2015 OU2 GROUNDWATER INVESTIGATION RE107D1, RE107D2, RE107D3 (VPB141) INSTALLATION REPORT

NAVAL WEAPONS INDUSTRIAL RESERVE PLANT (NWIRP) SITE 1 OU2 BETHPAGE, NY

Prepared for:



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

March 2016

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Prepared by:



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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

ESS Environmental Sequence Stratigraphy

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

NYS New York State

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
PPE Personal Protective Equipment

PVC Polyvinylchloride

SAP Sampling and Analysis Plan

SVOC Semivolatile Organic Compounds

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, one initial sampling event and one quarterly groundwater monitoring event (specifically at the Vertical Profile Boring [VPB] 141 location) in 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 three wells (RE107D1, RE107D2 and RE107D3) associated with the VPB141 location. The purpose of this investigation was to ascertain contaminant levels and depths south of the On-site Containment Treatment System (ONCT), west of Wantagh Avenue and northwest of the RE108 Hot Spot, to provide information on the effectiveness of the ONCT, and to evaluate the possibility of contaminant influx to the RE108 hotspot from the northwest. The locations of RE107D1, RE107D2 and RE107D3, 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 2015 in accordance with the *United Federal Programs Sampling and Analysis Plan (UFP SAP)*, Bethpage, New York (Resolution, 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 and Appendix B 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), as observed at the former NWIRP and extending southeast to areas south of Southern State Parkway. Locally at the RE107 locations, the bottom of the Magothy (top of the Raritan Clay) is encountered at approximately 838 feet bgs. 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, and above depths of approximately 500 feet bgs are laterally discontinuous. Preliminary cross section results from Environmental Sequence Stratigraphy (ESS) analysis suggests that below depths of approximately 500 feet bgs some clay layers may be persistent in a north-south direction; additional ESS cross section analyses are planned for the east-west direction.

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 (RE107 series wells) were installed in the vicinity of VPB141 between August and September 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 RE107D1, RE107D2 and RE107D3 were installed using mud rotary drilling techniques (Figure 2). Depths of monitoring wells RE107D1, RE107D2 and RE107D3 were 530 ft, 585 ft and 670 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. *2015 OU2 Groundwater Investigation VPB141* (Resolution Consultants, 2015) documents the installation of this VPB including detailed lithologic descriptions, continuous gamma plot and multiple Volatile Organic Compounds (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 RE107D1, RE107D2 and RE107D3 were initially sampled by Resolution Consultants on October 27-28, 2015 and were then included in the December 2015 quarterly monitoring event as part of the Navy's ongoing Environmental Restoration Program. Analytical results and stabilized field parameters for the October 2015 data are summarized in Table 3 and Table 4, respectively. Data validation is documented in Appendix A.

Analytical results and data validation for the December 2015 data are summarized in Appendix B. Complete documentation of the December sampling round is included in the *December 2015 Groundwater Sampling Data Summary Report* (Resolution Consultants, 2016).

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 NYS Net 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.

Resolution Consultants, 2013a. *United Federal Programs Sampling and Analysis Plan, Site OU-2 Offsite Trichloroethene (TCE) Groundwater Plume Investigation, Bethpage, New York.* April 2013.

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Resolution Consultants, 2015. 2015 OU2 Groundwater Investigation VPB141, Bethpage, NY. September 2015.

Resolution Consultants, 2016. *December 2015 Groundwater Sampling Data Summary Report, Bethpage, NY.* March 2016.

Smolensky, D., and Feldman, S., 1990. *Geohydrology of the Bethpage-Hicksville-Levittown Area, Long Island, New York, U.S.* Geological Survey Water-Resourced Investigations Report 88-4135, 25 pp.



TABLE 1 MONITORING WELL CONSTRUCTION SUMMARY 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)
RE107D1	9/25/2015	99.56	98.92	530	53	505 - 525	525 - 530	542
RE107D2	9/10/2015	99.57	98.99	585	53	560 - 580	580 - 585	597
RE107D3	8/25/2015	100.61	99.96	670	53	645 - 665	665 - 670	682

MSL - mean sea level

ft bgs - feet below ground surface

TABLE 2 MONITORING WELL DEVELOPMENT SUMMARY 2015 OU2 GROUNDWATER INVESTIGATION NWIRP BETHPAGE, NY

	AIR DEVEL	OPMENT.	PUM	1P DEVELOPME	APPROX. TOTAL	FINAL		
MONITORING WELL	DATE	APPROX. VOLUME (GAL)	DATE	FINAL PUMP DEPTH (FT BGS)	APPROX. VOLUME (GAL)	DEVELOPMENT VOLUME (GAL)	TURBIDITY (NTUs)	
RE107D1	10/1/2015	4,000	10/6/2015- 10/07/2015	505-525	8,000	12,000	26.46	
RE107D2	9/30/2015	2,200	10/5/2015	560-580	5,000	7,200	32.85	
RE107D3	10/2/2015	8,500	10/8/2015- 10/9/2015	645-665	5,000	13,500	27.02	

GAL - gallon

FT BGS - feet below ground surface NTUs - Nephelometric Turbidity Units

TABLE 3. ANALYTICAL DATA SUMMARY 2015 OU2 GROUNDWATER INVESTIGATION NWIRP BETHPAGE, NY

Location		RE107D1	RE107D2	RE107D2	RE107D3
Sample Date	NYSDEC Groundwater	10/27/2015	10/27/2015	10/27/2015	10/28/2015
Sample ID	Guidance or Standard Value	RE107DI-WG- 102715	RE107D2-WG- 102715	DUPLICATE-WG- 102715	RE107D3-WG- 102815
Sample type code	(Note 1)	N	N N	FD	N
VOC 8260C (ug/L)				1.5	
1,1,1-TRICHLOROETHANE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,1,2,2-TETRACHLOROETHANE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	5	0.82 J	14 J	15 J	1.9 J
1,1,2-TRICHLOROETHANE	1	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1.1-DICHLOROETHANE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,1-DICHLOROETHENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,2,4-TRICHLOROBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DIBROMO-3-CHLOROPROPANE	0.04	< 0.75 U	< 0.75 U	< 0.35 U	< 0.75 U
1,2-DIBROMOETHANE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DICHLOROBENZENE	3	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DICHLOROETHANE	5	< 0.50 U		< 0.50 U	< 0.50 U
1.2-DICHLOROETHANE 1.2-DICHLOROETHENE. TOTAL	5	< 0.50 U	< 0.50 U 2.2 J	< 0.50 U	< 0.50 U
1,2-DICHLOROPROPANE	1	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,3-DICHLOROPROPANE 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 < 0.50 U	< 0.50 U
•	1			†	
1,4-DIOXANE (Method 8270D_SIM) 2-BUTANONE	NL 50	8.2	10	12	< 0.18 U
	50	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 U
2-HEXANONE	50	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 U
4-METHYL-2-PENTANONE	NL	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ
ACETONE	50	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ
BENZENE	1	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
BROMODICHLOROMETHANE	50	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
BROMOFORM	50	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
BROMOMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
CARBON DISULFIDE	60	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CARBON TETRACHLORIDE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
CHLOROFORM	7	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROMETHANE	5	< 1.0 UJ	0.82 J	< 1.0 UJ	< 1.0 UJ
CIS-1,2-DICHLOROETHENE	5	< 0.50 U	2.2 J	2.2 J	< 0.50 U
CIS-1,3-DICHLOROPROPENE	0.4	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CYCLOHEXANE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
DIBROMOCHLOROMETHANE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
DICHLORODIFLUOROMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
ETHYLBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
ISOPROPYLBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
M- AND P-XYLENE	NL	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
METHYL ACETATE	NL	< 0.75 U	< 0.75 U	< 0.75 U	< 0.75 U
METHYL CYCLOHEXANE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
METHYL TERT-BUTYL ETHER	10	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
METHYLENE CHLORIDE	5	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ
O-XYLENE	NL	4.4 J	4.3 J	4.3 J	< 0.50 U
STYRENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TETRACHLOROETHENE	5	1.3 J	7.3 J	6.9 J	< 0.50 U
TOLUENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TRANS-1,2-DICHLOROETHENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TRANS-1,3-DICHLOROPROPENE	0.4	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TRICHLOROETHENE	5	14 J	140 J	150 J	< 0.50 U
TRICHLOROFLUOROMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
VINYL CHLORIDE	2	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
XYLENES, TOTAL	5	4.4 J	4.3 J	4.3 J	< 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 2015 OU2 GROUNDWATER INVESTIGATION NWIRP BETHPAGE, NY

Well	Date	Temperature (°C)	рН	Specific Conductance (µS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Depth to water (ft bgs)	Flow rate (ml/min)
RE107D1	10/27/2015	14.38	6.34	0.094	2.61	113.9	423	41.96	550
RE107D2	10/27/2015	14.79	5.73	81	3.51	179.2	>1000	42.52	600
RE107D3	10/28/2015	14.93	5.58	0.025	6.50	139.2	377	42.62	500

°C - degrees Celsius

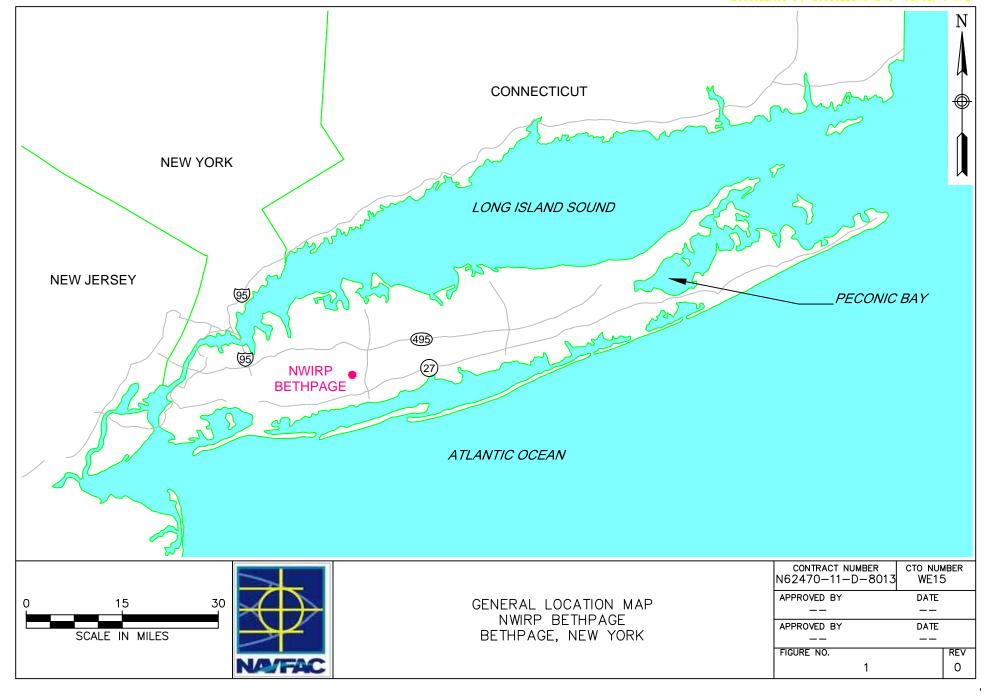
 $\mu S/cm$ - Microsiemens per Centimeter

