Method | Analyte | CAS No | Units | Result | Qual | RC |
| 8260C | 1,1,1-TRICHLOROETHANE | 71-55-6 | UG_L | 0.53 | J | |
| 8260C | 1,1,2,2-TETRACHLOROETHANE | 79-34-5 | UG_L | 0.5 | U | |
| 8260C | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 76-13-1 | UG_L | 16 | | |
| 8260C | 1,1,2-TRICHLOROETHANE | 79-00-5 | UG_L | 0.77 | J | |
| 8260C | 1,1-DICHLOROETHANE | 75-34-3 | UG_L | 1.1 | | |
| 8260C | 1,1-DICHLOROETHENE | 75-35-4 | UG_L | 6.8 | | |
| 8260C | 1,2,4-TRICHLOROBENZENE | 120-82-1 | UG_L | 0.5 | U | |
| 8260C | 1,2-DIBROMO-3-CHLOROPROPANE | 96-12-8 | UG_L | 0.75 | U | |
| 8260C | 1,2-DIBROMOETHANE | 106-93-4 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROBENZENE | 95-50-1 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROETHANE | 107-06-2 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROETHENE, TOTAL | 540-59-0 | UG_L | 3.9 | | |
| 8260C | 1,2-DICHLOROPROPANE | 78-87-5 | UG_L | 0.5 | U | |
| 8260C | 1,3-DICHLOROBENZENE | 541-73-1 | UG_L | 0.5 | U | |
| 8260C | 1,4-DICHLOROBENZENE | 106-46-7 | UG_L | 0.5 | U | |
| 8260C | 2-BUTANONE | 78-93-3 | UG_L | 2.5 | UJ | c |
| 8260C | 2-HEXANONE | 591-78-6 | UG_L | 2.5 | UJ | c |
| 8260C | 4-METHYL-2-PENTANONE | 108-10-1 | UG_L | 2.5 | U | |
| 8260C | ACETONE | 67-64-1 | UG_L | 2.5 | UJ | c |
| 8260C | BENZENE | 71-43-2 | UG_L | 0.5 | U | |
| 8260C | BROMODICHLOROMETHANE | 75-27-4 | UG_L | 0.5 | U | |
| 8260C | BROMOFORM | 75-25-2 | UG_L | 0.5 | U | |
| 8260C | BROMOMETHANE | 74-83-9 | UG_L | 1 | UJ | c |
| 8260C | CARBON DISULFIDE | 75-15-0 | UG_L | 0.56 | J | l,c |
| 8260C | CARBON TETRACHLORIDE | 56-23-5 | UG_L | 0.5 | U | |
| 8260C | CHLOROBENZENE | 108-90-7 | UG_L | 0.5 | U | |
| 8260C | CHLOROETHANE | 75-00-3 | UG_L | 1 | UJ | c |
| 8260C | CHLOROFORM | 67-66-3 | UG_L | 0.76 | J | |
| 8260C | CHLOROMETHANE | 74-87-3 | UG_L | 1 | UJ | c |
| 8260C | CIS-1,2-DICHLOROETHENE | 156-59-2 | UG_L | 3.9 | | |
| 8260C | CIS-1,3-DICHLOROPROPENE | 10061-01-5 | UG_L | 0.5 | U | |
| 8260C | CYCLOHEXANE | 110-82-7 | UG_L | 0.5 | UJ | c |
| 8260C | DIBROMOCHLOROMETHANE | 124-48-1 | UG_L | 0.5 | U | |
| 8260C | DICHLORODIFLUOROMETHANE | 75-71-8 | UG_L | 1 | UJ | c |
| 8260C | ETHYLBENZENE | 100-41-4 | UG_L | 0.5 | U | |
| 8260C | ISOPROPYLBENZENE | 98-82-8 | UG_L | 0.5 | U | |
| 8260C | M- AND P-XYLENE | 108-38-3/106-42 | UG_L | 1 | U | |
| 8260C | METHYL ACETATE | 79-20-9 | UG_L | 0.75 | U | |
| 8260C | METHYL CYCLOHEXANE | 108-87-2 | UG_L | 0.5 | U | |
| 8260C | METHYL TERT-BUTYL ETHER | 1634-04-4 | UG_L | 0.5 | U | |
| 8260C | METHYLENE CHLORIDE | 75-09-2 | UG_L | 2.5 | U | |
| 8260C | O-XYLENE | 95-47-6 | UG_L | 0.5 | U | |
| 8260C | STYRENE | 100-42-5 | UG_L | 0.5 | U | |
| 8260C | TETRACHLOROETHENE | 127-18-4 | UG_L | 4.6 | J | c |
| 8260C | TOLUENE | 108-88-3 | UG_L | 0.5 | U | |
| 8260C | TRANS-1,2-DICHLOROETHENE | 156-60-5 | UG_L | 0.5 | U | |
| 8260C | TRANS-1,3-DICHLOROPROPENE | 10061-02-6 | UG_L | 0.5 | U | |
| 8260C | TRICHLOROETHENE | 79-01-6 | UG_L | 900 | J | s |
| 8260C | TRICHLOROFLUOROMETHANE | 75-69-4 | UG_L | 1 | U | |
| 8260C | VINYL CHLORIDE | 75-01-4 | UG_L | 1 | U | |
| 8260C | XYLENES, TOTAL | 1330-20-7 | UG_L | 1.5 | U | |
| 8270D_SIM | 1,4-DIOXANE | 123-91-1 | UG_L | 19 | | |

| Sample Delivery Group | | | | SI1843 | | |
|-----------------------|---------------------------------------|-----------------|-------|-------------------|------|----|
| Lab ID | | | | SI1843-5 | | |
| Sample ID | | | | RE104D1-GW-032315 | | |
| Sample Date | | | | 3/23/2015 | | |
| Sample Type | | | | Groundwater | | |
| Method | Analyte | CAS No | Units | Result | Qual | RC |
| 8260C | 1,1,1-TRICHLOROETHANE | 71-55-6 | UG_L | 0.27 | J | |
| 8260C | 1,1,2,2-TETRACHLOROETHANE | 79-34-5 | UG_L | 0.5 | U | |
| 8260C | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 76-13-1 | UG_L | 6.2 | | |
| 8260C | 1,1,2-TRICHLOROETHANE | 79-00-5 | UG_L | 0.5 | U | |
| 8260C | 1,1-DICHLOROETHANE | 75-34-3 | UG_L | 0.5 | U | |
| 8260C | 1,1-DICHLOROETHENE | 75-35-4 | UG_L | 0.8 | J | |
| 8260C | 1,2,4-TRICHLOROBENZENE | 120-82-1 | UG_L | 0.5 | U | |
| 8260C | 1,2-DIBROMO-3-CHLOROPROPANE | 96-12-8 | UG_L | 0.75 | U | |
| 8260C | 1,2-DIBROMOETHANE | 106-93-4 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROBENZENE | 95-50-1 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROETHANE | 107-06-2 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROETHENE, TOTAL | 540-59-0 | UG_L | 1.5 | J | |
| 8260C | 1,2-DICHLOROPROPANE | 78-87-5 | UG_L | 0.5 | U | |
| 8260C | 1,3-DICHLOROBENZENE | 541-73-1 | UG_L | 0.5 | U | |
| 8260C | 1,4-DICHLOROBENZENE | 106-46-7 | UG_L | 0.5 | U | |
| 8260C | 2-BUTANONE | 78-93-3 | UG_L | 2.5 | UJ | c |
| 8260C | 2-HEXANONE | 591-78-6 | UG_L | 2.5 | UJ | c |
| 8260C | 4-METHYL-2-PENTANONE | 108-10-1 | UG_L | 2.5 | U | |
| 8260C | ACETONE | 67-64-1 | UG_L | 2.5 | UJ | c |
| 8260C | BENZENE | 71-43-2 | UG_L | 0.5 | U | |
| 8260C | BROMODICHLOROMETHANE | 75-27-4 | UG_L | 0.5 | U | |
| 8260C | BROMOFORM | 75-25-2 | UG_L | 0.5 | U | |
| 8260C | BROMOMETHANE | 74-83-9 | UG_L | 1 | UJ | c |
| 8260C | CARBON DISULFIDE | 75-15-0 | UG_L | 0.5 | UJ | c |
| 8260C | CARBON TETRACHLORIDE | 56-23-5 | UG_L | 0.5 | U | |
| 8260C | CHLOROBENZENE | 108-90-7 | UG_L | 0.5 | U | |
| 8260C | CHLOROETHANE | 75-00-3 | UG_L | 1 | UJ | c |
| 8260C | CHLOROFORM | 67-66-3 | UG_L | 0.5 | U | |
| 8260C | CHLOROMETHANE | 74-87-3 | UG_L | 1 | UJ | c |
| 8260C | CIS-1,2-DICHLOROETHENE | 156-59-2 | UG_L | 1.5 | | |
| 8260C | CIS-1,3-DICHLOROPROPENE | 10061-01-5 | UG_L | 0.5 | U | |
| 8260C | CYCLOHEXANE | 110-82-7 | UG_L | 0.5 | UJ | c |
| 8260C | DIBROMOCHLOROMETHANE | 124-48-1 | UG_L | 0.5 | U | |
| 8260C | DICHLORODIFLUOROMETHANE | 75-71-8 | UG_L | 1 | UJ | c |
| 8260C | ETHYLBENZENE | 100-41-4 | UG_L | 0.5 | U | |
| 8260C | ISOPROPYLBENZENE | 98-82-8 | UG_L | 0.5 | U | |
| 8260C | M- AND P-XYLENE | 108-38-3/106-42 | UG_L | 1 | U | |
| 8260C | METHYL ACETATE | 79-20-9 | UG_L | 0.75 | U | |
| 8260C | METHYL CYCLOHEXANE | 108-87-2 | UG_L | 0.5 | U | |
| 8260C | METHYL TERT-BUTYL ETHER | 1634-04-4 | UG_L | 0.5 | U | |
| 8260C | METHYLENE CHLORIDE | 75-09-2 | UG_L | 2.5 | U | |
| 8260C | O-XYLENE | 95-47-6 | UG_L | 0.5 | U | |
| 8260C | STYRENE | 100-42-5 | UG_L | 0.5 | U | |
| 8260C | TETRACHLOROETHENE | 127-18-4 | UG_L | 2.4 | J | c |
| 8260C | TOLUENE | 108-88-3 | UG_L | 0.5 | U | |
| 8260C | TRANS-1,2-DICHLOROETHENE | 156-60-5 | UG_L | 0.5 | U | |
| 8260C | TRANS-1,3-DICHLOROPROPENE | 10061-02-6 | UG_L | 0.5 | U | |
| 8260C | TRICHLOROETHENE | 79-01-6 | UG_L | 110 | | |
| 8260C | TRICHLOROFLUOROMETHANE | 75-69-4 | UG_L | 1 | U | |
| 8260C | VINYL CHLORIDE | 75-01-4 | UG_L | 1 | U | |
| 8260C | XYLENES, TOTAL | 1330-20-7 | UG_L | 1.5 | U | |
| 8270D_SIM | 1,4-DIOXANE | 123-91-1 | UG_L | 9.7 | | |