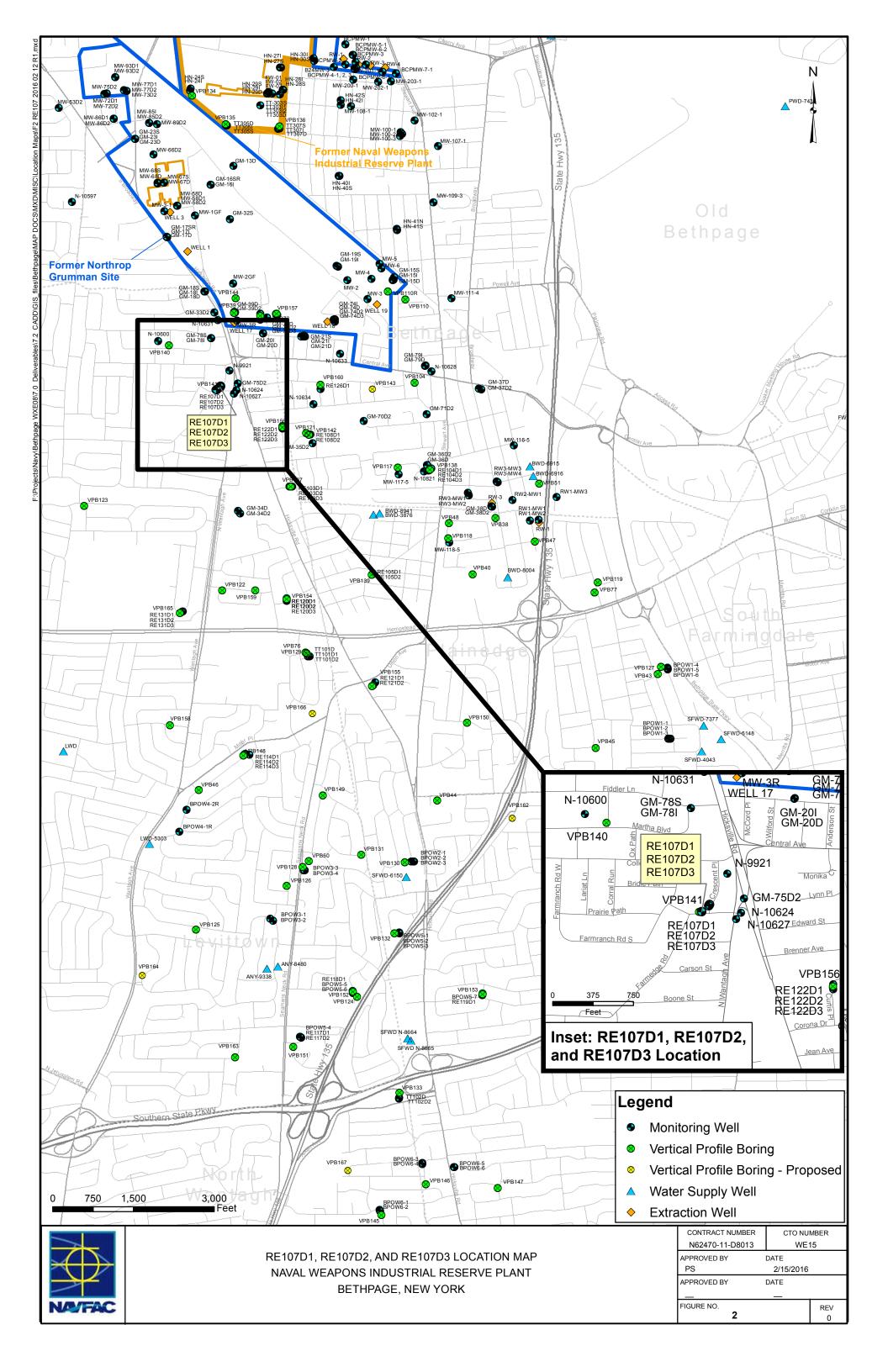
mg/L - milligrams per liter

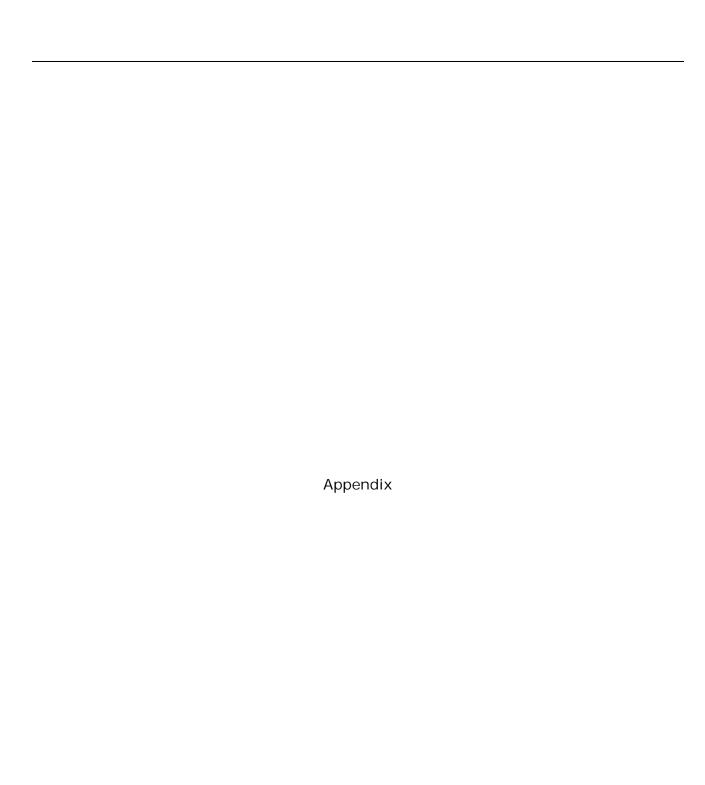
mV - Millivolts

NTU - Nephelometric Turbidity Unit ft bgs - feet below ground surface ml/min - milliliters per minute









Appendix A
RE107D1, RE107D2, RE107D3

Section 1

Boring Logs

Boring Log

BORING #: RE107D1 Sheet 1 of 2

Client: Department of the Navy, Naval Facilit	Logged By: V. Varricchio							
Location: Farmedge Rd. & Crescent PI, Levitt	Drilling Company: Delta Well & Pump							
Project #: 60266526	Ground Elevation (msl): 99.56	Well Screen Interval (ft): 505-525						
Start Date: 9/18/2015	Drilling Method: Auger (0-50' bgs) Mud Rotary (>50' bgs)	Water Level (ft):						
Finish Date: 9/25/2015	Northing: 208546.42 Easting: 1123823.36	Total Depth (ft): 542.0						

						T	
DEPTH (ft)	PID (ppm)	Formation	nscs	GRAPHIC LOG	MATERIAL DESCRIPTION	Well Completion	Well Construction
0					0-505 ft bgs: See VPB141 for Descriptions		10" Diameter Steel Casing
50							
100							
150							
200							
250						-	Bentonite Grout
300							
350							
400						-	4" Diameter Schedule 80 PVC Riser

Boring Log

BORING #: RE107D1 Sheet 2 of 2

Client: Department of the Navy, Naval Facilit	Logged By: V. Varricchio			
Location: Farmedge Rd. & Crescent PI, Levitt	Drilling Company: Delta Well & Pump			
Project #: 60266526	Ground Elevation (msl): 99.56	Well Screen Interval (ft): 505-525		
Start Date: 9/18/2015	Drilling Method: Auger (0-50' bgs) Mud Rotary (>50' bgs)	Water Level (ft):		
Finish Date: 9/25/2015	Northing: 208546.42 Easting: 1123823.36	Total Depth (ft): 542.0		

DEPTH (ft)	PID (ppm)	Formation	SOSN	GRAPHIC LOG	MATERIAL DESCRIPTION	Well	Well Construction
446 448 450 452 454					0-505 ft bgs: See VPB141 for Descriptions (continued)		4" Diameter Schedule 80 PVC Riser (continued)
456 458 460 462 464 466 468 470 472 474 476 478						•	#00 Filter Sand
482 484 486 488 490 492 494 496 498 500 502					Light gray (10VP 7/1) Clayou fine to medium well graded		#1 Filter Sand
506 508 510	0 ,		SC SP		Light gray (10YR 7/1) Clayey fine to medium well graded SAND with trace lignite Light gray (10YR 7/1) poorly graded medium subrounded SAND with trace Clay and bands of lignite		
514 516 518	0,		SP		Brownish yellow (10YR 6/6) poorly graded medium subangular SAND with trace Clay		4" Diameter Schedule 80 PVC, 10 Slot Well Screen
520 522 524	0 ,		SP		Brownish yellow (10YR 6/6) poorly graded medium subangular SAND with trace Clay		(505-525 ft bgs)
526 528 530 532 534							Sump
536 538 540							#1 Sand to Bottom
542					End of boring at 542.0 ft. bgs.		