| Sample Delivery Group | | | | SI1843 | | |
|-----------------------|---------------------------------------|-----------------|-------|-------------------|------|-----|
| Lab ID | | | | SI1843-6 | | |
| Sample ID | | | | RE104D2-GW-032315 | | |
| Sample Date | | | | 3/23/2015 | | |
| Sample Type | | | | Groundwater | | |
| Method | Analyte | CAS No | Units | Result | Qual | RC |
| 8260C | 1,1,1-TRICHLOROETHANE | 71-55-6 | UG_L | 0.5 | U | |
| 8260C | 1,1,2,2-TETRACHLOROETHANE | 79-34-5 | UG_L | 0.5 | U | |
| 8260C | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 76-13-1 | UG_L | 0.5 | U | |
| 8260C | 1,1,2-TRICHLOROETHANE | 79-00-5 | UG_L | 0.5 | U | |
| 8260C | 1,1-DICHLOROETHANE | 75-34-3 | UG_L | 0.5 | U | |
| 8260C | 1,1-DICHLOROETHENE | 75-35-4 | UG_L | 0.5 | U | |
| 8260C | 1,2,4-TRICHLOROBENZENE | 120-82-1 | UG_L | 0.5 | U | |
| 8260C | 1,2-DIBROMO-3-CHLOROPROPANE | 96-12-8 | UG_L | 0.75 | U | |
| 8260C | 1,2-DIBROMOETHANE | 106-93-4 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROBENZENE | 95-50-1 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROETHANE | 107-06-2 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROETHENE, TOTAL | 540-59-0 | UG_L | 1.2 | J | |
| 8260C | 1,2-DICHLOROPROPANE | 78-87-5 | UG_L | 0.5 | U | |
| 8260C | 1,3-DICHLOROBENZENE | 541-73-1 | UG_L | 0.5 | U | |
| 8260C | 1,4-DICHLOROBENZENE | 106-46-7 | UG_L | 0.5 | U | |
| 8260C | 2-BUTANONE | 78-93-3 | UG_L | 2.5 | UJ | c |
| 8260C | 2-HEXANONE | 591-78-6 | UG_L | 2.5 | UJ | c |
| 8260C | 4-METHYL-2-PENTANONE | 108-10-1 | UG_L | 2.5 | U | |
| 8260C | ACETONE | 67-64-1 | UG_L | 2.5 | UJ | c |
| 8260C | BENZENE | 71-43-2 | UG_L | 0.5 | U | |
| 8260C | BROMODICHLOROMETHANE | 75-27-4 | UG_L | 0.5 | U | |
| 8260C | BROMOFORM | 75-25-2 | UG_L | 0.5 | U | |
| 8260C | BROMOMETHANE | 74-83-9 | UG_L | 1 | UJ | c |
| 8260C | CARBON DISULFIDE | 75-15-0 | UG_L | 0.5 | UJ | c |
| 8260C | CARBON TETRACHLORIDE | 56-23-5 | UG_L | 0.5 | U | |
| 8260C | CHLOROBENZENE | 108-90-7 | UG_L | 0.5 | U | |
| 8260C | CHLOROETHANE | 75-00-3 | UG_L | 1 | UJ | c |
| 8260C | CHLOROFORM | 67-66-3 | UG_L | 0.5 | U | |
| 8260C | CHLOROMETHANE | 74-87-3 | UG_L | 1 | UJ | c |
| 8260C | CIS-1,2-DICHLOROETHENE | 156-59-2 | UG_L | 1.2 | | |
| 8260C | CIS-1,3-DICHLOROPROPENE | 10061-01-5 | UG_L | 0.5 | U | |
| 8260C | CYCLOHEXANE | 110-82-7 | UG_L | 0.5 | UJ | c |
| 8260C | DIBROMOCHLOROMETHANE | 124-48-1 | UG_L | 0.5 | U | |
| 8260C | DICHLORODIFLUOROMETHANE | 75-71-8 | UG_L | 1 | UJ | c |
| 8260C | ETHYLBENZENE | 100-41-4 | UG_L | 0.5 | U | |
| 8260C | ISOPROPYLBENZENE | 98-82-8 | UG_L | 0.5 | U | |
| 8260C | M- AND P-XYLENE | 108-38-3/106-42 | UG_L | 1 | U | |
| 8260C | METHYL ACETATE | 79-20-9 | UG_L | 0.75 | U | |
| 8260C | METHYL CYCLOHEXANE | 108-87-2 | UG_L | 0.5 | UJ | l,c |
| 8260C | METHYL TERT-BUTYL ETHER | 1634-04-4 | UG_L | 0.5 | U | |
| 8260C | METHYLENE CHLORIDE | 75-09-2 | UG_L | 2.5 | U | |
| 8260C | O-XYLENE | 95-47-6 | UG_L | 0.5 | U | |
| 8260C | STYRENE | 100-42-5 | UG_L | 0.5 | U | |
| 8260C | TETRACHLOROETHENE | 127-18-4 | UG_L | 0.5 | UJ | c |
| 8260C | TOLUENE | 108-88-3 | UG_L | 0.5 | U | |
| 8260C | TRANS-1,2-DICHLOROETHENE | 156-60-5 | UG_L | 0.5 | U | |
| 8260C | TRANS-1,3-DICHLOROPROPENE | 10061-02-6 | UG_L | 0.5 | U | |
| 8260C | TRICHLOROETHENE | 79-01-6 | UG_L | 3 | | |
| 8260C | TRICHLOROFLUOROMETHANE | 75-69-4 | UG_L | 1 | U | |
| 8260C | VINYL CHLORIDE | 75-01-4 | UG_L | 1 | U | |
| 8260C | XYLENES, TOTAL | 1330-20-7 | UG_L | 1.5 | U | |
| 8270D_SIM | 1,4-DIOXANE | 123-91-1 | UG_L | 0.12 | J | |

| Sample Delivery Group | | | | SI1843 | | |
|-----------------------|---------------------------------------|-----------------|-------|-------------------|------|----|
| Lab ID | | | | SI1843-7 | | |
| Sample ID | | | | RE104D3-GW-032315 | | |
| Sample Date | | | | 3/23/2015 | | |
| Sample Type | | | | Groundwater | | |
| Method | Analyte | CAS No | Units | Result | Qual | RC |
| 8260C | 1,1,1-TRICHLOROETHANE | 71-55-6 | UG_L | 0.5 | U | |
| 8260C | 1,1,2,2-TETRACHLOROETHANE | 79-34-5 | UG_L | 0.5 | U | |
| 8260C | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 76-13-1 | UG_L | 0.5 | U | |
| 8260C | 1,1,2-TRICHLOROETHANE | 79-00-5 | UG_L | 0.5 | U | |
| 8260C | 1,1-DICHLOROETHANE | 75-34-3 | UG_L | 0.5 | U | |
| 8260C | 1,1-DICHLOROETHENE | 75-35-4 | UG_L | 0.5 | U | |
| 8260C | 1,2,4-TRICHLOROBENZENE | 120-82-1 | UG_L | 0.5 | U | |
| 8260C | 1,2-DIBROMO-3-CHLOROPROPANE | 96-12-8 | UG_L | 0.75 | U | |
| 8260C | 1,2-DIBROMOETHANE | 106-93-4 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROBENZENE | 95-50-1 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROETHANE | 107-06-2 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROETHENE, TOTAL | 540-59-0 | UG_L | 1 | U | |
| 8260C | 1,2-DICHLOROPROPANE | 78-87-5 | UG_L | 0.5 | U | |
| 8260C | 1,3-DICHLOROBENZENE | 541-73-1 | UG_L | 0.5 | U | |
| 8260C | 1,4-DICHLOROBENZENE | 106-46-7 | UG_L | 0.5 | U | |
| 8260C | 2-BUTANONE | 78-93-3 | UG_L | 2.5 | UJ | c |
| 8260C | 2-HEXANONE | 591-78-6 | UG_L | 2.5 | UJ | c |
| 8260C | 4-METHYL-2-PENTANONE | 108-10-1 | UG_L | 2.5 | U | |
| 8260C | ACETONE | 67-64-1 | UG_L | 2.5 | UJ | c |
| 8260C | BENZENE | 71-43-2 | UG_L | 0.5 | U | |
| 8260C | BROMODICHLOROMETHANE | 75-27-4 | UG_L | 0.5 | U | |
| 8260C | BROMOFORM | 75-25-2 | UG_L | 0.5 | U | |
| 8260C | BROMOMETHANE | 74-83-9 | UG_L | 1 | UJ | c |
| 8260C | CARBON DISULFIDE | 75-15-0 | UG_L | 0.5 | UJ | c |
| 8260C | CARBON TETRACHLORIDE | 56-23-5 | UG_L | 0.5 | U | |
| 8260C | CHLOROBENZENE | 108-90-7 | UG_L | 0.5 | U | |
| 8260C | CHLOROETHANE | 75-00-3 | UG_L | 1 | U | |
| 8260C | CHLOROFORM | 67-66-3 | UG_L | 0.5 | U | |
| 8260C | CHLOROMETHANE | 74-87-3 | UG_L | 1 | UJ | c |
| 8260C | CIS-1,2-DICHLOROETHENE | 156-59-2 | UG_L | 0.5 | U | |
| 8260C | CIS-1,3-DICHLOROPROPENE | 10061-01-5 | UG_L | 0.5 | U | |
| 8260C | CYCLOHEXANE | 110-82-7 | UG_L | 0.5 | UJ | c |
| 8260C | DIBROMOCHLOROMETHANE | 124-48-1 | UG_L | 0.5 | U | |
| 8260C | DICHLORODIFLUOROMETHANE | 75-71-8 | UG_L | 1 | UJ | c |
| 8260C | ETHYLBENZENE | 100-41-4 | UG_L | 0.5 | U | |
| 8260C | ISOPROPYLBENZENE | 98-82-8 | UG_L | 0.5 | U | |
| 8260C | M- AND P-XYLENE | 108-38-3/106-42 | UG_L | 1 | U | |
| 8260C | METHYL ACETATE | 79-20-9 | UG_L | 0.75 | U | |
| 8260C | METHYL CYCLOHEXANE | 108-87-2 | UG_L | 0.5 | U | |
| 8260C | METHYL TERT-BUTYL ETHER | 1634-04-4 | UG_L | 0.5 | U | |
| 8260C | METHYLENE CHLORIDE | 75-09-2 | UG_L | 2.5 | U | |
| 8260C | O-XYLENE | 95-47-6 | UG_L | 0.5 | U | |
| 8260C | STYRENE | 100-42-5 | UG_L | 0.5 | U | |
| 8260C | TETRACHLOROETHENE | 127-18-4 | UG_L | 0.5 | UJ | c |
| 8260C | TOLUENE | 108-88-3 | UG_L | 0.5 | U | |
| 8260C | TRANS-1,2-DICHLOROETHENE | 156-60-5 | UG_L | 0.5 | U | |
| 8260C | TRANS-1,3-DICHLOROPROPENE | 10061-02-6 | UG_L | 0.5 | U | |
| 8260C | TRICHLOROETHENE | 79-01-6 | UG_L | 0.46 | J | |
| 8260C | TRICHLOROFLUOROMETHANE | 75-69-4 | UG_L | 1 | U | |
| 8260C | VINYL CHLORIDE | 75-01-4 | UG_L | 1 | U | |
| 8260C | XYLENES, TOTAL | 1330-20-7 | UG_L | 1.5 | U | |
| 8270D_SIM | 1,4-DIOXANE | 123-91-1 | UG_L | 0.19 | U | |

| Sample Delivery Group | | | | SI1843 | | |
|-----------------------|---------------------------------------|-----------------|-------|---------------------|------|----|
| Lab ID | | | | SI1843-8 | | |
| Sample ID | | | | DUPLICATE-GW-032315 | | |
| Sample Date | | | | 3/23/2015 | | |
| Sample Type | | | | Field Duplicate | | |
| Method | Analyte | CAS No | Units | Result | Qual | RC |
| 8260C | 1,1,1-TRICHLOROETHANE | 71-55-6 | UG_L | 0.5 | U | |
| 8260C | 1,1,2,2-TETRACHLOROETHANE | 79-34-5 | UG_L | 0.5 | U | |
| 8260C | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 76-13-1 | UG_L | 0.5 | U | |
| 8260C | 1,1,2-TRICHLOROETHANE | 79-00-5 | UG_L | 0.5 | U | |
| 8260C | 1,1-DICHLOROETHANE | 75-34-3 | UG_L | 0.5 | U | |
| 8260C | 1,1-DICHLOROETHENE | 75-35-4 | UG_L | 0.5 | U | |
| 8260C | 1,2,4-TRICHLOROBENZENE | 120-82-1 | UG_L | 0.5 | U | |
| 8260C | 1,2-DIBROMO-3-CHLOROPROPANE | 96-12-8 | UG_L | 0.75 | U | |
| 8260C | 1,2-DIBROMOETHANE | 106-93-4 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROBENZENE | 95-50-1 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROETHANE | 107-06-2 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROETHENE, TOTAL | 540-59-0 | UG_L | 1.3 | J | |
| 8260C | 1,2-DICHLOROPROPANE | 78-87-5 | UG_L | 0.5 | U | |
| 8260C | 1,3-DICHLOROBENZENE | 541-73-1 | UG_L | 0.5 | U | |
| 8260C | 1,4-DICHLOROBENZENE | 106-46-7 | UG_L | 0.5 | U | |
| 8260C | 2-BUTANONE | 78-93-3 | UG_L | 2.5 | UJ | c |
| 8260C | 2-HEXANONE | 591-78-6 | UG_L | 2.5 | UJ | c |
| 8260C | 4-METHYL-2-PENTANONE | 108-10-1 | UG_L | 2.5 | U | |
| 8260C | ACETONE | 67-64-1 | UG_L | 2.5 | UJ | c |
| 8260C | BENZENE | 71-43-2 | UG_L | 0.5 | U | |
| 8260C | BROMODICHLOROMETHANE | 75-27-4 | UG_L | 0.5 | U | |
| 8260C | BROMOFORM | 75-25-2 | UG_L | 0.5 | U | |
| 8260C | BROMOMETHANE | 74-83-9 | UG_L | 1 | UJ | c |
| 8260C | CARBON DISULFIDE | 75-15-0 | UG_L | 0.5 | UJ | c |
| 8260C | CARBON TETRACHLORIDE | 56-23-5 | UG_L | 0.5 | U | |
| 8260C | CHLOROBENZENE | 108-90-7 | UG_L | 0.5 | U | |
| 8260C | CHLOROETHANE | 75-00-3 | UG_L | 1 | UJ | c |
| 8260C | CHLOROFORM | 67-66-3 | UG_L | 0.5 | U | |
| 8260C | CHLOROMETHANE | 74-87-3 | UG_L | 1 | UJ | c |
| 8260C | CIS-1,2-DICHLOROETHENE | 156-59-2 | UG_L | 1.3 | | |
| 8260C | CIS-1,3-DICHLOROPROPENE | 10061-01-5 | UG_L | 0.5 | U | |
| 8260C | CYCLOHEXANE | 110-82-7 | UG_L | 0.5 | UJ | c |
| 8260C | DIBROMOCHLOROMETHANE | 124-48-1 | UG_L | 0.5 | U | |
| 8260C | DICHLORODIFLUOROMETHANE | 75-71-8 | UG_L | 1 | UJ | c |
| 8260C | ETHYLBENZENE | 100-41-4 | UG_L | 0.5 | U | |
| 8260C | ISOPROPYLBENZENE | 98-82-8 | UG_L | 0.5 | U | |
| 8260C | M- AND P-XYLENE | 108-38-3/106-42 | UG_L | 1 | U | |
| 8260C | METHYL ACETATE | 79-20-9 | UG_L | 0.75 | U | |
| 8260C | METHYL CYCLOHEXANE | 108-87-2 | UG_L | 0.5 | U | |
| 8260C | METHYL TERT-BUTYL ETHER | 1634-04-4 | UG_L | 0.5 | U | |
| 8260C | METHYLENE CHLORIDE | 75-09-2 | UG_L | 2.5 | U | |
| 8260C | O-XYLENE | 95-47-6 | UG_L | 0.5 | U | |
| 8260C | STYRENE | 100-42-5 | UG_L | 0.5 | U | |
| 8260C | TETRACHLOROETHENE | 127-18-4 | UG_L | 0.5 | UJ | c |
| 8260C | TOLUENE | 108-88-3 | UG_L | 0.5 | U | |
| 8260C | TRANS-1,2-DICHLOROETHENE | 156-60-5 | UG_L | 0.5 | U | |
| 8260C | TRANS-1,3-DICHLOROPROPENE | 10061-02-6 | UG_L | 0.5 | U | |
| 8260C | TRICHLOROETHENE | 79-01-6 | UG_L | 3.1 | | |
| 8260C | TRICHLOROFLUOROMETHANE | 75-69-4 | UG_L | 1 | U | |
| 8260C | VINYL CHLORIDE | 75-01-4 | UG_L | 1 | U | |
| 8260C | XYLENES, TOTAL | 1330-20-7 | UG_L | 1.5 | U | |
| 8270D_SIM | 1,4-DIOXANE | 123-91-1 | UG_L | 0.096 | J | |

| Sample Delivery Group | | | | SI1843 | | |
|-----------------------|---------------------------------------|-----------------|-------|-------------------|------|----|
| Lab ID | | | | SI1843-9 | | |
| Sample ID | | | | RE122D1-GW-032415 | | |
| Sample Date | | | | 3/24/2015 | | |
| Sample Type | | | | Groundwater | | |
| Method | Analyte | CAS No | Units | Result | Qual | RC |
| 8260C | 1,1,1-TRICHLOROETHANE | 71-55-6 | UG_L | 0.5 | U | |
| 8260C | 1,1,2,2-TETRACHLOROETHANE | 79-34-5 | UG_L | 0.5 | U | |
| 8260C | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 76-13-1 | UG_L | 7.2 | | |
| 8260C | 1,1,2-TRICHLOROETHANE | 79-00-5 | UG_L | 0.4 | J | |
| 8260C | 1,1-DICHLOROETHANE | 75-34-3 | UG_L | 0.5 | U | |
| 8260C | 1,1-DICHLOROETHENE | 75-35-4 | UG_L | 0.5 | U | |
| 8260C | 1,2,4-TRICHLOROBENZENE | 120-82-1 | UG_L | 0.5 | U | |
| 8260C | 1,2-DIBROMO-3-CHLOROPROPANE | 96-12-8 | UG_L | 0.75 | U | |
| 8260C | 1,2-DIBROMOETHANE | 106-93-4 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROBENZENE | 95-50-1 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROETHANE | 107-06-2 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROETHENE, TOTAL | 540-59-0 | UG_L | 2 | | |
| 8260C | 1,2-DICHLOROPROPANE | 78-87-5 | UG_L | 0.5 | U | |
| 8260C | 1,3-DICHLOROBENZENE | 541-73-1 | UG_L | 0.5 | U | |
| 8260C | 1,4-DICHLOROBENZENE | 106-46-7 | UG_L | 0.5 | U | |
| 8260C | 2-BUTANONE | 78-93-3 | UG_L | 2.5 | UJ | c |
| 8260C | 2-HEXANONE | 591-78-6 | UG_L | 2.5 | UJ | c |
| 8260C | 4-METHYL-2-PENTANONE | 108-10-1 | UG_L | 2.5 | U | |
| 8260C | ACETONE | 67-64-1 | UG_L | 2.5 | UJ | c |
| 8260C | BENZENE | 71-43-2 | UG_L | 0.5 | U | |
| 8260C | BROMODICHLOROMETHANE | 75-27-4 | UG_L | 0.5 | U | |
| 8260C | BROMOFORM | 75-25-2 | UG_L | 0.5 | U | |
| 8260C | BROMOMETHANE | 74-83-9 | UG_L | 1 | UJ | c |
| 8260C | CARBON DISULFIDE | 75-15-0 | UG_L | 0.5 | UJ | c |
| 8260C | CARBON TETRACHLORIDE | 56-23-5 | UG_L | 0.5 | U | |
| 8260C | CHLOROBENZENE | 108-90-7 | UG_L | 0.5 | U | |
| 8260C | CHLOROETHANE | 75-00-3 | UG_L | 1 | UJ | c |
| 8260C | CHLOROFORM | 67-66-3 | UG_L | 0.62 | J | |
| 8260C | CHLOROMETHANE | 74-87-3 | UG_L | 1 | UJ | c |
| 8260C | CIS-1,2-DICHLOROETHENE | 156-59-2 | UG_L | 2 | | |
| 8260C | CIS-1,3-DICHLOROPROPENE | 10061-01-5 | UG_L | 0.5 | U | |
| 8260C | CYCLOHEXANE | 110-82-7 | UG_L | 0.5 | UJ | c |
| 8260C | DIBROMOCHLOROMETHANE | 124-48-1 | UG_L | 0.5 | U | |
| 8260C | DICHLORODIFLUOROMETHANE | 75-71-8 | UG_L | 1 | UJ | c |
| 8260C | ETHYLBENZENE | 100-41-4 | UG_L | 0.5 | U | |
| 8260C | ISOPROPYLBENZENE | 98-82-8 | UG_L | 0.5 | U | |
| 8260C | M- AND P-XYLENE | 108-38-3/106-42 | UG_L | 1 | U | |
| 8260C | METHYL ACETATE | 79-20-9 | UG_L | 0.75 | U | |
| 8260C | METHYL CYCLOHEXANE | 108-87-2 | UG_L | 0.5 | U | |
| 8260C | METHYL TERT-BUTYL ETHER | 1634-04-4 | UG_L | 0.5 | U | |
| 8260C | METHYLENE CHLORIDE | 75-09-2 | UG_L | 2.5 | U | |
| 8260C | O-XYLENE | 95-47-6 | UG_L | 0.5 | U | |
| 8260C | STYRENE | 100-42-5 | UG_L | 0.5 | U | |
| 8260C | TETRACHLOROETHENE | 127-18-4 | UG_L | 1.3 | J | c |
| 8260C | TOLUENE | 108-88-3 | UG_L | 0.5 | U | |
| 8260C | TRANS-1,2-DICHLOROETHENE | 156-60-5 | UG_L | 0.5 | U | |
| 8260C | TRANS-1,3-DICHLOROPROPENE | 10061-02-6 | UG_L | 0.5 | U | |
| 8260C | TRICHLOROETHENE | 79-01-6 | UG_L | 570 | J | s |
| 8260C | TRICHLOROFLUOROMETHANE | 75-69-4 | UG_L | 1 | U | |
| 8260C | VINYL CHLORIDE | 75-01-4 | UG_L | 1 | U | |
| 8260C | XYLENES, TOTAL | 1330-20-7 | UG_L | 1.5 | U | |
| 8270D_SIM | 1,4-DIOXANE | 123-91-1 | UG_L | 8.1 | | |

| Sample Delivery Group | | | | SI1843 | | |
|-----------------------|---------------------------------------|-----------------|-------|-------------------|------|----|
| Lab ID | | | | SI1843-10 | | |
| Sample ID | | | | RE122D2-GW-032415 | | |
| Sample Date | | | | 3/24/2015 | | |
| Sample Type | | | | Groundwater | | |
| Method | Analyte | CAS No | Units | Result | Qual | RC |
| 8260C | 1,1,1-TRICHLOROETHANE | 71-55-6 | UG_L | 0.71 | J | |
| 8260C | 1,1,2,2-TETRACHLOROETHANE | 79-34-5 | UG_L | 0.5 | U | |
| 8260C | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 76-13-1 | UG_L | 31 | | |
| 8260C | 1,1,2-TRICHLOROETHANE | 79-00-5 | UG_L | 2.8 | | |
| 8260C | 1,1-DICHLOROETHANE | 75-34-3 | UG_L | 1.7 | | |
| 8260C | 1,1-DICHLOROETHENE | 75-35-4 | UG_L | 8.7 | | |
| 8260C | 1,2,4-TRICHLOROBENZENE | 120-82-1 | UG_L | 0.5 | U | |
| 8260C | 1,2-DIBROMO-3-CHLOROPROPANE | 96-12-8 | UG_L | 0.75 | U | |
| 8260C | 1,2-DIBROMOETHANE | 106-93-4 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROBENZENE | 95-50-1 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROETHANE | 107-06-2 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROETHENE, TOTAL | 540-59-0 | UG_L | 6 | | |
| 8260C | 1,2-DICHLOROPROPANE | 78-87-5 | UG_L | 0.5 | U | |
| 8260C | 1,3-DICHLOROBENZENE | 541-73-1 | UG_L | 0.5 | U | |
| 8260C | 1,4-DICHLOROBENZENE | 106-46-7 | UG_L | 0.5 | U | |
| 8260C | 2-BUTANONE | 78-93-3 | UG_L | 2.5 | UJ | c |
| 8260C | 2-HEXANONE | 591-78-6 | UG_L | 2.5 | UJ | c |
| 8260C | 4-METHYL-2-PENTANONE | 108-10-1 | UG_L | 2.5 | U | |
| 8260C | ACETONE | 67-64-1 | UG_L | 2.5 | UJ | c |
| 8260C | BENZENE | 71-43-2 | UG_L | 0.5 | U | |
| 8260C | BROMODICHLOROMETHANE | 75-27-4 | UG_L | 0.5 | U | |
| 8260C | BROMOFORM | 75-25-2 | UG_L | 0.5 | U | |
| 8260C | BROMOMETHANE | 74-83-9 | UG_L | 1 | UJ | c |
| 8260C | CARBON DISULFIDE | 75-15-0 | UG_L | 0.5 | UJ | c |
| 8260C | CARBON TETRACHLORIDE | 56-23-5 | UG_L | 1.5 | | |
| 8260C | CHLOROBENZENE | 108-90-7 | UG_L | 0.5 | U | |
| 8260C | CHLOROETHANE | 75-00-3 | UG_L | 1 | UJ | c |
| 8260C | CHLOROFORM | 67-66-3 | UG_L | 2.4 | | |
| 8260C | CHLOROMETHANE | 74-87-3 | UG_L | 1 | UJ | c |
| 8260C | CIS-1,2-DICHLOROETHENE | 156-59-2 | UG_L | 6 | | |
| 8260C | CIS-1,3-DICHLOROPROPENE | 10061-01-5 | UG_L | 0.5 | U | |
| 8260C | CYCLOHEXANE | 110-82-7 | UG_L | 0.5 | UJ | c |
| 8260C | DIBROMOCHLOROMETHANE | 124-48-1 | UG_L | 0.5 | U | |
| 8260C | DICHLOROFLUOROMETHANE | 75-71-8 | UG_L | 1 | UJ | c |
| 8260C | ETHYLBENZENE | 100-41-4 | UG_L | 0.5 | U | |
| 8260C | ISOPROPYLBENZENE | 98-82-8 | UG_L | 0.5 | U | |
| 8260C | M- AND P-XYLENE | 108-38-3/106-42 | UG_L | 1 | U | |
| 8260C | METHYL ACETATE | 79-20-9 | UG_L | 0.75 | U | |
| 8260C | METHYL CYCLOHEXANE | 108-87-2 | UG_L | 0.5 | U | |
| 8260C | METHYL TERT-BUTYL ETHER | 1634-04-4 | UG_L | 0.5 | U | |
| 8260C | METHYLENE CHLORIDE | 75-09-2 | UG_L | 2.5 | U | |
| 8260C | O-XYLENE | 95-47-6 | UG_L | 0.5 | U | |
| 8260C | STYRENE | 100-42-5 | UG_L | 0.5 | U | |
| 8260C | TETRACHLOROETHENE | 127-18-4 | UG_L | 2.7 | J | c |
| 8260C | TOLUENE | 108-88-3 | UG_L | 0.63 | J | |
| 8260C | TRANS-1,2-DICHLOROETHENE | 156-60-5 | UG_L | 0.5 | U | |
| 8260C | TRANS-1,3-DICHLOROPROPENE | 10061-02-6 | UG_L | 0.5 | U | |
| 8260C | TRICHLOROETHENE | 79-01-6 | UG_L | 4600 | J | s |
| 8260C | TRICHLOROFLUOROMETHANE | 75-69-4 | UG_L | 1 | U | |
| 8260C | VINYL CHLORIDE | 75-01-4 | UG_L | 1 | U | |
| 8260C | XYLENES, TOTAL | 1330-20-7 | UG_L | 1.5 | U | |
| 8270D_SIM | 1,4-DIOXANE | 123-91-1 | UG_L | 14 | | |