Boring Log

BORING #: RE107D2 Sheet 1 of 2

Client: Department of the Navy, Naval Facilit	Logged By: V. Varricchio						
Location: Farmedge Rd. & Crescent PI, Levitt	Drilling Company: Delta Well & Pump						
Project #: 60266526	Ground Elevation (msl): 99.57	Well Screen Interval (ft): 560-580					
Start Date: 9/2/2015	Drilling Method: Auger (0-50' bgs) Mud Rotary (>50' bgs)	Water Level (ft):					
Finish Date: 9/10/2015	Northing: 208558.7 Easting: 1123830.8	Total Depth (ft): 597.0					

ОЕРТН (ft)	PID (ppm)	Formation	nscs	GRAPHIC LOG	MATERIAL DESCRIPTION	Well	Well Construction
0					0-563 ft bgs: See VPB141 for Descriptions		10" Diameter Steel Casing
50							
100							
150							
200							
250						→	Bentonite Grout
300							
350							
400							
450						•	4" Diameter Schedule 80 PVC Riser
							80 PVC Riser

Boring Log

BORING #: RE107D2 Sheet 2 of 2

Client: Department of the Navy, Naval Facilit	Logged By: V. Varricchio						
Location: Farmedge Rd. & Crescent PI, Levitt	Drilling Company: Delta Well & Pump						
Project #: 60266526	Ground Elevation (msl): 99.57	Well Screen Interval (ft): 560-580					
Start Date: 9/2/2015	Drilling Method: Auger (0-50' bgs) Mud Rotary (>50' bgs)	Water Level (ft):					
Finish Date: 9/10/2015	Northing: 208558.7 Easting: 1123830.8	Total Depth (ft): 597.0					

#1 Filter Sand #1 Filter Sand	DEPTH (ft)	PID (ppm)	Formation	nscs	GRAPHIC LOG	MATERIAL DESCRIPTION	Well	Well Construction
#100 Filter Sand #00 Filte	500 502 504 506					0-563 ft bgs: See VPB141 for Descriptions (continued)		Schedule 80 PVC
Signature Sign	510 512 514 516 518 520 522 524 526						4	#00 Filter Sand
Secondary Seco	530 532 534 536 538 540 542 544 546 548 550 552							#1 Filter Sand
576 578 580 580 582 584 586 588 590 590 5P SC Light gray (10YR 7/1) Clayey SAND with lamination Sump Sump ### Send to Pottors ### Send to Pottors	560 562 564 566 568 570	_						4" Diameter
582 584 586 588 590	574 576 578							Schedule 80 PVC, 10 Slot Well Screen (560-580 ft bgs)
	582 584 586							Sump
592 ————————————————————————————————————	590 592 594							#1 Sand to Bottom

Boring Log

BORING #: RE107D3
Sheet 1 of 2

Client: Department of the Navy, Naval Facilit	Logged By: V. Varricchio		
Location: Farmedge Rd. & Crescent PI, Levitt	Drilling Company: Delta Well & Pump		
Project #: 60266526	Ground Elevation (msl): 100.61	Well Screen Interval (ft): 645-665	
Start Date: 8/14/2015	Drilling Method: Auger (0-50' bgs) Mud Rotary (>50' bgs)	Water Level (ft):	
Finish Date: 8/25/2015	Northing: 208495.32 Easting: 1123759.78	Total Depth (ft): 682.0	

			1			1		
DEPTH (ft)	PID (ppm)	Formation	nscs	GRAPHIC LOG	MATERIAL DESCRIPTION	Well	Completion	Well Construction
0					0-647 ft bgs: See VPB141 for Descriptions			10" Diameter Steel Casing
50								Caong
100								
150							1000	
200								
250								
300							-	Bentonite Grout
350								
400								
450								
500								
550							*	4" Diameter Schedule 80 PVC Riser

Boring Log

BORING #: RE107D3 Sheet 2 of 2

Client: Department of the Navy, Naval Facilit	Logged By: V. Varricchio		
Location: Farmedge Rd. & Crescent PI, Levitt	Drilling Company: Delta Well & Pump		
Project #: 60266526	Ground Elevation (msl): 100.61	Well Screen Interval (ft): 645-665	
Start Date: 8/14/2015	Drilling Method: Auger (0-50' bgs) Mud Rotary (>50' bgs)	Water Level (ft):	
Finish Date: 8/25/2015	Northing: 208495.32 Easting: 1123759.78	Total Depth (ft): 682.0	

DEPTH (ft) PID (ppm) Formation		Formation USCS GRAPHIC LOG	MATERIAL DESCRIPTION	Well	Well Construction		
586 588 590 592					0-647 ft bgs: See VPB141 for Descriptions (continued)		4" Diameter Schedule 80 PVC Riser (continued)
596 598 600 602 604 606 608							#00 Filter Sand
612 614 616 618 620 622 624							
628 630 632 634 636 638 640 642							#1 Filter Sand
646 648 650	0		SP		Light gray (10YR 7/1) poorly graded fine SAND with iron banding		
652 654 656	0		GW		Light gray (10YR 7/1) well graded subrounded fine to coarse GRAVEL with some well graded fine to coarse subrounded Sand		4" Diameter
658	0		SW-GW		White (10YR 8/1) well graded fine to coarse subrounded SAND mixed with fine to coarse subrounded Gravel		Schedule 80 PVC, 10 Slot Well Screen (645-665 ft bgs)
662	0		GW		Very pale brown (10YR 7/3) well graded fine to coarse subrounded GRAVEL with trace lean Clay and well graded fine		
666 668 670 672					to coarse subangular sand		Sump
674 676 678							#1 Sand to Bottom
680					End of boring at 682.0 ft. bgs.		

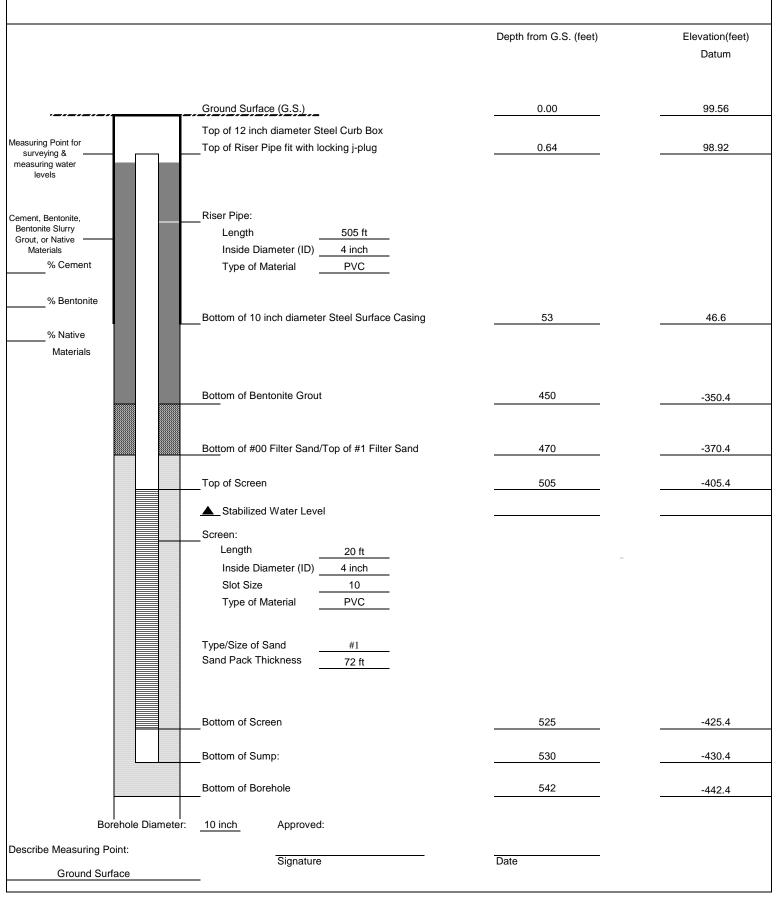
Section 2

Monitoring Well Construction Logs



Client: NAVFAC	Project Number: 60266526	WELL ID: RE107D1
Site Location: NWIRP	BETHPAGE, NY	
Well Location: Farmedg	e Rd. & Crescent Pl, Bethpage, NY	Date Installed: 9/18/2015 - 9/25/2015
Method: MUD ROTA	RY	Inspector: V. Varricchio
Coords: Northing: 20	8546.42 Fasting: 1123823.36	Contractor: DELTA WELL & PLIMP

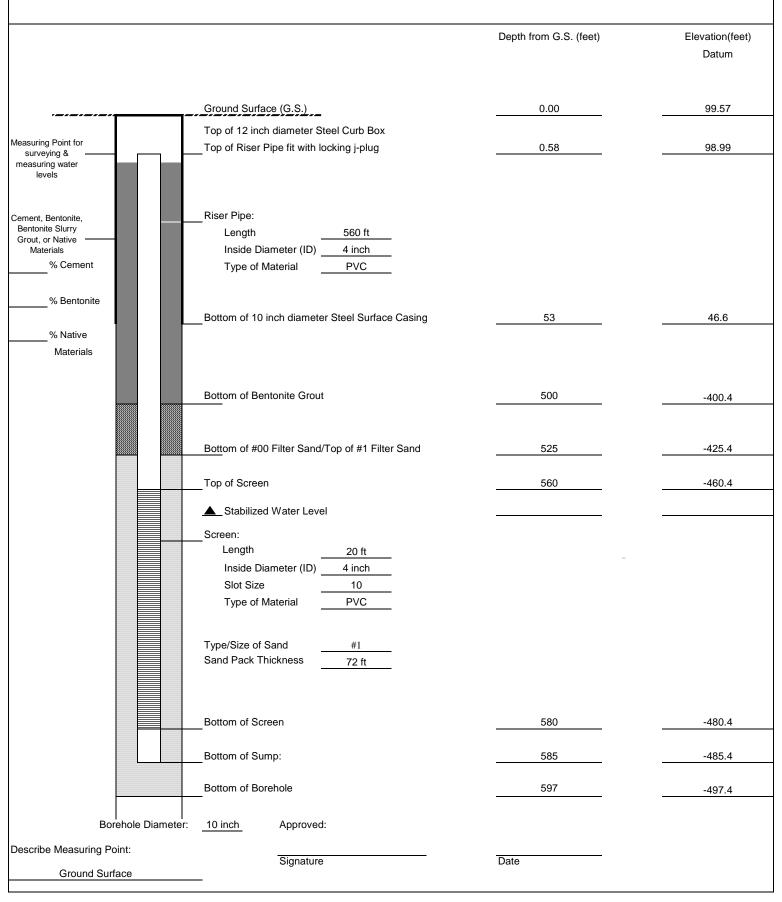
MONITORING WELL CONSTRUCTION DETAIL





Client:	NAVFAC	Project Number:	60266526	WELL	ID: RE107D2
Site Location	: NWIRP BETHPAC	E, NY			
Well Locatio	n: Farmedge Rd. & Cr	Date Installed:	9/2/2015 - 9/10/2015		
Method:	MUD ROTARY	Inspector:	V. Varricchio		
Coords:	Northing: 208558 70	Fasting: 1123830.8	0	Contractor	DELTA WELL & PLIMP