| Sample Delivery Group | | | | SI1843 | | |
|-----------------------|---------------------------------------|-----------------|-------|-------------------|------|-----|
| Lab ID | | | | SI1843-11 | | |
| Sample ID | | | | RE122D3-GW-032415 | | |
| Sample Date | | | | 3/24/2015 | | |
| Sample Type | | | | Groundwater | | |
| Method | Analyte | CAS No | Units | Result | Qual | RC |
| 8260C | 1,1,1-TRICHLOROETHANE | 71-55-6 | UG_L | 0.5 | U | |
| 8260C | 1,1,2,2-TETRACHLOROETHANE | 79-34-5 | UG_L | 0.5 | U | |
| 8260C | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 76-13-1 | UG_L | 0.5 | U | |
| 8260C | 1,1,2-TRICHLOROETHANE | 79-00-5 | UG_L | 0.5 | U | |
| 8260C | 1,1-DICHLOROETHANE | 75-34-3 | UG_L | 0.5 | U | |
| 8260C | 1,1-DICHLOROETHENE | 75-35-4 | UG_L | 0.5 | U | |
| 8260C | 1,2,4-TRICHLOROBENZENE | 120-82-1 | UG_L | 0.5 | U | |
| 8260C | 1,2-DIBROMO-3-CHLOROPROPANE | 96-12-8 | UG_L | 0.75 | U | |
| 8260C | 1,2-DIBROMOETHANE | 106-93-4 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROBENZENE | 95-50-1 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROETHANE | 107-06-2 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROETHENE, TOTAL | 540-59-0 | UG_L | 1 | U | |
| 8260C | 1,2-DICHLOROPROPANE | 78-87-5 | UG_L | 0.5 | U | |
| 8260C | 1,3-DICHLOROBENZENE | 541-73-1 | UG_L | 0.5 | U | |
| 8260C | 1,4-DICHLOROBENZENE | 106-46-7 | UG_L | 0.5 | U | |
| 8260C | 2-BUTANONE | 78-93-3 | UG_L | 2.5 | UJ | c |
| 8260C | 2-HEXANONE | 591-78-6 | UG_L | 2.5 | UJ | c |
| 8260C | 4-METHYL-2-PENTANONE | 108-10-1 | UG_L | 2.5 | U | |
| 8260C | ACETONE | 67-64-1 | UG_L | 2.5 | UJ | c |
| 8260C | BENZENE | 71-43-2 | UG_L | 0.5 | U | |
| 8260C | BROMODICHLOROMETHANE | 75-27-4 | UG_L | 0.5 | U | |
| 8260C | BROMOFORM | 75-25-2 | UG_L | 0.5 | U | |
| 8260C | BROMOMETHANE | 74-83-9 | UG_L | 1 | UJ | c |
| 8260C | CARBON DISULFIDE | 75-15-0 | UG_L | 0.5 | UJ | c |
| 8260C | CARBON TETRACHLORIDE | 56-23-5 | UG_L | 0.5 | U | |
| 8260C | CHLOROBENZENE | 108-90-7 | UG_L | 0.5 | U | |
| 8260C | CHLOROETHANE | 75-00-3 | UG_L | 1 | U | |
| 8260C | CHLOROFORM | 67-66-3 | UG_L | 0.5 | U | |
| 8260C | CHLOROMETHANE | 74-87-3 | UG_L | 1 | UJ | c |
| 8260C | CIS-1,2-DICHLOROETHENE | 156-59-2 | UG_L | 0.5 | U | |
| 8260C | CIS-1,3-DICHLOROPROPENE | 10061-01-5 | UG_L | 0.5 | U | |
| 8260C | CYCLOHEXANE | 110-82-7 | UG_L | 0.5 | UJ | c |
| 8260C | DIBROMOCHLOROMETHANE | 124-48-1 | UG_L | 0.5 | U | |
| 8260C | DICHLOROFLUOROMETHANE | 75-71-8 | UG_L | 1 | UJ | c |
| 8260C | ETHYLBENZENE | 100-41-4 | UG_L | 0.5 | U | |
| 8260C | ISOPROPYLBENZENE | 98-82-8 | UG_L | 0.5 | U | |
| 8260C | M- AND P-XYLENE | 108-38-3/106-42 | UG_L | 1 | U | |
| 8260C | METHYL ACETATE | 79-20-9 | UG_L | 0.75 | U | |
| 8260C | METHYL CYCLOHEXANE | 108-87-2 | UG_L | 0.5 | UJ | l,c |
| 8260C | METHYL TERT-BUTYL ETHER | 1634-04-4 | UG_L | 0.5 | U | |
| 8260C | METHYLENE CHLORIDE | 75-09-2 | UG_L | 2.5 | U | |
| 8260C | O-XYLENE | 95-47-6 | UG_L | 0.5 | U | |
| 8260C | STYRENE | 100-42-5 | UG_L | 0.5 | U | |
| 8260C | TETRACHLOROETHENE | 127-18-4 | UG_L | 0.5 | UJ | c |
| 8260C | TOLUENE | 108-88-3 | UG_L | 0.37 | J | |
| 8260C | TRANS-1,2-DICHLOROETHENE | 156-60-5 | UG_L | 0.5 | U | |
| 8260C | TRANS-1,3-DICHLOROPROPENE | 10061-02-6 | UG_L | 0.5 | U | |
| 8260C | TRICHLOROETHENE | 79-01-6 | UG_L | 6.8 | | |
| 8260C | TRICHLOROFLUOROMETHANE | 75-69-4 | UG_L | 1 | U | |
| 8260C | VINYL CHLORIDE | 75-01-4 | UG_L | 1 | U | |
| 8260C | XYLENES, TOTAL | 1330-20-7 | UG_L | 1.5 | U | |
| 8270D_SIM | 1,4-DIOXANE | 123-91-1 | UG_L | 0.19 | U | |

| Sample Delivery Group | | | | SI1843 | | |
|-----------------------|---------------------------------------|-----------------|-------|-------------------|------|-----|
| Lab ID | | | | SI1843-14 | | |
| Sample ID | | | | TT101D1-GW-032415 | | |
| Sample Date | | | | 3/24/2015 | | |
| Sample Type | | | | Groundwater | | |
| Method | Analyte | CAS No | Units | Result | Qual | RC |
| 8260C | 1,1,1-TRICHLOROETHANE | 71-55-6 | UG_L | 0.68 | J | |
| 8260C | 1,1,2,2-TETRACHLOROETHANE | 79-34-5 | UG_L | 0.5 | U | |
| 8260C | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 76-13-1 | UG_L | 21 | | |
| 8260C | 1,1,2-TRICHLOROETHANE | 79-00-5 | UG_L | 0.49 | J | |
| 8260C | 1,1-DICHLOROETHANE | 75-34-3 | UG_L | 0.71 | J | |
| 8260C | 1,1-DICHLOROETHENE | 75-35-4 | UG_L | 4.9 | | |
| 8260C | 1,2,4-TRICHLOROBENZENE | 120-82-1 | UG_L | 0.5 | U | |
| 8260C | 1,2-DIBROMO-3-CHLOROPROPANE | 96-12-8 | UG_L | 0.75 | U | |
| 8260C | 1,2-DIBROMOETHANE | 106-93-4 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROBENZENE | 95-50-1 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROETHANE | 107-06-2 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROETHENE, TOTAL | 540-59-0 | UG_L | 1.8 | J | |
| 8260C | 1,2-DICHLOROPROPANE | 78-87-5 | UG_L | 0.5 | U | |
| 8260C | 1,3-DICHLOROBENZENE | 541-73-1 | UG_L | 0.5 | U | |
| 8260C | 1,4-DICHLOROBENZENE | 106-46-7 | UG_L | 0.5 | U | |
| 8260C | 2-BUTANONE | 78-93-3 | UG_L | 2.5 | UJ | c |
| 8260C | 2-HEXANONE | 591-78-6 | UG_L | 2.5 | UJ | c |
| 8260C | 4-METHYL-2-PENTANONE | 108-10-1 | UG_L | 2.5 | U | |
| 8260C | ACETONE | 67-64-1 | UG_L | 2.5 | UJ | c |
| 8260C | BENZENE | 71-43-2 | UG_L | 0.5 | U | |
| 8260C | BROMODICHLOROMETHANE | 75-27-4 | UG_L | 0.5 | U | |
| 8260C | BROMOFORM | 75-25-2 | UG_L | 0.5 | U | |
| 8260C | BROMOMETHANE | 74-83-9 | UG_L | 1 | UJ | c |
| 8260C | CARBON DISULFIDE | 75-15-0 | UG_L | 0.5 | UJ | c |
| 8260C | CARBON TETRACHLORIDE | 56-23-5 | UG_L | 0.89 | J | |
| 8260C | CHLOROBENZENE | 108-90-7 | UG_L | 0.5 | U | |
| 8260C | CHLOROETHANE | 75-00-3 | UG_L | 1 | U | |
| 8260C | CHLOROFORM | 67-66-3 | UG_L | 0.92 | J | |
| 8260C | CHLOROMETHANE | 74-87-3 | UG_L | 1 | UJ | c |
| 8260C | CIS-1,2-DICHLOROETHENE | 156-59-2 | UG_L | 1.8 | | |
| 8260C | CIS-1,3-DICHLOROPROPENE | 10061-01-5 | UG_L | 0.5 | U | |
| 8260C | CYCLOHEXANE | 110-82-7 | UG_L | 0.5 | UJ | c |
| 8260C | DIBROMOCHLOROMETHANE | 124-48-1 | UG_L | 0.5 | U | |
| 8260C | DICHLOROFLUOROMETHANE | 75-71-8 | UG_L | 2.1 | J | c |
| 8260C | ETHYLBENZENE | 100-41-4 | UG_L | 0.5 | U | |
| 8260C | ISOPROPYLBENZENE | 98-82-8 | UG_L | 0.5 | U | |
| 8260C | M- AND P-XYLENE | 108-38-3/106-42 | UG_L | 1 | U | |
| 8260C | METHYL ACETATE | 79-20-9 | UG_L | 0.75 | U | |
| 8260C | METHYL CYCLOHEXANE | 108-87-2 | UG_L | 0.5 | UJ | l,c |
| 8260C | METHYL TERT-BUTYL ETHER | 1634-04-4 | UG_L | 0.5 | U | |
| 8260C | METHYLENE CHLORIDE | 75-09-2 | UG_L | 2.5 | U | |
| 8260C | O-XYLENE | 95-47-6 | UG_L | 0.5 | U | |
| 8260C | STYRENE | 100-42-5 | UG_L | 0.5 | U | |
| 8260C | TETRACHLOROETHENE | 127-18-4 | UG_L | 0.5 | UJ | c |
| 8260C | TOLUENE | 108-88-3 | UG_L | 0.5 | U | |
| 8260C | TRANS-1,2-DICHLOROETHENE | 156-60-5 | UG_L | 0.5 | U | |
| 8260C | TRANS-1,3-DICHLOROPROPENE | 10061-02-6 | UG_L | 0.5 | U | |
| 8260C | TRICHLOROETHENE | 79-01-6 | UG_L | 170 | | |
| 8260C | TRICHLOROFLUOROMETHANE | 75-69-4 | UG_L | 1 | U | |
| 8260C | VINYL CHLORIDE | 75-01-4 | UG_L | 1 | U | |
| 8260C | XYLENES, TOTAL | 1330-20-7 | UG_L | 1.5 | U | |
| 8270D_SIM | 1,4-DIOXANE | 123-91-1 | UG_L | 8.7 | | |

| Sample Delivery Group | | | | SI1843 | | |
|-----------------------|---------------------------------------|-----------------|-------|-------------------|------|----|
| Lab ID | | | | SI1843-12 | | |
| Sample ID | | | | TT101D2-GW-032415 | | |
| Sample Date | | | | 3/24/2015 | | |
| Sample Type | | | | Groundwater | | |
| Method | Analyte | CAS No | Units | Result | Qual | RC |
| 8260C | 1,1,1-TRICHLOROETHANE | 71-55-6 | UG_L | 0.44 | J | |
| 8260C | 1,1,2,2-TETRACHLOROETHANE | 79-34-5 | UG_L | 0.5 | U | |
| 8260C | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 76-13-1 | UG_L | 25 | | |
| 8260C | 1,1,2-TRICHLOROETHANE | 79-00-5 | UG_L | 0.62 | J | |
| 8260C | 1,1-DICHLOROETHANE | 75-34-3 | UG_L | 0.75 | J | |
| 8260C | 1,1-DICHLOROETHENE | 75-35-4 | UG_L | 4 | | |
| 8260C | 1,2,4-TRICHLOROBENZENE | 120-82-1 | UG_L | 0.5 | U | |
| 8260C | 1,2-DIBROMO-3-CHLOROPROPANE | 96-12-8 | UG_L | 0.75 | U | |
| 8260C | 1,2-DIBROMOETHANE | 106-93-4 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROBENZENE | 95-50-1 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROETHANE | 107-06-2 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROETHENE, TOTAL | 540-59-0 | UG_L | 2 | | |
| 8260C | 1,2-DICHLOROPROPANE | 78-87-5 | UG_L | 0.5 | U | |
| 8260C | 1,3-DICHLOROBENZENE | 541-73-1 | UG_L | 0.5 | U | |
| 8260C | 1,4-DICHLOROBENZENE | 106-46-7 | UG_L | 0.5 | U | |
| 8260C | 2-BUTANONE | 78-93-3 | UG_L | 2.5 | UJ | c |
| 8260C | 2-HEXANONE | 591-78-6 | UG_L | 2.5 | UJ | c |
| 8260C | 4-METHYL-2-PENTANONE | 108-10-1 | UG_L | 2.5 | U | |
| 8260C | ACETONE | 67-64-1 | UG_L | 2.5 | UJ | c |
| 8260C | BENZENE | 71-43-2 | UG_L | 0.5 | U | |
| 8260C | BROMODICHLOROMETHANE | 75-27-4 | UG_L | 0.5 | U | |
| 8260C | BROMOFORM | 75-25-2 | UG_L | 0.5 | U | |
| 8260C | BROMOMETHANE | 74-83-9 | UG_L | 1 | UJ | c |
| 8260C | CARBON DISULFIDE | 75-15-0 | UG_L | 0.5 | UJ | c |
| 8260C | CARBON TETRACHLORIDE | 56-23-5 | UG_L | 0.36 | J | |
| 8260C | CHLOROBENZENE | 108-90-7 | UG_L | 0.5 | U | |
| 8260C | CHLOROETHANE | 75-00-3 | UG_L | 1 | UJ | c |
| 8260C | CHLOROFORM | 67-66-3 | UG_L | 0.82 | J | |
| 8260C | CHLOROMETHANE | 74-87-3 | UG_L | 1 | UJ | c |
| 8260C | CIS-1,2-DICHLOROETHENE | 156-59-2 | UG_L | 2 | | |
| 8260C | CIS-1,3-DICHLOROPROPENE | 10061-01-5 | UG_L | 0.5 | U | |
| 8260C | CYCLOHEXANE | 110-82-7 | UG_L | 0.5 | UJ | c |
| 8260C | DIBROMOCHLOROMETHANE | 124-48-1 | UG_L | 0.5 | U | |
| 8260C | DICHLORODIFLUOROMETHANE | 75-71-8 | UG_L | 1 | UJ | c |
| 8260C | ETHYLBENZENE | 100-41-4 | UG_L | 0.5 | U | |
| 8260C | ISOPROPYLBENZENE | 98-82-8 | UG_L | 0.5 | U | |
| 8260C | M- AND P-XYLENE | 108-38-3/106-42 | UG_L | 1 | U | |
| 8260C | METHYL ACETATE | 79-20-9 | UG_L | 0.75 | U | |
| 8260C | METHYL CYCLOHEXANE | 108-87-2 | UG_L | 0.5 | U | |
| 8260C | METHYL TERT-BUTYL ETHER | 1634-04-4 | UG_L | 0.5 | U | |
| 8260C | METHYLENE CHLORIDE | 75-09-2 | UG_L | 2.5 | U | |
| 8260C | O-XYLENE | 95-47-6 | UG_L | 0.5 | U | |
| 8260C | STYRENE | 100-42-5 | UG_L | 0.5 | U | |
| 8260C | TETRACHLOROETHENE | 127-18-4 | UG_L | 0.47 | J | c |
| 8260C | TOLUENE | 108-88-3 | UG_L | 0.5 | U | |
| 8260C | TRANS-1,2-DICHLOROETHENE | 156-60-5 | UG_L | 0.5 | U | |
| 8260C | TRANS-1,3-DICHLOROPROPENE | 10061-02-6 | UG_L | 0.5 | U | |
| 8260C | TRICHLOROETHENE | 79-01-6 | UG_L | 480 | J | s |
| 8260C | TRICHLOROFLUOROMETHANE | 75-69-4 | UG_L | 1 | U | |
| 8260C | VINYL CHLORIDE | 75-01-4 | UG_L | 1 | U | |
| 8260C | XYLENES, TOTAL | 1330-20-7 | UG_L | 1.5 | U | |
| 8270D_SIM | 1,4-DIOXANE | 123-91-1 | UG_L | 2.4 | | |