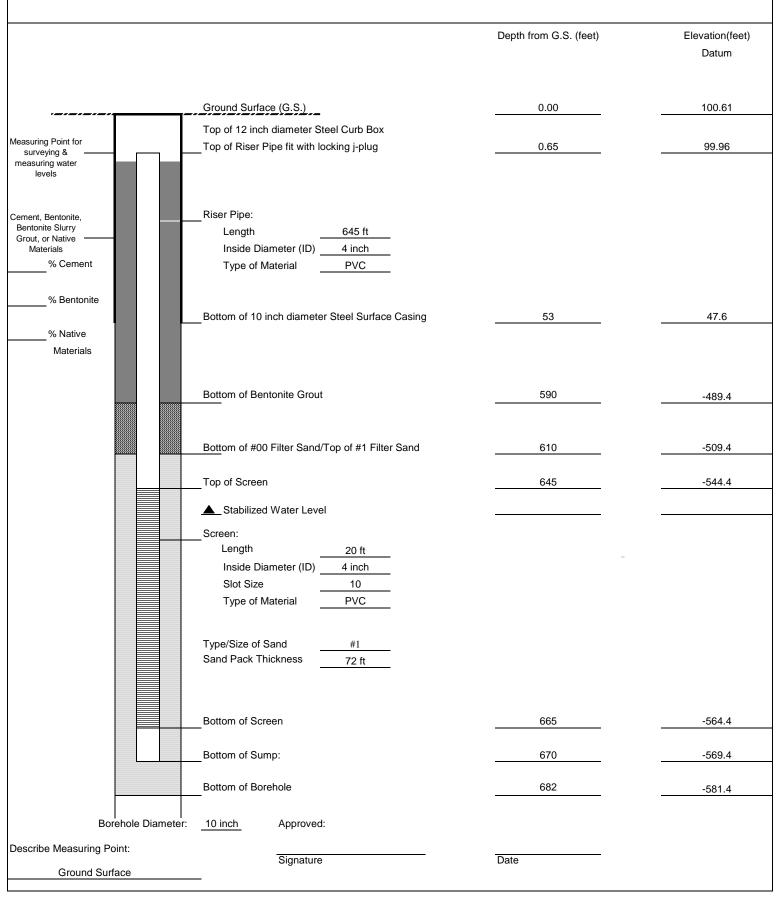
MONITORING WELL CONSTRUCTION DETAIL





Client: NAV	'FAC	Project Number:	60266526	WELL	ID: RE107D3
Site Location:	NWIRP BETHPAC	E, NY			
Well Location:	Farmedge Rd. & Cr	Date Installed:	8/14/2015 - 8/25/2015		
Method: MUI	O ROTARY	Inspector:	V. Varricchio		
Coords: Nort	hing: 208495 32	Fasting: 1123759 7	8	Contractor	DELTA WELL & PLIMP

MONITORING WELL CONSTRUCTION DETAIL



Section 3

Groundwater Sample Log Sheets

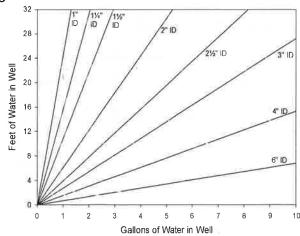


Well ID:	REIOT	Δ1
	110101	M-

Low Flow Ground Water Sample Collection Record

CONSUL	IANIS									
Client:		IWIRP Be				Date: 10) /	<u>/ 15</u>	Time: Start	
Project N	- 33	6026652							Finish	_ <i>[830]</i> am/pm
Site Loca Weather	125	Cre	0.1	600		Col	lector(s):	Pa.	1 Kareth	
vveatrier	Conds.	00	creast,	DU		001	160101(3).	- flax	INACICIMI	
1., WATE	R LEVEL	DATA: (n	neasured	from Top	of Casing	1)				
a. Tot	al Well Ler	ngth	530	.c. Length	of Water	Column		(a-b)	Casing Dia 4-inch PVC	meter/Material
b. Wa	ater Table D	Pepth -	42.05	d. Calcula	ated Syste	m Volume	e (see back)			1gal
	PURGE D		Geotech	bladder pu	mp with di	rop tube a	assembly			
	ceptance C							13		
	emperature		ilieu (see		Turbidity	± 10%		- D.O.	± 10% (val	ues >0.5 mg/L)
	•	± 0.1 ur	nit		- ORP		•		·	15
-	Sp. Cond.	± 3%		- C	rawdown	< 0.3'		Remove a	minimum 1	screen volume
c. Fie	ld Testing E	Equipmen	t used:		Make		Model		Serial Num	nber
					Hanna		18703		22335	
					y57.		556		73188)
	Volume			Spec.					,	
<u>Time</u>	Removed	Temp.	pН	Cond.	DO	ORP	Turbidity	Flow Rate	Depth to	Color/Odor
(24hr)	(Liters)	(°C)	10	(mS/cm)	(mg/L)	(mV)	(NTU)	(ml/min)	water (ft)	-
1615									, , , , , , , , , , , , , , , , , , ,	
17:00	Scal	14.70	(0.32	0.093	5,48	115.9	295	500	41.98	Grey elone
1720	-	14.75	6.34	0.094	2 F18	1151				,,,,,,
17.30	enr T	121.56		0.094	z.49	115.1	390	550	411.96	Great March
17: 33	10 Gast		6.32	0.095	2.69	114.9	429	550	41.96	/ /
17:48	gasthi.	14.49	631	0.095	2.66	114 4	434	550	41,94	Grey
	ceptance c	•	ss/fail		Yes	No	N/A	327		(continued on back)
	as required									
	as required ave parame	-		ieu		H	H			
- 10	If no or N					ш				
3. SAMF	PLE COLLE	ECTION:		Method:	Geotech	bladder p	ump with o	drop tube a	ssembly	
Sample I	D		Contai	ner Type	No. of Co	ontainers	Prese	ervation	Analys	is Reg. Time
				L vials	3			-ICI	VC	Cs
			1-L :	amber	2		n	one	1,4-Di	oxane
			. 3 7							
Commer	nts	USCO	1210	7						
Signature	2								Date	
Signature	- Je								-	
Signature	Day.							I	LowFlow-GWa	- Oct 2015 rev1.xls
Mac										

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

One screen volume (4-inch well)

15 ft = 37.1 L / 9.8 G 20 ft = 49.4 L / 13.1 G 25 ft = 61.8 L / 16.3 G

Well ID:

Continued from Time (24 hr)	volume Removed (Liters)	Temp (°C)	рН	Specific Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
17 55	- Martin	141.39	6.32	0.095	Z,59	11 4.4	426	550	211.99	Plantoly
18:00	_	14.39	6334			uH.1		550	41,96	
18:05	(BA)	14:40	10.28		2.58	114.0	409	550	411.98	Claudy
18:15	15 cm	14.38	6.24	0.094	2.61	112.9	4123	5.5.0	211.96	
								16		
									37	[[4]
										The state of the s
						*	3 -			
						-				
						9	17			
				8						
						9.				
							V.			, I
									747	

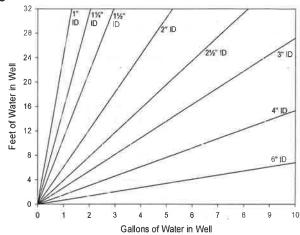


Well ID:	REINT	ומ	
	101		

Low Flow Ground Water Sample Collection Record

CONSUL	TANTS										
Client:	Navy N	IWIRP Be	thpage			Date: 10	1127	/ 15	Time: Start _		_am/pm
Project N		6026652			#					1830	_am/pm
Site Loca			nesemb	w 674		9 0 "		1211	1 10	- 11	
Weather	Conds:	_ grace	cast 6	0		Coll	ector(s):	I Kar	ch, F. Be	211	_
1. WATE	1. WATER LEVEL DATA: (measured from Top of Casing)										
a. Tot	al Well Ler	ngth	585	c. Length	of Water	Column		(a-b)	Casing Dia 4-inch PVC		iterial
b. Wa	ater Table D	Depth	42.35	d. Calcul	ated Syste	m Volume	e (see back)		13.19		
2. WELL	2. WELL PURGE DATA										
a. Pui	a. Purge Method: Geotech bladder pump with drop tube assembly										
	ceptance C mperature		•		- Turbidity - ORP			- D.O.	± 10% (val	ues >0.5	mg/L)
-	Sp. Cond.		iit.	- [Drawdown			Remove a	minimum 1	screen vo	olume
c. Fie	ld Testing I	Equipmen	t used:	e T	Make	.3	Model		Serial Num	1	
					Heana		98703		22335		
	1.										
Time	Volume Removed	Temp.	pН	Spec. Cond.	DO	ORP	Turbidity	Flow Rate	Depth to	Color	/Odor
(24hr)	(Liters)	(°C)	<u> </u>	(mS/cm)	(mg/L)	(mV)	(NTU)	(ml/min)	water (ft)	COIOI	<u>/Odoi</u>
1500		/								OH.	
16/4		15.09	5.38	0.092	7.30	183.9				2.0	
1600	5 Gal	15 18	5.69	0.098	5.31	174.9	359	500	42.26		
17.10	*****	15,08	5 69	0.098		175.15	1		42.26	Grey (Joud .
1720		14.75	6.34	0.094	304	114.8	,		41.98		
1715			black	er bro	he D	ell on	-0 d-4	chlace	Cost	10 14	n.
	ceptance c	•		•	Yes	N6	N/A	7		(continued or	back)
	as required as required										
	ve parame			lou	3						
	If no or N				^	1 1. 1	L				
			2 hour	Mugo	you !	whole	Tay1				
3. SAMF	PLE COLLE	ECTION:		Method:	Geotech	bladder p	ump with	drop tube a	ssembly		
Sample I	D			ner Type	No. of Co	ontainers		ervation	Analysi	-	Time
REIOI	12-104-	1023/		L vials	3			HCI	VO 1 4 Di		1815
- 0	plick	00-102	1/2 I-L 8	amber			Ŋ	one	1,4-Di	UNAI IE	1800
Commen	ıts	410	uinto								
		-/	1,000								
		7									
Signature	9	Pecal	Karel))					_Date	10/27	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

One screen volume (4-inch well)

15 ft = 37.1 L / 9.8 G 20 ft = 49.4 L / 13.1 G 25 ft = 61.8 L / 16.3 G

Well ID: REIOZ - DZ

(continued fro	om front)						Y			
	Volume			Specific				Flow		
Time	Removed		рН	Cond.	DO	ORP	Turbidity	Rate	Depth to	Color/Odor
(24 hr)	(Liters)	(°C)		(mS/cm)	(mg/L)	(mV)	(NTU)	(ml/min)	water (ft)	
1725		1480	5.52	79	6.2.3	187.4	- 40	600		
1730		1484	5.60	74	5.12	183.0	393	(4)	42,34	
1735		1483	5.65	79	4.77	18/6				
1248		14.83	5.62	78	4.86	182.7	511			
1745	logal	14.82	5.69	80	4.41	182.3	659			
1250		1482	5.70	80	4.14	18/2	873			
1755		181	5.72	81	3.87	180.1	71000		4256	
1200	13gal	14.80	5.73	81	3.75	179.7		600		
1805	/ 10	14.80	5.74	8/	3.58	1789	71000		42.52	TA TA
1810	1596	14.79	573	8/	3.51	1792	4			Those purge.
1818	Lend	e	Buk	Lite						
	1		10 W/	·/Ch						
			-							
			-							
						•				
						0				

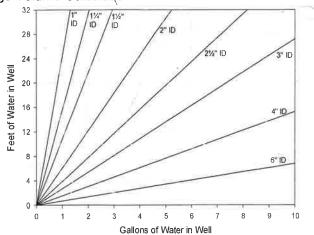


Well ID:	RZ	10703	
	1	1000	

Low Flow Ground Water Sample Collection Record

CONSUL	TANTS										
Client:	` Navy N	WIRP Be	thpage			Date: 10	1128	15	Γime: Start _	800	_am/pm
Project N		6026652		+			×		Finish	1830	am/pm
Site Loca			seub,	w pp ma				100	J -	,0 .1	
Weather	Conds:	OVERCA	CH, HILL	14,550		Col	lector(s):	/ X/	erch F.	hell	-
1. WATE	ER LEVEL	DATA: (n	neasured	from Top	of Casing	1)					
a. To	tal Well Ler	ngth	43.45	c. Length	of Water	Column		(a-b)	Casing Dia	meter/Ma	terial
7							_		4-inch PVC		
b. Wa	ater Table [Depth	4335	d. Calcula	ated Syste	m Volume	e (see back)		13.	1901	
2. WELI	PURGE D	ATA									
a. Pu	rge Method	<u>:</u>	Geotech	bladder pu	mp with d	rop tube a	assembly				
b. Ac	ceptance C	riteria def	ined (see	workplan)							
- Te	emperature			-	Turbidity			- D.O.	± 10% (valu	ues >0.5 r	ng/L)
	- pH Sp. Cond.	± 0.1 ur	nit		ORP - Orawdown	± 10mV		Pemove a	minimum 1	screen vo	dume
_	Sp. Cond.	I 370		- L	Hawuowii	~ 0.3		Nelliove a	IIIIIIIIIIIIIIIIIIIIIIIIIIIIIIIIIIIIII	SCIECTI VC	nume
c. Fie	ld Testing I	Equipmen	t used:		Make		Model	cur of	Serial Num		
				1	VSI		YSI.5		220	-535	
)	Hanna		7.8/0)	Sagar No.	7.5.3	
	Volume			Spec.							
<u>Time</u>	Removed	Temp.	<u>pH</u>	Cond.	<u>DO</u>	<u>ORP</u>	Control at the Control of the Control	Flow Rate	Depth to	Color/	<u>Odor</u>
(24hr)	(Liters)	(°C)		(mS/cm)	(mg/L)	(mV)	(NTU)	(ml/min)	water (ft)	h	
815							i i	500		ORI	
825		1490	5.36	0.033	5.92	163.3			43.55		
8:30		14.86	5.60	0.038	5.70	140.8					
8.35		14.86	5.67	6.030	5.68	136.7	66.3				
840	Þ.	14.87	5.77	0.028	5.47	125.7			43.15	Clo	nedy
845		1490	6.20	0.064	4.23	132.3	71004			reddist	brecon
d. Ac	ceptance o				Yes	No	N/A			(continued on	back)
	as required				P						
	as required ave parame	•		nea			H				
- 110	If no or N				45		- T	1. 75	2 (1		
				ž	Ehr 1	rural					
		FOTION			0 - 1 - 1						
3. SAMI	PLE COLLI	ECTION:		Method:	Geotech	bladder p	ump with c	drop tube as	ssembly		
Sample I	ID		Contai	ner Type	No. of Co	ontainers	Prese	ervation	Analysi	s Req.	Time
	13-104-10	2815	40-m	ıL vials	3		H	HCI	vo		1036
	V		1-L	amber	2		n	one	1,4-Die	oxane	1050
		-									
Commer	nts										
	.//	1 0 12	1.	_							
Signatur	e //a	ad N	ueth						Date		

Purge Volume Calculation



Volume / Linear Ft. of Pipe							
Gallon	Liter						
0.0025	0.0097						
0.0057	0.0217						
0.0102	0.0386						
0.0229	0.0869						
0.0408	0.1544						
0.0637	0.2413						
0.0918	0.3475						
0.1632	0.6178						
0.2550	0.9653						
0.3672	1.3900						
0.6528	2.4711						
1.4688	5.5600						
	Gallon 0.0025 0.0057 0.0102 0.0229 0.0408 0.0637 0.0918 0.1632 0.2550 0.3672 0.6528						

One screen volume (4-inch well)

15 ft = 37.1 L / 9.8 G 20 ft = 49.4 L / 13.1 G 25 ft = 61.8 L / 16.3 G

Well ID:	RE10713	2815
		u 0 0

(continued fro		3%	tal.	#33	T10	±10ml	£10	1		1
Time (24 hr)	Volume Removed (Liters)	Temp (°C)	рН	Specific Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
750		1190	634	4.090	2,06	132.6	7/000			
855	V	1490	6.33	11.084	1.95	1294	7/000	500		
150	100al	14.89	1.28	0.674	234	125.7	71000			
9.05	1	1489	6.23	0.065	2.77	12-3.1	71000		42.98	
915		14.89	6.12	0.054	3.48	119.8	7/000	-		
9.15		14.89	6.01	0.044	4.12	1129	7/1000			
920		14.92	5.79	0.032	6.08	1185	47 8			drain Slowcell
1.25	1	1493	5.73	0.030	G. Z.	119. 2	63Z			to remove si
1:30		1492	5,6%	0.827	6.47	123,1	493			
135		1992	5.63	6.027	6.49	124.7				
941)	15aul	14.92	5.62	0.626	654	1263	449		42.62	
9:45		1496	5.61	0.025	550	128.7				
950		14.93	5.61	0.026	6.52	129.7	462	500		
455		14.93	5.60	0.025	5,53	131.2	416	8		
1005		14.95	5.61	0.025	6.53	1348	376	500		
1010		14.96	556	0.025	6.55	136.8	354			*
1015		1493	5.60	0.025	4.56	137.2	358			
1020	17.504	14.93	5.58	0.025	5.50	139.2	-377			5
1020										sample
38Y										
									-	

Section 4 Analytical Data Validation



DATA VALIDATION REPORT

Project:	Regional Groundwater Investigation — NWIRP Bethpage						
Laboratory:	Katahdin Analytical						
Sample Delivery Group:	SI8613						
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: 12/5/2016						
Reviewed by:	Tina Clemmey/Resolution Consultants File Name: SI8613_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 27 to 28 October 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
RE107D1-WG-102715	Groundwater	8260C/8270D_SIM
RE107D2-WG-102715	Groundwater	8260C/8270D_SIM
RE107D3-WG-102815	Groundwater	8260C/8270D_SIM
TB-102815	Trip Blank	8260C
DUPLICATE-WG-102715	Field Duplicate of RE107D2-WG-102715	8260C/8270D_SIM

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* (United States Environmental Protection Agency [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 Contract Laboratory Program National Functional Guidelines for Superfund Organic Methods Data Review (U.S. EPA, 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
- X Initial calibration /continuing calibration verification
- ✓ Laboratory blanks/trip blanks/field blanks
- **X** Surrogate spike recoveries
- **X** Matrix spike and/or matrix spike duplicate results
- ✓ Laboratory control sample/laboratory control sample duplicate results
- ✓ Field duplicates
- ✓ Internal standards
- ✓ Sample results/reporting issues

The symbol (\checkmark) indicates that no validation qualifiers were applied based on this parameter. 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 (ICAL) 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 (CCV) method percent difference or percent drift and response factor acceptance criteria were met
- The retention time method acceptance criteria were met

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

ICAL Linearity Non-conformance:

Critoria	Actions				
Criteria	Detected Results	Non-detected Results			
%RSD >15% and quantitation based on mean response factor	J	W			

Notes:

%RSD = Relative standard deviation

J = Estimated

UJ = Undetected and estimated

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

ICV Recovery Non-conformance:

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

Notes:

J = Estimated

UJ = Undetected and estimated

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

CCV Linearity Non-conformance:

Γ	Criteria	Actions				
	Criteria	Detected Results	Non-detected Results			
Γ	%Difference or %Drift > 20%	ı	111			

Notes:

J = Estimated

UJ = Undetected and estimated



ICAL, ICV and CCV non-conformances are summarized in Attachment A in Tables A-1, A-2, and A-3.

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		
Criteria	Detected	Non-detected	
% R > Upper Limit	J	No qualification	
20% < %R < Lower Limit	J	UJ	
% R < 20%	J	Rejected	

Notes:

%R = Percent recovery
J = Estimated

UJ = Undetected and estimated

Surrogate recovery non-conformance is summarized in Attachment A in Table A-4.

Matrix Spike/Matrix Spike Duplicate (MS/MSD) Results

MS/MSDs are generated to provide information about the effect of each sample matrix on the sample preparation and the measurement methodology. MS/MSD percent recoveries (%Rs) assess the effect of the sample matrix on the accuracy of the analytical results and %Rs above the laboratory control limit could indicate a potential high result bias while %Rs below QC limits could indicate a potential low result bias. The relative percent differences (RPDs) between the MS and MSD results are evaluated to assess sample precision. The MS/MSD %Rs and RPDs were reviewed for conformance with the QC acceptance criteria. Data qualification to the analytes associated with the specific MS/MSD non-conformances were as follows:

MS/MSD Non-conformances Chart:

Critorio	Action				
Criteria	Detected Compounds	Non-detected Compounds			
%R>Upper Limit	J	No qualification			
20% < %R < Lower Limit	J	UJ			
%R <20%	J	Rejected			



%R = Percent recovery

RPD = Relative percent difference

J = Estimated

UJ = Undetected and estimated

MS/MSD non-conformances are summarized in Attachment A in Table A-5.