| Sample Delivery Group | | | | S11843 | | |
|-----------------------|---------------------------------------|-----------------|-------|------------------|------|-----|
| Lab ID | | | | S11843-13 | | |
| Sample ID | | | | TT101D-GW-032415 | | |
| Sample Date | | | | 3/24/2015 | | |
| Sample Type | | | | Groundwater | | |
| Method | Analyte | CAS No | Units | Result | Qual | RC |
| 8260C | 1,1,1-TRICHLOROETHANE | 71-55-6 | UG_L | 0.36 | J | |
| 8260C | 1,1,2,2-TETRACHLOROETHANE | 79-34-5 | UG_L | 0.5 | U | |
| 8260C | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 76-13-1 | UG_L | 22 | | |
| 8260C | 1,1,2-TRICHLOROETHANE | 79-00-5 | UG_L | 0.5 | U | |
| 8260C | 1,1-DICHLOROETHANE | 75-34-3 | UG_L | 0.78 | J | |
| 8260C | 1,1-DICHLOROETHENE | 75-35-4 | UG_L | 3.5 | | |
| 8260C | 1,2,4-TRICHLOROBENZENE | 120-82-1 | UG_L | 0.5 | U | |
| 8260C | 1,2-DIBROMO-3-CHLOROPROPANE | 96-12-8 | UG_L | 0.75 | U | |
| 8260C | 1,2-DIBROMOETHANE | 106-93-4 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROBENZENE | 95-50-1 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROETHANE | 107-06-2 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROETHENE, TOTAL | 540-59-0 | UG_L | 3 | | |
| 8260C | 1,2-DICHLOROPROPANE | 78-87-5 | UG_L | 0.5 | U | |
| 8260C | 1,3-DICHLOROBENZENE | 541-73-1 | UG_L | 0.5 | U | |
| 8260C | 1,4-DICHLOROBENZENE | 106-46-7 | UG_L | 0.5 | U | |
| 8260C | 2-BUTANONE | 78-93-3 | UG_L | 2.5 | UJ | c |
| 8260C | 2-HEXANONE | 591-78-6 | UG_L | 2.5 | UJ | c |
| 8260C | 4-METHYL-2-PENTANONE | 108-10-1 | UG_L | 2.5 | U | |
| 8260C | ACETONE | 67-64-1 | UG_L | 2.5 | UJ | c |
| 8260C | BENZENE | 71-43-2 | UG_L | 0.5 | U | |
| 8260C | BROMODICHLOROMETHANE | 75-27-4 | UG_L | 0.5 | U | |
| 8260C | BROMOFORM | 75-25-2 | UG_L | 0.5 | U | |
| 8260C | BROMOMETHANE | 74-83-9 | UG_L | 1 | UJ | c |
| 8260C | CARBON DISULFIDE | 75-15-0 | UG_L | 0.5 | UJ | c |
| 8260C | CARBON TETRACHLORIDE | 56-23-5 | UG_L | 0.5 | U | |
| 8260C | CHLOROBENZENE | 108-90-7 | UG_L | 0.5 | U | |
| 8260C | CHLOROETHANE | 75-00-3 | UG_L | 1 | U | |
| 8260C | CHLOROFORM | 67-66-3 | UG_L | 0.5 | J | |
| 8260C | CHLOROMETHANE | 74-87-3 | UG_L | 1 | UJ | c |
| 8260C | CIS-1,2-DICHLOROETHENE | 156-59-2 | UG_L | 3 | | |
| 8260C | CIS-1,3-DICHLOROPROPENE | 10061-01-5 | UG_L | 0.5 | U | |
| 8260C | CYCLOHEXANE | 110-82-7 | UG_L | 0.5 | UJ | c |
| 8260C | DIBROMOCHLOROMETHANE | 124-48-1 | UG_L | 0.5 | U | |
| 8260C | DICHLORODIFLUOROMETHANE | 75-71-8 | UG_L | 2.1 | J | c |
| 8260C | ETHYLBENZENE | 100-41-4 | UG_L | 0.5 | U | |
| 8260C | ISOPROPYLBENZENE | 98-82-8 | UG_L | 0.5 | U | |
| 8260C | M- AND P-XYLENE | 108-38-3/106-42 | UG_L | 1 | U | |
| 8260C | METHYL ACETATE | 79-20-9 | UG_L | 0.75 | U | |
| 8260C | METHYL CYCLOHEXANE | 108-87-2 | UG_L | 0.5 | UJ | l,c |
| 8260C | METHYL TERT-BUTYL ETHER | 1634-04-4 | UG_L | 0.5 | U | |
| 8260C | METHYLENE CHLORIDE | 75-09-2 | UG_L | 2.5 | U | |
| 8260C | O-XYLENE | 95-47-6 | UG_L | 0.5 | U | |
| 8260C | STYRENE | 100-42-5 | UG_L | 0.5 | U | |
| 8260C | TETRACHLOROETHENE | 127-18-4 | UG_L | 0.5 | UJ | c |
| 8260C | TOLUENE | 108-88-3 | UG_L | 0.5 | U | |
| 8260C | TRANS-1,2-DICHLOROETHENE | 156-60-5 | UG_L | 0.5 | U | |
| 8260C | TRANS-1,3-DICHLOROPROPENE | 10061-02-6 | UG_L | 0.5 | U | |
| 8260C | TRICHLOROETHENE | 79-01-6 | UG_L | 61 | | |
| 8260C | TRICHLOROFLUOROMETHANE | 75-69-4 | UG_L | 1 | U | |
| 8260C | VINYL CHLORIDE | 75-01-4 | UG_L | 1 | U | |
| 8260C | XYLENES, TOTAL | 1330-20-7 | UG_L | 1.5 | U | |
| 8270D_SIM | 1,4-DIOXANE | 123-91-1 | UG_L | 9.7 | | |

| Sample Delivery Group | | | | S11843 | | |
|-----------------------|---------------------------------------|-----------------|-------|-----------------|------|-----|
| Lab ID | | | | S11843-15 | | |
| Sample ID | | | | TRIPBLANK031615 | | |
| Sample Date | | | | 3/24/2015 | | |
| Sample Type | | | | Trip Blank | | |
| Method | Analyte | CAS No | Units | Result | Qual | RC |
| 8260C | 1,1,1-TRICHLOROETHANE | 71-55-6 | UG_L | 0.5 | U | |
| 8260C | 1,1,2,2-TETRACHLOROETHANE | 79-34-5 | UG_L | 0.5 | U | |
| 8260C | 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE | 76-13-1 | UG_L | 0.5 | U | |
| 8260C | 1,1,2-TRICHLOROETHANE | 79-00-5 | UG_L | 0.5 | U | |
| 8260C | 1,1-DICHLOROETHANE | 75-34-3 | UG_L | 0.5 | U | |
| 8260C | 1,1-DICHLOROETHENE | 75-35-4 | UG_L | 0.5 | U | |
| 8260C | 1,2,4-TRICHLOROBENZENE | 120-82-1 | UG_L | 0.5 | U | |
| 8260C | 1,2-DIBROMO-3-CHLOROPROPANE | 96-12-8 | UG_L | 0.75 | U | |
| 8260C | 1,2-DIBROMOETHANE | 106-93-4 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROBENZENE | 95-50-1 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROETHANE | 107-06-2 | UG_L | 0.5 | U | |
| 8260C | 1,2-DICHLOROETHENE, TOTAL | 540-59-0 | UG_L | 1 | U | |
| 8260C | 1,2-DICHLOROPROPANE | 78-87-5 | UG_L | 0.5 | U | |
| 8260C | 1,3-DICHLOROBENZENE | 541-73-1 | UG_L | 0.5 | U | |
| 8260C | 1,4-DICHLOROBENZENE | 106-46-7 | UG_L | 0.5 | U | |
| 8260C | 2-BUTANONE | 78-93-3 | UG_L | 2.5 | UJ | c |
| 8260C | 2-HEXANONE | 591-78-6 | UG_L | 2.5 | UJ | c |
| 8260C | 4-METHYL-2-PENTANONE | 108-10-1 | UG_L | 2.5 | U | |
| 8260C | ACETONE | 67-64-1 | UG_L | 2.5 | UJ | c |
| 8260C | BENZENE | 71-43-2 | UG_L | 0.5 | U | |
| 8260C | BROMODICHLOROMETHANE | 75-27-4 | UG_L | 0.5 | U | |
| 8260C | BROMOFORM | 75-25-2 | UG_L | 0.5 | U | |
| 8260C | BROMOMETHANE | 74-83-9 | UG_L | 1 | UJ | c |
| 8260C | CARBON DISULFIDE | 75-15-0 | UG_L | 0.5 | UJ | c |
| 8260C | CARBON TETRACHLORIDE | 56-23-5 | UG_L | 0.5 | U | |
| 8260C | CHLOROBENZENE | 108-90-7 | UG_L | 0.5 | U | |
| 8260C | CHLOROETHANE | 75-00-3 | UG_L | 1 | U | |
| 8260C | CHLOROFORM | 67-66-3 | UG_L | 0.5 | U | |
| 8260C | CHLOROMETHANE | 74-87-3 | UG_L | 1 | UJ | c |
| 8260C | CIS-1,2-DICHLOROETHENE | 156-59-2 | UG_L | 0.5 | U | |
| 8260C | CIS-1,3-DICHLOROPROPENE | 10061-01-5 | UG_L | 0.5 | U | |
| 8260C | CYCLOHEXANE | 110-82-7 | UG_L | 0.5 | UJ | c |
| 8260C | DIBROMOCHLOROMETHANE | 124-48-1 | UG_L | 0.5 | U | |
| 8260C | DICHLORODIFLUOROMETHANE | 75-71-8 | UG_L | 1 | UJ | c |
| 8260C | ETHYLBENZENE | 100-41-4 | UG_L | 0.5 | U | |
| 8260C | ISOPROPYLBENZENE | 98-82-8 | UG_L | 0.5 | U | |
| 8260C | M- AND P-XYLENE | 108-38-3/106-42 | UG_L | 1 | U | |
| 8260C | METHYL ACETATE | 79-20-9 | UG_L | 0.75 | U | |
| 8260C | METHYL CYCLOHEXANE | 108-87-2 | UG_L | 0.5 | UJ | l,c |
| 8260C | METHYL TERT-BUTYL ETHER | 1634-04-4 | UG_L | 0.5 | U | |
| 8260C | METHYLENE CHLORIDE | 75-09-2 | UG_L | 2.5 | U | |
| 8260C | O-XYLENE | 95-47-6 | UG_L | 0.5 | U | |
| 8260C | STYRENE | 100-42-5 | UG_L | 0.5 | U | |
| 8260C | TETRACHLOROETHENE | 127-18-4 | UG_L | 0.5 | UJ | c |
| 8260C | TOLUENE | 108-88-3 | UG_L | 0.5 | U | |
| 8260C | TRANS-1,2-DICHLOROETHENE | 156-60-5 | UG_L | 0.5 | U | |
| 8260C | TRANS-1,3-DICHLOROPROPENE | 10061-02-6 | UG_L | 0.5 | U | |
| 8260C | TRICHLOROETHENE | 79-01-6 | UG_L | 0.5 | U | |
| 8260C | TRICHLOROFLUOROMETHANE | 75-69-4 | UG_L | 1 | U | |
| 8260C | VINYL CHLORIDE | 75-01-4 | UG_L | 1 | U | |
| 8260C | XYLENES, TOTAL | 1330-20-7 | UG_L | 1.5 | U | |
| 8270D_SIM | 1,4-DIOXANE | 123-91-1 | UG_L | | | |

Notes:

UG_L = Micrograms per liter
Qual = Final qualifier (Refer to Attachment B)
RC = Reason code (Refer to Attachment C)



DATA VALIDATION REPORT

| | | |
|-------------------------|--|--|
| Project: | Regional Groundwater Investigation — NWIRP Bethpage | |
| Laboratory: | Katahdin Analytical | |
| Sample Delivery Groups: | SI0198 | |
| Analyses/Method: | Total Organic Carbon (TOC) by U.S. EPA SW-846 Method 9060A and Standard Method 5310B for Total Organic Carbon by High-Temperature Combustion | |
| Validation Level: | 3 | |
| Project Number: | 0888812477.SA.DV | |
| Prepared by: | Dana Miller/Resolution Consultants | Completed on: 02/26/2015 Revised on: 10/27/2015 |
| Reviewed by: | Tina Cantwell/Resolution Consultants | File Name: SI0198_9060A_5310B |

SUMMARY

This report summarizes data review findings for samples listed below, collected by Resolution Consultants from the Regional Groundwater Investigation — NWIRP Bethpage site on 9 January 2015 in accordance with the following Sampling and Analysis Plans:

- *Sampling and Analysis Plan, Bethpage, New York.* (Resolution Consultants April 2013).
- *UFP SAP Addendum, Installation of Vertical Profile Borings and Monitoring Wells, Operable Unit 2, NWIRP Bethpage, New York.* (Resolution Consultants November 2013).
- *UFP SAP Addendum, Inclusion of Additional Target Analytes for Volatile Organics Analyses, NWIRP Bethpage OU2, Bethpage, New York.* (Resolution Consultants August 2014).

| Sample ID | Lab ID | Matrix/Sample Type | Analysis |
|-----------------------------|----------|--------------------|----------|
| RE122D2-SOIL-010915-593-595 | SI0198-1 | Soil | 9060A |
| RE122D2-EB-010915 | SI0198-2 | Equipment Blank | 5310B |

Data validation activities were conducted using the following guidance documents: *Test Methods for Evaluating Solid Waste, Physical/Chemical Methods SW-846, specifically Method 9060A, Total Organic Carbon* (U.S. EPA, 1996), *Method SM5310B, Total Organic Carbon by High-Temperature Combustion*, *U.S. Environmental Protection Agency (U.S. EPA) Contract Laboratory Program National Functional Guidelines for Inorganic Superfund Data Review* (NFG, January 2010, and Department of Defense (DoD) Quality Systems Manual (QSM) for Environmental Laboratories, Version 4.2 (October 2010). In

the absence of method-specific information, laboratory quality control (QC) limits, project-specific requirements and/or professional judgment were used as appropriate.