Qualifications Actions

The data were 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. EPA 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 Non-Conformance							
Method								
8260C	CHLOROMETHANE	15.39851	<15%	SI8613-1, -2, -3, -4, -5, -6, -7, -8, -9, -1RA, -2RA, -3RA, -4RA, 5RA, -6RA	Detects: J Non-detects: UJ			
8260C	METHELYENE CHLORIDE	15.35476	<15%	SI8613-1, -2, -3, -4, -5, -6, -7, -8, -9, -1RA, -2RA, -3RA, -4RA, 5RA, -6RA	Detects: J Non-detects: UJ			
8260C	ACETONE	15.98868	<15%	SI8613-1, -2, -3, -4, -5, -6, -7, -8, -9, -1RA, -2RA, -3RA, -4RA, 5RA, -6RA	Detects: J Non-detects: UJ			
8260C	4-METHYL-2-PENTANONE	15.14674	<15%	SI8613-1, -2, -3, -4, -5, -6, -7, -8, -9, -1RA, -2RA, -3RA, -4RA, 5RA, -6RA	Detects: J Non-detects: UJ			
8260C	TRICHLOROFLUOROMETHANE	15.30044	<15%	SI8613-7RA, -8RA	Detects: J Non-detects: UJ			
8260C	ACETONE	15.32971	<15%	SI8613-7RA, -8RA	Detects: J Non-detects: UJ			

%RSD Relative standard deviation UJ = Non-detect estimated value

Estimated value

	Table A-2 Initial Calibration Verification Non-Conformance						
Method	Method Analytes ICVID %R Limit Associated Samples Qualifier						
8260C	ACETONE	W4194.D	71.82	80-120	SI8613-1, -2, -3, -4, -5, -6, -7, -8, -9, -1RA, -2RA, -3RA, - 4RA, 5RA, -6RA	Detects: J Non-detects: UJ	
8260C	DICHLORODIFLUOROMETHANE	W4559A.D	55.29	80-120	SI8613-7RA, -8RA	Detects: J Non-detects: UJ	

Notes:

ICV ID Initial calibration verification identification

ID = Identification %R

Percent recovery Non-detect estimated value UJ

Estimated value

Table A-3 Continuing Calibration Verification Non-Conformance							
Lab ID /Calibration ID	Analytes	%D	%D Limit	Associated Samples	Qualifier		
WG173408-4 / W4297.D	BROMOMETHANE	24.83948	+/- 20	SI8613-1RA, -2RA, -3RA, -4RA, 5RA, - 6RA, -7RA, -8RA	Detects: J Non-detects: UJ		
WG173408-4 / W4297.D	CHLOROETHANE	30.94867	+/- 20	SI8613-1RA, -2RA, -3RA, -4RA, 5RA, - 6RA, -7RA, -8RA	Detects: J Non-detects: UJ		
WG173296-4 / W4248.D	ACETONE	-20.22537	+/- 20	SI8613-1, -2, -3, -4, -5, -6, -7, -8, -9	Detects: J Non-detects: UJ		
WG173296-4 / W4248.D	CHLOROMETHANE	-25.72419	+/- 20	SI8613-1, -2, -3, -4, -5, -6, -7, -8, -9	Detects: J Non-detects: UJ		

Notes:

ID Identification %D Percent difference

UJ Non-detect estimated value Detected estimated value

	Table A-4 Surrogate Non-Conformance							
Method	Surrogate	%R	Limits	Associated Sample	Qualifier			
8260C	1,2-DICHLOROETHANE-D4	138	70-120	DUPLICATE-WG-102715	Detects: J			
8260C	DIBROMOFLUOROMETHANE	124	85-115	DUPLICATE-WG-102715	Detects: J			
8260C	1,2-DICHLOROETHANE-D4	138	70-120	RE107D2-WG-102715	Detects: J			
8260C	DIBROMOFLUOROMETHANE	123	85-115	RE107D2-WG-102715	Detects: J			
8260C	1,2-DICHLOROETHANE-D4	140	70-120	RE107D3-WG-102815	Detects: J			
8260C	DIBROMOFLUOROMETHANE	125	85-115	RE107D3-WG-102815	Detects: J			
8260C	1,2-DICHLOROETHANE-D4	137	70-120	RE107DI-WG-102715	Detects: J			
8260C	DIBROMOFLUOROMETHANE	121	85-115	RE107DI-WG-102715	Detects: J			

%R Percent recovery Detected estimated value =

Table A-5 Matrix Spike/Matrix Spike Duplicate Non-Conformance (Micrograms per liter)							
Spiked Sample	Analytes	Sample Result	Spike Added	MS %R	MSD %R	%R Limits	Qualifier
RE107DI-WG-102715	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	0.82	50.0	129	123	73-126	J
RE107DI-WG-102715	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	0.74	50.0	129	123	73-126	J

Notes:

MS

MSD %R

Bold

Matrix spike
Matrix spike duplicate
Percent recovery
Percent recovery not within control limit
Detected analytes in associated sample qualified estimated "J" because %R is lower than the control limit.

Attachment B

Qualifier Codes and Explanations

Qualifier	Explanation
J	The analytes was positively identified; the associated numerical value is the approximate concentration of the analytes in the sample.
UJ	The analytes 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 analytes in the sample.
U	The analytes 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
bm	Missing blank information
bt	Trip blank contamination
С	Calibration issue
cr	Chromatographic resolution
d	Reporting limit raised due to chromatographic interference
dt	Dissolved result > total over limit
е	Ether interference
ej	Above calibration range; result estimated.
f	Presumed contamination from FB or ER.
fd	Field duplicate RPDs
h	Holding times
hs	Headspace greater than 6mm in all sample vials
i	Internal standard areas
ii	Injection internal standard area or retention time exceedance
it	Instrument tune
k	Estimated maximum possible concentrations (EMPC)
l	LCS recoveries
Ic	Labeled compound recovery
ld	Laboratory duplicate RPDs
lp	Laboratory control sample/laboratory control sample duplicate RPDs
m	Matrix spike recovery
mc	Deviation from the method
md	MS/MSD RPDs
nb	Negative laboratory blank contamination
р	Chemical preservation issue
p-h	Uncertainty near detection limit (< Reporting Limit), historical reason code applied.
pe	Post Extraction Spike
q	Quantitation issue
r	Dual column RPD
rt	SIM ions not within + 2 seconds
S	Surrogate recovery
sp	Sample preparation issue
SU	Evidence of ion suppression
t	Temperature Preservation Issue
X	Low % solids
У	Serial dilution results
Z	ICS results

Attachment D Final Results after Data Review

		Sample Deliv	ery Group Lab ID		18613 8613-1	
			Sample ID mple Date		I-WG-10271 27/2015	5
			mple Type		undwater	
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	0.82	J	s,m
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	U	
8260C	2-HEXANONE	591-78-6	UG_L	2.5	U	
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	UJ	С
8260C	ACETONE	67-64-1	UG_L	2.5	UJ	С
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	U	
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	U	
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	С
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	U	
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG_L	0.5	U	
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	1	U	
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	UJ	С
8260C	O-XYLENE	95-47-6	UG_L	4.4	J	S
8260C	STYRENE	100-42-5	UG_L	0.5	U	
8260C	TETRACHLOROETHENE	127-18-4	UG_L	1.3	J	S
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	14	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	4.4	J	S
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	8.2		

UG_L = Micrograms per liter NA

		Sample Deliv	ery Group Lab ID		518613 8613-2	
			Sample ID		2-WG-10271	5
			mple Date		27/2015	
<u> </u>	T		mple Type		undwater	
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 14	U J	
8260C 8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L UG L	0.5	U	S
8260C	1,1,2-TRICHLOROETHANE 1,1-DICHLOROETHANE	79-00-5 75-34-3	UG_L UG L	0.5	U	
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L 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-DIBROMO-3-CHEOKOFKOFANE	106-93-4	UG_L	0.75	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	2.2	J	S
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	U	3
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	U	
8260C	2-HEXANONE	591-78-6	UG L	2.5	U	
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	UJ	С
8260C	ACETONE	67-64-1	UG L	2.5	UJ	С
8260C	BENZENE	71-43-2	UG L	0.5	U	
8260C	BROMODICHLOROMETHANE	75-27-4	UG L	0.5	Ü	
8260C	BROMOFORM	75-25-2	UG L	0.5	U	
8260C	BROMOMETHANE	74-83-9	UG_L	1	Ü	
8260C	CARBON DISULFIDE	75-15-0	UG L	0.5	Ü	
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	0.82	J	S,C
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	2.2	J	S
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG_L	0.5	U	
8260C	CYCLOHEXANE	110-82-7	UG_L	0.5	U	
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG_L	0.5	U	
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	1	U	
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	UJ	С
8260C	O-XYLENE	95-47-6	UG_L	4.3	J	S
8260C	STYRENE	100-42-5	UG_L	0.5	U	
8260C	TETRACHLOROETHENE	127-18-4	UG_L	7.3	J	S
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	140	J	S
8260C	TRICHLOROFLUOROMETHANE	75-69-4	UG_L	1	U	
8260C	VINYL CHLORIDE	75-01-4	UG_L	1 1 2	U	
8260C	XYLENES, TOTAL	1330-20-7	UG_L	4.3	J	S
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	10		

UG_L = Micrograms per liter NA

	Sample Delivery Group SI8613					
		Sample Deliv	Lab ID			
					8613-3	745
			Sample ID		ΓE-WG-102	/15
			mple Date		27/2015	
			mple Type		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	15	J	S
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	2.2	J	S
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	U	
8260C	2-HEXANONE	591-78-6	UG_L	2.5	U	
8260C	4-METHYL-2-PENTANONE	108-10-1	UG L	2.5	UJ	С
8260C	ACETONE	67-64-1	UG L	2.5	UJ	С
8260C	BENZENE	71-43-2	UG_L	0.5	U	
8260C	BROMODICHLOROMETHANE	75-27-4	UG L	0.5	Ü	
8260C	BROMOFORM	75-25-2	UG L	0.5	Ü	
8260C	BROMOMETHANE	74-83-9	UG L	1	Ü	
8260C	CARBON DISULFIDE	75-15-0	UG L	0.5	Ü	
8260C	CARBON TETRACHLORIDE	56-23-5	UG L	0.5	Ü	
8260C	CHLOROBENZENE	108-90-7	UG L	0.5	U	
8260C	CHLOROETHANE	75-00-3	UG L	1	Ü	
8260C	CHLOROFORM	67-66-3	UG L	0.5	Ü	
8260C	CHLOROMETHANE	74-87-3	UG L	1	UJ	С
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	2.2	J	S
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG L	0.5	U	
8260C	CYCLOHEXANE	110-82-7	UG L	0.5	Ü	
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG L	0.5	Ü	
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG L	1	Ü	
8260C	ETHYLBENZENE	100-41-4	UG L	0.5	Ü	
8260C	ISOPROPYLBENZENE	98-82-8	UG L	0.5	Ü	
8260C	M- AND P-XYLENE	108-38-3/106-42	UG L	1	Ü	
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	UJ	С
8260C	O-XYLENE	95-47-6	UG L	4.3	J	S
8260C	STYRENE	100-42-5	UG_L	0.5	U	
8260C	TETRACHLOROETHENE	127-18-4	UG_L	6.9	J	S
8260C	TOLUENE	108-88-3	UG_L	0.5	U	3
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	150	J	S
8260C	TRICHLOROFLUOROMETHANE	75-69-4	UG_L	1	U	3
8260C	VINYL CHLORIDE	75-09-4	UG_L	<u> </u>	U	
8260C 8260C	XYLENES, TOTAL	1330-20-7	UG_L	4.3	J	S
	1,4-DIOXANE				J	3
UZ / UD_SIIV	I I,T-DIOAANL	123-91-1	UG_L	12		

UG_L = Micrograms per liter NA

		Sample Deliv	ery Group	S	518613	
			Lab ID	SI	8613-8	
			Sample ID	RE107D	3-WG-1028	15
		Sai	mple Date	10/	28/2015	
			mple Type	Gro	undwater	
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	1.9	J	S
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG L	0.5	U	3
8260C	1.1-DICHLOROETHANE	75-34-3	UG_L	0.5	U	
		+		0.5	U	
8260C 8260C	1,1-DICHLOROETHENE 1,2,4-TRICHLOROBENZENE	75-35-4 120-82-1	UG_L 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	U	
8260C	2-HEXANONE	591-78-6	UG_L	2.5	U	
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	UJ	С
8260C	ACETONE	67-64-1	UG_L	2.5	UJ	С
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	U	
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	U	
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	С
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	U	
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG L	0.5	U	
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	1	Ü	
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	0.75	U	
00/00	METING TEST SUTIO ETUES	1634-04-4	UG_L			
8260C 8260C	METHYL TERT-BUTYL ETHER METHYLENE CHLORIDE	75-09-2	UG_L	0.5 2.5	UJ	С
8260C	O-XYLENE	95-47-6	UG_L	0.5	U	· ·
8260C 8260C	STYRENE STYRENE	100-42-5	UG_L	0.5	U	
8260C	TETRACHLOROETHENE	127-18-4	UG_L UG L	0.5	U	
8260C 8260C						
	TOLUENE TRANS 1.2 DICHLODOETHENE	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 1 5	U	
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	U	
8270D_SIN	1,4-DIOXANE	123-91-1	UG_L	0.18	U	

UG_L = Micrograms per liter NA

	Sample Delivery Group Lab I D Sample I D Sample Date			S TB	SI8613 SI8613-9 TB-102815 10/28/2015		
			mple Type		ip 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	Ü		
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	U		
8260C	2-HEXANONE	591-78-6	UG_L	2.5	U		
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	UJ	С	
8260C	ACETONE	67-64-1	UG_L	2.5	UJ	С	
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	U		
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	U		
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	С	
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	U		
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG L	0.5	U		
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG L	1	U		
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	UJ	С	
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	U		
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		
	1,4-DIOXANE	123-91-1	UG_L	NA			

UG_L = Micrograms per liter NA



DATA VALIDATION REPORT

Project:	Regional Groundwater Investigation — NWIRP Bethpage		
Laboratory:	Katahdin Analytical		
Sample Delivery Groups:	SI6392		
Analyses/Method:		by U.S. EPA SW-846 Method 9060A and Standard anic Carbon by High-Temperature Combustion	
Validation Level:	3		
Project Number:	0888812477.SA.DV		
Prepared by:	Dana Miller/Resolution Consultants	Completed on: 10/25/2015	
Reviewed by:	Tina Clemmey/Resolution Consultants	File Name: SI6392_ 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 20 August 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
RE107D3-SOIL-082015-647-649	SI6392-1	Soil	9060A
RE107D3-EB-082015	SI6392-2	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
- ✓ 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. 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. Attachment A, Table A-1 provides final results after data review.

ATTACHMENTS

Attachment A: Table A-1 Final Results after Data Review

Attachment A Final Results after Data Review

Table A-1
Final Results after Data Review
Regional Groundwater Investigation NWIRP Bethpage

	Sample Delivery Group			SI6392	SI6392
	Lab ID			SI6392-1	SI6392-2
	Sample ID F			RE107D3-SOIL-082015-647-649	RE107D3-EB-082015
	Sample Date			8/20/2015	8/20/2015
			Sample Type	Soil	Equipment Blank
Method	Analyte	CAS No	Units	Result	Result
2540G	TOTAL SOLIDS	-29	PCT	84	NA
5310B	TOTAL ORGANIC CARBON	-28	MG_L	NA	0.12
9060A	TOTAL ORGANIC CARBON	-28	UG_G	940	NA

ID = Identification PCT = Percent

 $MG_L = Milligrams per liter$ $UG_G = Micrograms per gram$

NA = Not analyzed



DATA VALIDATION REPORT

Regional Groundwater Investigation — NWIRP Bethpage		
Katahdin Analytical		
SI6938		
	by U.S. EPA SW-846 Method 9060A and Standard anic Carbon by High-Temperature Combustion	
3		
0888812477.SA.DV		
Dana Miller/Resolution Consultants	Completed on: 10/8/2015	
Tina Clemmey/Resolution Consultants	File Name: SI6938_9060A_5310B	
	Katahdin Analytical SI6938 Total Organic Carbon (TOC) Method 5310B for Total Orga 3 0888812477.SA.DV Dana Miller/Resolution Consultants Tina Clemmey/Resolution	

SUMMARY

This report summarizes data review findings for samples listed below, collected by Resolution Consultants from the Regional Groundwater Investigation — NWIRP Bethpage site on 8 September 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	
RE107D2-SOIL-090815-563-565	SI6938-1	Soil	9060A	
RE107D2-EB-090815	SI6938-2	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
- ✓ 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 were 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. Attachment A, Table A-1 provides final results after data review.



ATTACHMENTS

Attachment A: Table A-1 Final Results after Data Review

Attachment A Final Results after Data Review

Table A-1
Final Results after Data Review
Regional Groundwater Investigation NWIRP Bethpage

Sample Delivery Group			SI6	938	SI6	938	
Lab ID			SI6938-1		SI6938-2		
Sample ID			RE107D2-SOIL-090815-563-565		RE107D2-EB-090815		
Sample Date		9/8/2015		9/8/2015			
	Sample Type		Soil		Equipment Blank		
Method	Analyte	CAS No	Units	Result	Qual	Result	Qual
2540G	TOTAL SOLIDS	-29	PCT	85		NA	
5310B	TOTAL ORGANIC CARBON	-28	MG_L	NA		0.29	
9060A	TOTAL ORGANIC CARBON	-28	UG_G	300	J	NA	

ID = Identification PCT = Percent

MG_L = Milligrams per liter
UG_G = Micrograms per gram
Qual = Final interpreted qualifier

NA = Not analyzed

J = Estimated value – Value was below the limit of quantitation.



DATA VALIDATION REPORT

Regional Groundwater Investigation — NWIRP Bethpage				
Katahdin Analytical				
SI7430				
Total Organic Carbon (TOC) by U.S. EPA SW-846 Method 9060A and Standard Method 5310B for Total Organic Carbon by High-Temperature Combustion				
3				
0888812477.SA.DV				
Dana Miller/Resolution Consultants	Completed on: 10/25/2015			
Tina Clemmey/Resolution Consultants	File Name: SI7430_ 9060A_5310B			
	Katahdin Analytical SI7430 Total Organic Carbon (TOC) Method 5310B for Total Orga 3 0888812477.SA.DV Dana Miller/Resolution Consultants Tina Clemmey/Resolution			

SUMMARY

This report summarizes data review findings for samples listed below, collected by Resolution Consultants from the Regional Groundwater Investigation — NWIRP Bethpage site on 22 September 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	
RE107D1-SOIL-515-517-092215	SI7430-1	Soil	9060A	
RE107D1-EB-515-517-092215	SI7430-2	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
- ✓ 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. 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. Attachment A, Table A-1 provides final results after data review.

ATTACHMENTS

Attachment A: Table A-1 Final Results after Data Review

Attachment A Final Results after Data Review

Table A-1 Final Results after Data Review Regional Groundwater Investigation NWIRP Bethpage

		SI7430	SI7430		
	Lab ID			SI7430-1	SI7430-2
				RE107D1-SOIL-515-517-	RE107D1-EB-515-517-
			Sample ID	092215	092215
			Sample Date	9/22/2015	9/22/2015
			Sample Type	Soil	Equipment Blank
Method	Analyte	CAS No	Units	Result	Result
2540G	TOTAL SOLIDS	-29	PCT	83	NA
5310B	TOTAL ORGANIC CARBON	-28	MG_L	NA	0.24
9060A	TOTAL ORGANIC CARBON	-28	UG_G	720	NA

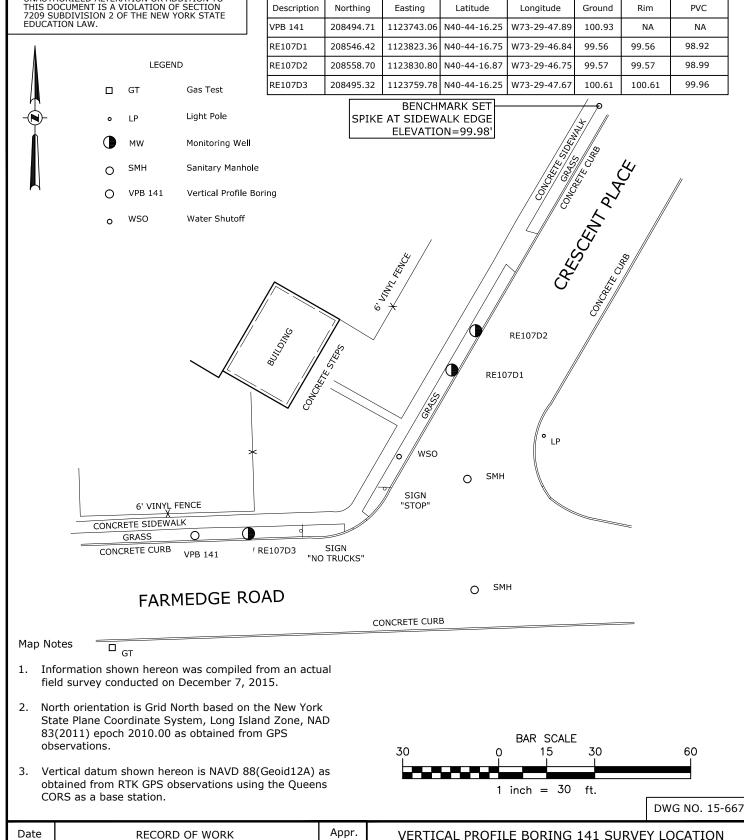
ID Identification PCT = Percent

MG_L = UG_G = NA = Milligrams per liter Micrograms per gram Not analyzed

Section 5

Survey

UNAUTHORIZED ALTERATION OR ADDITION TO



Drafter: LMK Checker: JFC

Appr. by: JFC Proj. No. 14.4121

VERTICAL PROFILE BORING 141 SURVEY LOCATION RE107D1-RE107D2-RE107D3 FARMEDGE ROAD AND CRESCENT 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: D

DATE: DECEMBER 7, 2015

Appendix B.

RE107D1, RE107D2, RE107D3

December 2015 Analytical Data Summary

[Attachment D of the Data Validation report for the December 2015 Quarterly Sampling included here contains only results tables for RE107D1, RE107D2 and RE107D3; for the December 2015 Quarterly Sampling Data Validation report with complete Attachment D with all well results tables see *December 2015 Groundwater Sampling Data Summary Report, Bethpage, NY*, Resolution Consultants, 2016.]