REVIEW ELEMENTS

The data were evaluated based on the following parameters (where applicable to the method):

- ✓ Data completeness (chain-of-custody)/sample integrity
- ✓ Holding times and sample preservation
- NA Gas chromatography/Mass spectrometer performance checks
- ✓ Initial calibration/continuing calibration verification
- ✓ Laboratory blanks/equipment blanks/field blanks/trip blanks
- NA Surrogate spike recoveries
- ✓ Matrix spike and/or matrix spike duplicate results
- ✓ Laboratory control sample laboratory control sample duplicate results
- NA Field duplicates
- NA Internal standards
- ✓ Sample results/reporting issues

The symbol (✓) indicates that no validation qualifiers were applied based on this parameter. NA indicates that the parameter was not included as part of this data set or was not applicable to this validation and therefore not reviewed. Acceptable data parameters for which all criteria were met and no qualification was performed and non-conformance or other issues that were noted during validation, but did not result in qualification of data are not discussed further.

Qualifications Actions

The data was reviewed independently from the laboratory to assess data quality. All compounds detected at concentrations less than the limit of quantitation but greater than the method detection limit were qualified by the laboratory as estimated (J). This "J" qualifier was retained during data validation. Any sample that was analyzed at a dilution because of high concentrations of target or non-target analytes was checked to confirm that the results and/or sample-specific limit of quantitation and limit of detections were adjusted accordingly by the laboratory.

No results were qualified during this review. Analytical completeness was calculated to be 100% and the data are usable for their intended purpose, according to U.S. Environmental Protection Agency and Department of Defense guidelines. Katahdin Analytical discovered a QC error in Total Organic Carbon calculations, informed us, and has implemented corrective action provided in Attachment A. Attachment B provides final results after data review.



ATTACHMENTS

Attachment A: Katahdin Analytical Corrective Action Report

Attachment B: Final Results after Data Review

Attachment A
Katahdin Analytical Corrective Action Report

KATAHDIN ANALYTICAL SERVICES, INC. – CORRECTIVE ACTION REPORT

| | | |
|--|------------------------------------|------------------|
| Problem Identification (Person initiating CAR) Name: Leslie Dimond | | Date: 08/28/2015 |
| Discovered by Laboratory | X Discovered by Client (Complaint) | Other |
| Details of Problem: | | |
| <p>On August 28, 2015, a client called regarding several TOC results. They thought the results for their samples seemed to be off by a factor of ten. George Brewer looked into the matter and found that their sample results were calculated incorrectly. Calculations are set up to be performed automatically in KIMS. Each product code for a test contains information such as the method name, LOQ, LOD, MDL, limits and a specific process chain for reporting to the either the LOQ, LOD or MDL (i.e. three process chains). This process chain contains all of the calculations associated with a test including separate calculations for sample results, adjusted LOQs, LODs and MDLs, and results for QC samples. In October of 2014, a client asked why the lab was not adjusting LOQs, LODs and/or MDLs for the sample amount used, but only for an instrument dilution and total solids. The lab had felt that there wasn't really a standard weight amount, so this was not incorporated. After numerous discussions, the lab decided to change this and use 500 mg as the standard and to incorporate any different amount into the LOQ/LOD/MDL adjustment. MIS was asked to revise this TOC calculation in our Laboratory Information System (KIMS). The incorrect information was given to MIS, so the adjusted LOQ/LOD/MDL calculation was wrong.</p> <p>The formula was: $ADJ. LOQ/LOD/MDL = LOQ/LOD/MDL * DF * (Sample\ amount / 1000) * (100/TS)$.</p> <p>During data review, sample results that are calculated automatically by KIMS, are checked at a frequency of 10% of all calculations. Sample results are checked in this manner, but adjusted limits are not necessarily checked. In December of 2014, a data reviewer noticed that some adjusted limits were not correct. This issue was reviewed with MIS, and it was discovered that the calculation was incorrect. It was corrected at this time and changed to:</p> <p>$ADJ. LOQ/LOD/MDL = LOQ/LOD/MDL * DF * (500 / Sample\ Amount) * (100/TS)$ (where 500 is the standard sample amount).</p> <p>At the time this change was made, MIS was under the impression that the sample calculation was also incorrect and also needed to be corrected.</p> <p>The formula for the sample result was: $Results\ (ug/g) = \frac{Total\ carbon}{(sample\ amount / 1000)} * (100/TS)$ (where 1000 is a conversion factor)</p> <p>This was incorrectly changed to: $Results\ (ug/g) = \frac{Total\ carbon}{(500 / Sample\ Amount)} * TS$</p> <p>From December 29, 2014 to April 13, 2015, this error was often discovered during data review and manually corrected. The problem was not addressed with the MIS department. However, there were several cases where this error was not caught during data review and results were reported incorrectly. On April 13, 2015, this issue was brought to the attention of MIS and the sample calculation was corrected back to:</p> <p>$Results\ (ug/g) = \frac{Total\ carbon}{(sample\ amount / 1000)} * (100/TS)$ (where 1000 is a conversion factor)</p> <p>Blanks and LCSs were not affected by these changes because they are calculated using different calculations since total solids is not used in these situations. Duplicates and MS/MSD samples are also not affected because they use different calculations since these calculation strings involve recoveries, RPDs, etc.</p> | | |
| Associated Non-Conformances: List logbook and page numbers | | |
| There are no non-conformances associated with this corrective action. | | |
| Root Cause Investigation & Determination (To be completed by Department Manager, Operations Manager and/or QA Officer) | | |
| Review the 6 "M's" below and investigate to determine whether one of them, or more than one, could be the cause of the problem. | | |
| Possible Causes | Details | |

KATAHDIN ANALYTICAL SERVICES, INC. – CORRECTIVE ACTION REPORT

| | |
|--|--|
| Machine (Instrument) | KIMS – incorrect formulas were entered into the process chains for TOC in soil. |
| Method (or Process) | <p>Katahdin's policy for data review: From the QAM: "For data that are reduced via computer, calculations are checked by the analyst (or designee) assigned to this task at a frequency designed to assure that the final data generation is valid." From SOP SD-904, Data Reduction, Review and Reporting: "All manual integrations, calculations and transcriptions are checked and 10% of all spreadsheet calculations are checked. The remainder of spreadsheet calculations is spot checked for potential anomalies."</p> <p>Katahdin does not have a formal process for handling calculation changes to process changes in KIMS. A verbal request is made to MIS and the change is made. Although the KIMS system does have an audit trail to track dates of changes and calculation changes, the request for these is not documented.</p> |
| Materials | Not Applicable |
| Maintenance (Is something not working correctly?) | Not Applicable |
| Man (training, human error) | All personnel involved with data review are aware of the requirement to hand check 10% of the sample results. In some cases this appears to have been done and when the results did not calculate correctly, they were corrected in KIMS. There was no communication that the error may affect other batches of samples. In other sample batches, the 10% hand check must not have occurred. |
| Mother Nature (accidents, power issues, beyond our control) | Not Applicable |

Corrective Action Plan: Name: Leslie Dimond

Date: 08/28/2015

Details of Corrective Action Plan –

Queries were run through KIMS to generate lists for TOC in soil data entered into KIMS between two timeframes:

10/23/14 to 12/29/14 (incorrect LOQ/LOD/MDL adjustment) – 14 Work Orders were found in this timeframe

12/29/14 and 4/13/15 (incorrect sample calculation) - 16 Work Orders were found in this timeframe

The lab has reviewed all TOC in soil data from these time periods for accuracy. Some inaccurate data was found (as expected from the incorrect formulas). Some data was found to be correct. In these cases, MIS was able to determine, through the KIMS audit trail, that the data had been manually corrected.

A new form has been created for personnel to fill out when requesting calculation changes in KIMS. This form will ensure that MIS is clear on exactly what needs to be changed in KIMS. Through KIMS we are currently able to track formula changes (ie. the formula before and after the change) and when the change occurred. This new form will allow Katahdin to track the changes made by person requesting the change and why the change was necessary. Supervisor and QA/Management approval are required on these forms.

There also will be a mandatory retraining for all employees on Katahdin's policy for data review, stressing that 10% of all results generated from KIMS and spreadsheets need to be confirmed. Also, this retraining will stress the need for communication. If a problem is discovered with one batch of samples, then it might also be affecting other batches.

Additionally, when calculation changes are made, Katahdin must establish a time frame of 30 days or 10 workorders where

KATAHDIN ANALYTICAL SERVICES, INC. – CORRECTIVE ACTION REPORT

senior management or the Quality Assurance Officer must also check the new calculations.

Review & Approval of Corrective Action Plan

Supervisor Approval: *[Signature]* Date: 09.11.15

Operations Manager Approval: *Deborah J. Hadeau* Date: 9.11.15

Quality Assurance Officer: *Lisee Dimond* Date: 09.11.15

Monitoring of Corrective Action (To be completed by QA Officer and/or Operations Manager): List details of follow-up

[Empty space for follow-up details]

| Corrective Action Effective | Return to Control – | Yes | No | Further Monitoring Needed/Additional Corrective Action |
|--|---------------------|-----|----|--|
| <i>[Empty space for monitoring data]</i> | | | | |

QA Approval: _____ Date: _____

Additional Information:

[Empty space for additional information]

Attachment B
Final Results after Data Review

| Sample Delivery Group | | | | SI0198 | SI0198 |
|-----------------------|----------------------|--------|-------|-----------------------------|-------------------|
| Lab ID | | | | SI0198-1 | SI0198-2 |
| Sample ID | | | | RE122D2-SOIL-010915-593-595 | RE122D2-EB-010915 |
| Sample Date | | | | 1/9/2015 | 1/9/2015 |
| Sample Type | | | | Soil | Equipment Blank |
| Method | Analyte | CAS No | Units | Result | Result |
| 5310B | TOTAL ORGANIC CARBON | -28 | MG_L | NA | 0.27 J |
| 9060A | TOTAL ORGANIC CARBON | -28 | UG_G | 630 | NA |

Notes:

- ID = Identification
- MG_L = Milligrams per liter
- UG_G = Micrograms per gram
- NA = Not analyzed
- J = The analyte was positively identified. The numerical value is the estimated concentration of the analyte in the sample.



DATA VALIDATION REPORT

| | | |
|-------------------------|--|--|
| Project: | Regional Groundwater Investigation — NWIRP Bethpage | |
| Laboratory: | Katahdin Analytical | |
| Sample Delivery Groups: | SI0594 | |
| Analyses/Method: | Total Organic Carbon (TOC) by U.S. EPA SW-846 Method 9060A and Standard Method 5310B for Total Organic Carbon by High-Temperature Combustion | |
| Validation Level: | 3 | |
| Project Number: | 0888812477.SA.DV | |
| Prepared by: | Dana Miller/Resolution Consultants | Completed on: 02/26/2015 Revised on: 10/27/2015 |
| Reviewed by: | Tina Cantwell/Resolution Consultants | File Name: SI0594_9060A_5310B |

SUMMARY

This report summarizes data review findings for samples listed below, collected by Resolution Consultants from the Regional Groundwater Investigation — NWIRP Bethpage site on 23 January 2015 in accordance with the following Sampling and Analysis Plans:

- *Sampling and Analysis Plan, Bethpage, New York.* (Resolution Consultants April 2013).
- *UFP SAP Addendum, Installation of Vertical Profile Borings and Monitoring Wells, Operable Unit 2, NWIRP Bethpage, New York.* (Resolution Consultants November 2013).
- *UFP SAP Addendum, Inclusion of Additional Target Analytes for Volatile Organics Analyses, NWIRP Bethpage OU2, Bethpage, New York.* (Resolution Consultants August 2014).

| Sample ID | Lab ID | Matrix/Sample Type | Analysis |
|-----------------------------|----------|--------------------|----------|
| RE122D1-SOIL-012315-528-530 | SI0594-1 | Soil | 9060A |
| RE122D1-EB-012315 | SI0594-2 | Equipment Blank | 5310B |

Data validation activities were conducted using the following guidance documents: *Test Methods for Evaluating Solid Waste, Physical/Chemical Methods SW-846, specifically Method 9060A, Total Organic Carbon* (U.S. EPA, 1996), *Method SM5310B, Total Organic Carbon by High-Temperature Combustion*, *U.S. Environmental Protection Agency (U.S. EPA) Contract Laboratory Program National Functional Guidelines for Inorganic Superfund Data Review* (NFG, January 2010, and Department of Defense (DoD) Quality Systems Manual (QSM) for Environmental Laboratories, Version 4.2 (October 2010). In

the absence of method-specific information, laboratory quality control (QC) limits, project-specific requirements and/or professional judgment were used as appropriate.

REVIEW ELEMENTS

The data were evaluated based on the following parameters (where applicable to the method):

- ✓ Data completeness (chain-of-custody)/sample integrity
- ✓ Holding times and sample preservation
- NA Gas chromatography/Mass spectrometer performance checks
- ✓ Initial calibration/continuing calibration verification
- ✓ Laboratory blanks/equipment blanks/field blanks/trip blanks
- NA Surrogate spike recoveries
- ✓ Matrix spike and/or matrix spike duplicate results
- ✓ Laboratory control sample laboratory control sample duplicate results
- NA Field duplicates
- NA Internal standards
- ✓ Sample results/reporting issues

The symbol (✓) indicates that no validation qualifiers were applied based on this parameter. NA indicates that the parameter was not included as part of this data set or was not applicable to this validation and therefore not reviewed. Acceptable data parameters for which all criteria were met and no qualification was performed and non-conformance or other issues that were noted during validation, but did not result in qualification of data are not discussed further.

Qualifications Actions

The data was reviewed independently from the laboratory to assess data quality. All compounds detected at concentrations less than the limit of quantitation but greater than the method detection limit were qualified by the laboratory as estimated (J). This "J" qualifier was retained during data validation. Any sample that was analyzed at a dilution because of high concentrations of target or non-target analytes was checked to confirm that the results and/or sample-specific limit of quantitation and limit of detections were adjusted accordingly by the laboratory.

No results were qualified during this review. Analytical completeness was calculated to be 100% and the data are usable for their intended purpose, according to U.S. Environmental Protection Agency and Department of Defense guidelines. Katahdin Analytical discovered a QC error in Total Organic Carbon calculations, informed us, and has implemented corrective action provided in Attachment A. Attachment B provides final results after data review.



ATTACHMENTS

Attachment A: Katahdin Analytical Corrective Action Report

Attachment B: Final Results after Data Review

Attachment A
Katahdin Analytical Corrective Action Report

KATAHDIN ANALYTICAL SERVICES, INC. – CORRECTIVE ACTION REPORT

| | | |
|--|------------------------------------|------------------|
| Problem Identification (Person initiating CAR) Name: Leslie Dimond | | Date: 08/28/2015 |
| Discovered by Laboratory | X Discovered by Client (Complaint) | Other |
| Details of Problem: | | |
| <p>On August 28, 2015, a client called regarding several TOC results. They thought the results for their samples seemed to be off by a factor of ten. George Brewer looked into the matter and found that their sample results were calculated incorrectly. Calculations are set up to be performed automatically in KIMS. Each product code for a test contains information such as the method name, LOQ, LOD, MDL, limits and a specific process chain for reporting to the either the LOQ, LOD or MDL (i.e. three process chains). This process chain contains all of the calculations associated with a test including separate calculations for sample results, adjusted LOQs, LODs and MDLs, and results for QC samples. In October of 2014, a client asked why the lab was not adjusting LOQs, LODs and/or MDLs for the sample amount used, but only for an instrument dilution and total solids. The lab had felt that there wasn't really a standard weight amount, so this was not incorporated. After numerous discussions, the lab decided to change this and use 500 mg as the standard and to incorporate any different amount into the LOQ/LOD/MDL adjustment. MIS was asked to revise this TOC calculation in our Laboratory Information System (KIMS). The incorrect information was given to MIS, so the adjusted LOQ/LOD/MDL calculation was wrong.</p> <p>The formula was: $ADJ. LOQ/LOD/MDL = LOQ/LOD/MDL * DF * (Sample\ amount / 1000) * (100/TS)$.</p> <p>During data review, sample results that are calculated automatically by KIMS, are checked at a frequency of 10% of all calculations. Sample results are checked in this manner, but adjusted limits are not necessarily checked. In December of 2014, a data reviewer noticed that some adjusted limits were not correct. This issue was reviewed with MIS, and it was discovered that the calculation was incorrect. It was corrected at this time and changed to:</p> <p>$ADJ. LOQ/LOD/MDL = LOQ/LOD/MDL * DF * (500 / Sample\ Amount) * (100/TS)$ (where 500 is the standard sample amount).</p> <p>At the time this change was made, MIS was under the impression that the sample calculation was also incorrect and also needed to be corrected.</p> <p>The formula for the sample result was: $Results\ (ug/g) = \frac{Total\ carbon}{(sample\ amount / 1000)} * (100/TS)$ (where 1000 is a conversion factor)</p> <p>This was incorrectly changed to: $Results\ (ug/g) = \frac{Total\ carbon}{(500 / Sample\ Amount)} * TS$</p> <p>From December 29, 2014 to April 13, 2015, this error was often discovered during data review and manually corrected. The problem was not addressed with the MIS department. However, there were several cases where this error was not caught during data review and results were reported incorrectly. On April 13, 2015, this issue was brought to the attention of MIS and the sample calculation was corrected back to:</p> <p>$Results\ (ug/g) = \frac{Total\ carbon}{(sample\ amount / 1000)} * (100/TS)$ (where 1000 is a conversion factor)</p> <p>Blanks and LCSs were not affected by these changes because they are calculated using different calculations since total solids is not used in these situations. Duplicates and MS/MSD samples are also not affected because they use different calculations since these calculation strings involve recoveries, RPDs, etc.</p> | | |
| Associated Non-Conformances: List logbook and page numbers | | |
| There are no non-conformances associated with this corrective action. | | |
| Root Cause Investigation & Determination (To be completed by Department Manager, Operations Manager and/or QA Officer) | | |
| Review the 6 "M's" below and investigate to determine whether one of them, or more than one, could be the cause of the problem. | | |
| Possible Causes | Details | |

KATAHDIN ANALYTICAL SERVICES, INC. – CORRECTIVE ACTION REPORT

| | |
|--|--|
| Machine (Instrument) | KIMS – incorrect formulas were entered into the process chains for TOC in soil. |
| Method (or Process) | <p>Katahdin's policy for data review: From the QAM: "For data that are reduced via computer, calculations are checked by the analyst (or designee) assigned to this task at a frequency designed to assure that the final data generation is valid." From SOP SD-904, Data Reduction, Review and Reporting: "All manual integrations, calculations and transcriptions are checked and 10% of all spreadsheet calculations are checked. The remainder of spreadsheet calculations is spot checked for potential anomalies."</p> <p>Katahdin does not have a formal process for handling calculation changes to process changes in KIMS. A verbal request is made to MIS and the change is made. Although the KIMS system does have an audit trail to track dates of changes and calculation changes, the request for these is not documented.</p> |
| Materials | Not Applicable |
| Maintenance (Is something not working correctly?) | Not Applicable |
| Man (training, human error) | All personnel involved with data review are aware of the requirement to hand check 10% of the sample results. In some cases this appears to have been done and when the results did not calculate correctly, they were corrected in KIMS. There was no communication that the error may affect other batches of samples. In other sample batches, the 10% hand check must not have occurred. |
| Mother Nature (accidents, power issues, beyond our control) | Not Applicable |

Corrective Action Plan: Name: Leslie Dimond

Date: 08/28/2015

Details of Corrective Action Plan –

Queries were run through KIMS to generate lists for TOC in soil data entered into KIMS between two timeframes:

10/23/14 to 12/29/14 (incorrect LOQ/LOD/MDL adjustment) – 14 Work Orders were found in this timeframe

12/29/14 and 4/13/15 (incorrect sample calculation) - 16 Work Orders were found in this timeframe

The lab has reviewed all TOC in soil data from these time periods for accuracy. Some inaccurate data was found (as expected from the incorrect formulas). Some data was found to be correct. In these cases, MIS was able to determine, through the KIMS audit trail, that the data had been manually corrected.

A new form has been created for personnel to fill out when requesting calculation changes in KIMS. This form will ensure that MIS is clear on exactly what needs to be changed in KIMS. Through KIMS we are currently able to track formula changes (ie. the formula before and after the change) and when the change occurred. This new form will allow Katahdin to track the changes made by person requesting the change and why the change was necessary. Supervisor and QA/Management approval are required on these forms.

There also will be a mandatory retraining for all employees on Katahdin's policy for data review, stressing that 10% of all results generated from KIMS and spreadsheets need to be confirmed. Also, this retraining will stress the need for communication. If a problem is discovered with one batch of samples, then it might also be affecting other batches.

Additionally, when calculation changes are made, Katahdin must establish a time frame of 30 days or 10 workorders where

KATAHDIN ANALYTICAL SERVICES, INC. – CORRECTIVE ACTION REPORT

senior management or the Quality Assurance Officer must also check the new calculations.

Review & Approval of Corrective Action Plan

Supervisor Approval: *[Signature]* Date: 09.11.15

Operations Manager Approval: *Deborah J. Hadeau* Date: 9.11.15

Quality Assurance Officer: *Lisee Dimond* Date: 09.11.15

Monitoring of Corrective Action (To be completed by QA Officer and/or Operations Manager): List details of follow-up

[Empty space for monitoring details]

| Corrective Action Effective | Return to Control – | Yes | No | Further Monitoring Needed/Additional Corrective Action |
|--|---------------------|-----|----|--|
| <i>[Empty space for monitoring data]</i> | | | | |

QA Approval: _____ Date: _____

Additional Information:

[Empty space for additional information]

Attachment B
Final Results after Data Review

| Sample Delivery Group | | | | SI0594 | | SI0594 | |
|-----------------------|----------------------|--------|-------|-----------------------------|------|-------------------|------|
| Lab ID | | | | SI0594-1 | | SI0594-2 | |
| Sample ID | | | | RE122D1-SOIL-012315-528-530 | | RE122D1-EB-012315 | |
| Sample Date | | | | 1/23/2015 | | 1/23/2015 | |
| Sample Type | | | | Soil | | Equipment Blank | |
| Method | Analyte | CAS No | Units | Result | Qual | Result | Qual |
| 5310B | TOTAL ORGANIC CARBON | -28 | MG_L | NA | | 0.24 | J |
| 9060A | TOTAL ORGANIC CARBON | -28 | UG_G | 250 | J | NA | |

Notes:

ID = Identification
MG_L = Milligrams per liter
UG_G = Micrograms per gram
NA = Not analyzed
Qual = Final qualifier

Final Qualifier:

J = The analyte was positively identified. The numerical value is the estimated concentration of the analyte in the sample.



DATA VALIDATION REPORT

| | | |
|-------------------------|--|---|
| Project: | Regional Groundwater Investigation — NWIRP Bethpage | |
| Laboratory: | Katahdin Analytical | |
| Sample Delivery Groups: | TH0533 | |
| Analyses/Method: | Total Organic Carbon (TOC) by U.S. EPA SW-846 Method 9060A and Standard Method 5310B for Total Organic Carbon by High-Temperature Combustion | |
| Validation Level: | 3 | |
| Project Number: | 0888812477.SA.DV | |
| Prepared by: | Dana Miller/Resolution Consultants | Completed on: 02/3/2015 Revised on: 10/27/2015 |
| Reviewed by: | Tina Cantwell/Resolution Consultants | File Name: TH0533_9060A_5310B |

SUMMARY

This report summarizes data review findings for samples listed below, collected by Resolution Consultants from the Regional Groundwater Investigation — NWIRP Bethpage site on 5 December 2014 in accordance with the following Sampling and Analysis Plans:

- *Sampling and Analysis Plan, Bethpage, New York.* (Resolution Consultants April 2013).
- *UFP SAP Addendum, Installation of Vertical Profile Borings and Monitoring Wells, Operable Unit 2, NWIRP Bethpage, New York.* (Resolution Consultants November 2013).
- *UFP SAP Addendum, Inclusion of Additional Target Analytes for Volatile Organics Analyses, NWIRP Bethpage OU2, Bethpage, New York.* (Resolution Consultants August 2014).

| Sample ID | Lab ID | Matrix/Sample Type | Analysis |
|-----------------------------|----------|--------------------|----------|
| RE122D3-SOIL-120514-718-720 | TH0533-1 | Soil | 9060A |
| RE122D3-EB-120514 | TH0533-2 | Equipment Blank | 5310B |

Data validation activities were conducted using the following guidance documents: *Test Methods for Evaluating Solid Waste, Physical/Chemical Methods SW-846, specifically Method 9060A, Total Organic Carbon* (U.S. EPA, 1996), *Method SM5310B, Total Organic Carbon by High-Temperature Combustion*, *U.S. Environmental Protection Agency (U.S. EPA) Contract Laboratory Program National Functional Guidelines for Inorganic Superfund Data Review* (NFG, January 2010, and Department of Defense (DoD) Quality Systems Manual (QSM) for Environmental Laboratories, Version 4.2 (October 2010). In

the absence of method-specific information, laboratory quality control (QC) limits, project-specific requirements and/or professional judgment were used as appropriate.

REVIEW ELEMENTS

The data were evaluated based on the following parameters (where applicable to the method):

- ✓ Data completeness (chain-of-custody)/sample integrity
- ✓ Holding times and sample preservation
- NA Gas chromatography/Mass spectrometer performance checks
- ✓ Initial calibration/continuing calibration verification
- ✓ Laboratory blanks/equipment blanks/field blanks/trip blanks
- NA Surrogate spike recoveries
- ✓ Matrix spike and/or matrix spike duplicate results
- ✓ Laboratory control sample laboratory control sample duplicate results
- NA Field duplicates
- NA Internal standards
- ✓ Sample results/reporting issues

The symbol (✓) indicates that no validation qualifiers were applied based on this parameter. NA indicates that the parameter was not included as part of this data set or was not applicable to this validation and therefore not reviewed. Acceptable data parameters for which all criteria were met and no qualification was performed and non-conformance or other issues that were noted during validation, but did not result in qualification of data are not discussed further.

Qualifications Actions

The data was reviewed independently from the laboratory to assess data quality. All compounds detected at concentrations less than the limit of quantitation but greater than the method detection limit were qualified by the laboratory as estimated (J). This "J" qualifier was retained during data validation. Any sample that was analyzed at a dilution because of high concentrations of target or non-target analytes was checked to confirm that the results and/or sample-specific limit of quantitation and limit of detections were adjusted accordingly by the laboratory.

No results were qualified during this review. Analytical completeness was calculated to be 100% and the data are usable for their intended purpose, according to U.S. Environmental Protection Agency and Department of Defense guidelines. Katahdin Analytical discovered a QC error in Total Organic Carbon calculations, informed us, and has implemented corrective action provided in Attachment A. Attachment B provides final results after data review.



ATTACHMENTS

Attachment A: Katahdin Analytical Corrective Action Report

Attachment B: Final Results after Data Review

Attachment A
Katahdin Analytical Corrective Action Report

KATAHDIN ANALYTICAL SERVICES, INC. – CORRECTIVE ACTION REPORT

| | | |
|--|------------------------------------|------------------|
| Problem Identification (Person initiating CAR) Name: Leslie Dimond | | Date: 08/28/2015 |
| Discovered by Laboratory | X Discovered by Client (Complaint) | Other |
| Details of Problem: | | |
| <p>On August 28, 2015, a client called regarding several TOC results. They thought the results for their samples seemed to be off by a factor of ten. George Brewer looked into the matter and found that their sample results were calculated incorrectly. Calculations are set up to be performed automatically in KIMS. Each product code for a test contains information such as the method name, LOQ, LOD, MDL, limits and a specific process chain for reporting to the either the LOQ, LOD or MDL (i.e. three process chains). This process chain contains all of the calculations associated with a test including separate calculations for sample results, adjusted LOQs, LODs and MDLs, and results for QC samples. In October of 2014, a client asked why the lab was not adjusting LOQs, LODs and/or MDLs for the sample amount used, but only for an instrument dilution and total solids. The lab had felt that there wasn't really a standard weight amount, so this was not incorporated. After numerous discussions, the lab decided to change this and use 500 mg as the standard and to incorporate any different amount into the LOQ/LOD/MDL adjustment. MIS was asked to revise this TOC calculation in our Laboratory Information System (KIMS). The incorrect information was given to MIS, so the adjusted LOQ/LOD/MDL calculation was wrong.</p> <p>The formula was: $ADJ. LOQ/LOD/MDL = LOQ/LOD/MDL * DF * (Sample\ amount / 1000) * (100/TS)$.</p> <p>During data review, sample results that are calculated automatically by KIMS, are checked at a frequency of 10% of all calculations. Sample results are checked in this manner, but adjusted limits are not necessarily checked. In December of 2014, a data reviewer noticed that some adjusted limits were not correct. This issue was reviewed with MIS, and it was discovered that the calculation was incorrect. It was corrected at this time and changed to:</p> <p>$ADJ. LOQ/LOD/MDL = LOQ/LOD/MDL * DF * (500 / Sample\ Amount) * (100/TS)$ (where 500 is the standard sample amount).</p> <p>At the time this change was made, MIS was under the impression that the sample calculation was also incorrect and also needed to be corrected.</p> <p>The formula for the sample result was: $Results\ (ug/g) = \frac{Total\ carbon}{(sample\ amount / 1000)} * (100/TS)$ (where 1000 is a conversion factor)</p> <p>This was incorrectly changed to: $Results\ (ug/g) = \frac{Total\ carbon}{(500 / Sample\ Amount)} * TS$</p> <p>From December 29, 2014 to April 13, 2015, this error was often discovered during data review and manually corrected. The problem was not addressed with the MIS department. However, there were several cases where this error was not caught during data review and results were reported incorrectly. On April 13, 2015, this issue was brought to the attention of MIS and the sample calculation was corrected back to:</p> <p>$Results\ (ug/g) = \frac{Total\ carbon}{(sample\ amount / 1000)} * (100/TS)$ (where 1000 is a conversion factor)</p> <p>Blanks and LCSs were not affected by these changes because they are calculated using different calculations since total solids is not used in these situations. Duplicates and MS/MSD samples are also not affected because they use different calculations since these calculation strings involve recoveries, RPDs, etc.</p> | | |
| Associated Non-Conformances: List logbook and page numbers | | |
| There are no non-conformances associated with this corrective action. | | |
| Root Cause Investigation & Determination (To be completed by Department Manager, Operations Manager and/or QA Officer) | | |
| Review the 6 "M's" below and investigate to determine whether one of them, or more than one, could be the cause of the problem. | | |
| Possible Causes | Details | |

KATAHDIN ANALYTICAL SERVICES, INC. – CORRECTIVE ACTION REPORT

| | |
|--|--|
| Machine (Instrument) | KIMS – incorrect formulas were entered into the process chains for TOC in soil. |
| Method (or Process) | <p>Katahdin's policy for data review: From the QAM: "For data that are reduced via computer, calculations are checked by the analyst (or designee) assigned to this task at a frequency designed to assure that the final data generation is valid." From SOP SD-904, Data Reduction, Review and Reporting: "All manual integrations, calculations and transcriptions are checked and 10% of all spreadsheet calculations are checked. The remainder of spreadsheet calculations is spot checked for potential anomalies."</p> <p>Katahdin does not have a formal process for handling calculation changes to process changes in KIMS. A verbal request is made to MIS and the change is made. Although the KIMS system does have an audit trail to track dates of changes and calculation changes, the request for these is not documented.</p> |
| Materials | Not Applicable |
| Maintenance (Is something not working correctly?) | Not Applicable |
| Man (training, human error) | All personnel involved with data review are aware of the requirement to hand check 10% of the sample results. In some cases this appears to have been done and when the results did not calculate correctly, they were corrected in KIMS. There was no communication that the error may affect other batches of samples. In other sample batches, the 10% hand check must not have occurred. |
| Mother Nature (accidents, power issues, beyond our control) | Not Applicable |

Corrective Action Plan: Name: Leslie Dimond

Date: 08/28/2015

Details of Corrective Action Plan –

Queries were run through KIMS to generate lists for TOC in soil data entered into KIMS between two timeframes:

10/23/14 to 12/29/14 (incorrect LOQ/LOD/MDL adjustment) – 14 Work Orders were found in this timeframe

12/29/14 and 4/13/15 (incorrect sample calculation) - 16 Work Orders were found in this timeframe

The lab has reviewed all TOC in soil data from these time periods for accuracy. Some inaccurate data was found (as expected from the incorrect formulas). Some data was found to be correct. In these cases, MIS was able to determine, through the KIMS audit trail, that the data had been manually corrected.

A new form has been created for personnel to fill out when requesting calculation changes in KIMS. This form will ensure that MIS is clear on exactly what needs to be changed in KIMS. Through KIMS we are currently able to track formula changes (ie. the formula before and after the change) and when the change occurred. This new form will allow Katahdin to track the changes made by person requesting the change and why the change was necessary. Supervisor and QA/Management approval are required on these forms.

There also will be a mandatory retraining for all employees on Katahdin's policy for data review, stressing that 10% of all results generated from KIMS and spreadsheets need to be confirmed. Also, this retraining will stress the need for communication. If a problem is discovered with one batch of samples, then it might also be affecting other batches.

Additionally, when calculation changes are made, Katahdin must establish a time frame of 30 days or 10 workorders where

Attachment B
Final Results after Data Review

| Sample Delivery Group | | | | TH0533 | | TH0533 | |
|-----------------------|----------------------|--------|-------|-----------------------------|------|-------------------|------|
| Lab ID | | | | TH0533-1 | | TH0533-1 | |
| Sample ID | | | | RE122D3-SOIL-120514-718-720 | | RE122D3-EB-120514 | |
| Sample Date | | | | 12/5/2014 | | 12/5/2014 | |
| Sample Type | | | | Soil | | Equipment Blank | |
| Method | Analyte | CAS No | Units | Result | Qual | Result | Qual |
| 5310B | TOTAL ORGANIC CARBON | -28 | MG_L | NA | | 0.30 | J |
| 9060A | TOTAL ORGANIC CARBON | -28 | UG_G | 280 | J | NA | |

Notes:

ID = Identification
MG_L = Milligrams per liter
UG_G = Micrograms per gram
NA = Not analyzed
Qual = Final qualifier

Final Qualifier:

J = The analyte was positively identified. The numerical value is the estimated concentration of the analyte in the sample.

Section 5

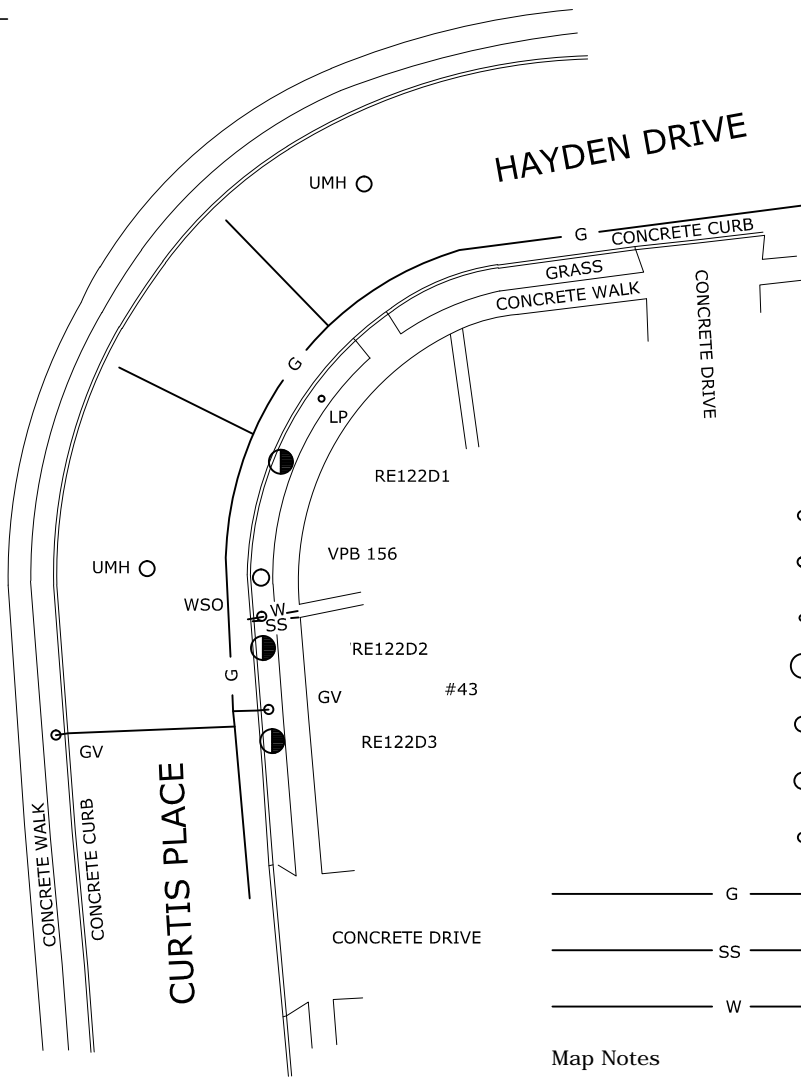
Survey

UNAUTHORIZED ALTERATION OR ADDITION TO THIS DOCUMENT IS A VIOLATION OF SECTION 7209 SUBDIVISION 2 OF THE NEW YORK STATE EDUCATION LAW.

| Description | Northing | Easting | Latitude | Longitude | Ground | Rim | PVC |
|-------------|-----------|------------|--------------|--------------|--------|-------|-------|
| VPB 156 | 207800.08 | 1124978.89 | N40-44-09.31 | W73-29-31.89 | 97.43 | NA | NA |
| RE122D1 | 207818.19 | 1124982.01 | N40-44-09.49 | W73-29-31.85 | 97.74 | 97.74 | 97.42 |
| RE122D2 | 207789.07 | 1124979.19 | N40-44-09.20 | W73-29-31.89 | 97.70 | 97.73 | 97.35 |
| RE122D3 | 207774.52 | 1124980.63 | N40-44-09.06 | W73-29-31.87 | 97.62 | 97.66 | 97.27 |



BENCHMARK FOUND
"X-CUT" NHOA ON HYDRANT
ELEVATION=100.40'

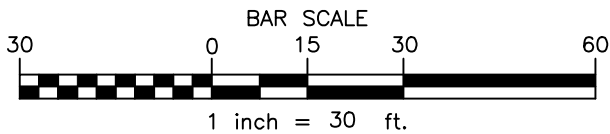


Legend

- GV Gas Valve
- ⊕ HYD Hydrant
- LP Light Post
- MW Monitoring Well
- UMH Unknown Manhole
- VPB 156 Vertical Profile Boring
- WSO Water Shutoff
- G — Underground Gas Line
- SS — Underground Sanitary Line
- W — Underground Water Line

Map Notes

1. Information shown hereon was compiled from an actual field survey conducted from September 23-25, 2014.
2. North orientation is Grid North based on the New York State Plane Coordinate System, Long Island Zone, NAD 83 as obtained from GPS observations.
3. Vertical datum shown hereon is NAVD 88 as obtained from GPS observations.



DWG NO. 14-498

| Date | RECORD OF WORK | Appr. |
|---------------|-----------------------|-------|
| 7-7-15 | ADDITIONAL FIELD WORK | JFC |
| | | |
| | | |
| | | |
| | | |
| Drafter: LMK | Checker: JFC | |
| Appr. by: JFC | Proj. No. 14.4121 | |

VERTICAL PROFILE BORING 156 SURVEY LOCATION
43 CURTIS PLACE

TOWN OF BETHPAGE NASSAU COUNTY, NEW YORK

C.T. MALE ASSOCIATES
Engineering, Surveying, Architecture & Landscape Architecture, D.P.C.

50 CENTURY HILL DRIVE, LATHAM, NY 12110
518.786.7400 * FAX 518.786.7299



SCALE: 1"=30'

DATE: SEPT. 24, 2014