Location	NVCDEC	RE107D1	RE107D2	RE107D3
Sample Date	NYSDEC Groundwater	12/18/2015	12/18/2015	12/29/2015
Sample ID	Guidance or Standard Value	RE107D1-GW-121815	RE107D2-GW-121815	RE107D3-GW-122915
Sample type code	(Note 1)	N	N	N
VOC 8260C (ug/L)				
1,1,1-TRICHLOROETHANE	5	< 0.50 U	< 0.50 U	< 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	0.95 J	15	4.9
1,1,2-TRICHLOROETHANE	1	< 0.50 U	< 0.50 U	< 0.50 U
1,1-DICHLOROETHANE	5	< 0.50 U	< 0.50 U	< 0.50 U
1,1-DICHLOROETHENE	5	< 0.50 U	< 0.50 U	< 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	0.21 J	2.7	< 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 NL	6.9	9.3	< 0.17 U
2-BUTANONE	50	< 2.5 U	< 2.5 UJ	< 2.5 U
2-HEXANONE	50	< 2.5 U	< 2.5 UJ	< 2.5 U
4-METHYL-2-PENTANONE	NL NL	< 2.5 U	< 2.5 U	< 2.5 U
ACETONE	50	1		< 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 U	< 1.0 U	< 1.0 U
CARBON DISULFIDE	60	< 0.50 U	< 0.50 U	< 0.50 U
CARBON TETRACHLORIDE	5	< 0.50 U	< 0.50 U	< 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	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROMETHANE	5	< 1.0 UJ	< 1.0 U	< 1.0 U
CIS-1,2-DICHLOROETHENE	5	0.21 J	2.7	< 0.50 U
CIS-1,3-DICHLOROPROPENE	0.4	< 0.50 U	< 0.50 U	< 0.50 U
CYCLOHEXANE	NL	< 0.50 U	< 0.50 U	< 0.50 U
DIBROMOCHLOROMETHANE	5	< 0.50 U	< 0.50 U	< 0.50 U
DICHLORODIFLUOROMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U
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 UJ	< 0.75 U	< 0.75 U
METHYL CYCLOHEXANE	NL	< 0.50 U	< 0.50 U	< 0.50 U
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	1.6 J	6.4	< 0.50 UJ
TOLUENE	5	< 0.50 U	< 0.50 U	< 0.50 U
TRANS-1,2-DICHLOROETHENE	5	< 0.50 U	< 0.50 U	< 0.50 U
TRANS-1,3-DICHLOROPROPENE	0.4	< 0.50 UJ	< 0.50 UJ	< 0.50 U
TRICHLOROETHENE	5	17	140	0.36 J
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

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.



DATA VALIDATION REPORT

Project:	Regional Groundwater Investigation — NWIRP Bethpage		
Laboratory:	Katahdin Analytical		
Sample Delivery Group:	BETHPAGE-3		
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: 02/5/2016		
Reviewed by:	Tina Clemmey/Resolution Consultants File Name: BETHPAGE 3_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 14 to 29 December 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
FIELD1-FB-121615	Field Blank	8260C/8270D_SIM
RE103D1-GW-121415	Groundwater	8260C/8270D_SIM
RE103D2-GW-121415	Groundwater	8260C/8270D_SIM
RE103D3-GW-121415	Groundwater	8260C/8270D_SIM
RE104D1-GW-121515	Groundwater	8260C/8270D_SIM
RE104D2-GW-121515	Groundwater	8260C/8270D_SIM
DUPLICATE1-GW-121515	Field Duplicate of RE104D2-GW-121515	8260C/8270D_SIM
RE104D3-GW-121515	Groundwater	8260C/8270D_SIM
RE105D1-GW-121715	Groundwater	8260C/8270D_SIM
RE105D2-GW-121715	Groundwater	8260C/8270D_SIM
RE107D1-GW-121815	Groundwater	8260C/8270D_SIM



Sample ID	Matrix/Sample Type	Analysis
RE107D2-GW-121815	Groundwater	8260C/8270D_SIM
RE107D3-GW-122915	Groundwater	8260C/8270D_SIM
RE108D1-GW-122215	Groundwater	8260C/8270D_SIM
RE108D2-GW-122215	Groundwater	8260C/8270D_SIM
RE114D1-GW-122115	Groundwater	8260C/8270D_SIM
RE114D2-GW-121615	Groundwater	8260C/8270D_SIM
RE114D3-GW-121615	Groundwater	8260C/8270D_SIM
RE120D1-GW-121815	Groundwater	8260C/8270D_SIM
RE120D2-GW-122915	Groundwater	8260C/8270D_SIM
RE120D3-GW-122915	Groundwater	8260C/8270D_SIM
RE121D1-GW-122115	Groundwater	8260C/8270D_SIM
RE121D2-GW-122115	Groundwater	8260C/8270D_SIM
RE122D1-GW-121515	Groundwater	8260C/8270D_SIM
RE122D2-GW-121515	Groundwater	8260C/8270D_SIM
RE122D3-GW-121515	Groundwater	8260C/8270D_SIM
RE123D1-GW-122115	Groundwater	8260C/8270D_SIM
RE123D2-GW-122115	Groundwater	8260C/8270D_SIM
RE123D3-GW-122115	Groundwater	8260C/8270D_SIM
TRIP BLANK 121415	Trip Blank	8260C
TRIP BLANK 121615	Trip Blank	8260C
TRIP BLANK-121815	Trip Blank	8260C
TRIP BLANK-122915	Trip Blank	8260C
TT101D1-GW-121715	Groundwater	8260C/8270D_SIM
TT101D2-GW-122115	Groundwater	8260C/8270D_SIM
DUPLICATE-GW-122115	Field Duplicate of TT101D2-GW-122115	8260C/8270D_SIM
TT101D-GW-121715	Groundwater	8260C/8270D_SIM

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* (United States Environmental Protection Agency [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 Contract Laboratory Program National Functional Guidelines for Superfund Organic Methods Data Review* (U.S. EPA, 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 verification (ICV)/continuing calibration verification (CCV)
- ✓ Laboratory blanks/trip blanks/field blanks
- **X** Surrogate spike recoveries
- X Matrix spike and/or matrix spike duplicate results
- ✓ Laboratory control sample/laboratory control sample duplicate results
- X Field duplicates
- ✓ Internal standards
- ✓ Sample results/reporting issues

The symbol (\checkmark) indicates that no validation qualifiers were applied based on this parameter. 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 (\checkmark) 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 ICV standard percent recovery acceptance criteria were met
- The CCV method percent difference or percent drift and response factor acceptance criteria were met
- The retention time method acceptance criteria were met



Data qualification to the analytes associated with the specific initial calibration (ICAL) was as follows:

ICAL Linearity Non-conformance:

Cultonia	Actions		
Criteria	Detected Results	Non-detected Results	
%RSD >15% and quantitation based on mean response factor	J	m	

Notes:

%RSD = Relative standard deviation

J = Estimated

UJ = Undetected and estimated

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

ICV Recovery Non-conformance:

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

Notes:

J = Estimated

UJ = Undetected and estimated

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

CCV Linearity Non-conformance:

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

Notes:

= Estimated

UJ = Undetected and estimated

ICAL, ICV and CCV non-conformances are summarized in Attachment A in Tables A-1, A-2, and A-3.

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	Acti	on
Criteria	Detected	Non-detected
% R > Upper Limit	J	No qualification
20% < %R < Lower Limit	J	UJ
% R < 20%	J	Rejected

Notes:

%R = Percent recovery J = Estimated

UJ = Undetected and estimated

Surrogate recovery non-conformance is summarized in Attachment A in Table A-4.

Matrix Spike/Matrix Spike Duplicate (MS/MSD) Results

MS/MSDs are generated to provide information about the effect of each sample matrix on the sample preparation and the measurement methodology. MS/MSD percent recoveries (%Rs) assess the effect of the sample matrix on the accuracy of the analytical results and %Rs above the laboratory control limit could indicate a potential high result bias while %Rs below QC limits could indicate a potential low result bias. The relative percent differences (RPDs) between the MS and MSD results are evaluated to assess sample precision. The MS/MSD %Rs and RPDs were reviewed for conformance with the QC acceptance criteria. Data qualification to the analytes associated with the specific MS/MSD non-conformances were as follows:

MS/MSD Non-conformances Chart:

Criteria	Action		
Criteria	Detected Compounds	Non-detected Compounds	
%R>Upper Limit	J	No qualification	
20% < %R < Lower Limit	J	UJ	
%R <20%	J	Rejected	

Notes:

%R = Percent recovery

RPD = Relative percent difference

J = Estimated

UJ = Undetected and estimated

MS/MSD non-conformances are summarized in Attachment A in Table A-5.



Field Duplicate

Two field duplicate pairs were collected to assess precision: RE104D2-GW-121515/ DUPLICATE1-GW-121515 and TT101D2-GW-122115/DUPLICATE-GW-122115. Field duplicate RPDs were reviewed for conformance with the Resolution Consultants QC criteria of \leq 30% for aqueous matrices and \leq 50% for solid matrices. These criteria apply if both results were greater than two times the limit of quantitation (LOQ). Data qualification to the analytes associated with the specific field duplicate RPDs was as follows:

Field Duplicate Non-conformances Chart:

Criteria	RPD	Action		
Criteria	INI D	Detected	Non-detected	
Sample and duplicate are nondetect	Not calculable (NC)	No qualification	No qualification	
Sample and duplicate results ≥2x LOQ	>30 (aqueous)	ı	Not Applicable	
Sample and duplicate results >2x LOQ	>50 (solids)	J	Not Applicable	
If sample or duplicate result is >2x LOQ and the other is not detected	NC	J	ΟΊ	
If sample or duplicate result is <2x LOQ and the other is not detected	NC	No qualification	No qualification	

Notes:

LOQ = Limit of quantitation

J = Estimated

UJ = Undetected and estimated

Field duplicate non-conformances are summarized in Attachment A in Table A-6.

Qualifications Actions

The data were 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. Trichloroethene in sample RE108D2-GW-121215 result value reported above the calibration range and was qualified estimated "J" because the value was off-scale.

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. EPA 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 Non-Conformance						
Method	Method Analyte %RSD Limit Associated Samples Qualifier						
■ 8260C CHLOROETHANE 17 909/1 >15%				Detects: J Non-detects: UJ			
8260C	TETRACHLOROETHENE	15.13611	<15%	TI0428-4, -1, -3RA, -1DL, -2RA, and TI0330- 17DL	Detects: J Non-detects: UJ		

%RSD = Relative standard deviation UJ = Non-detect estimated value

J = Estimated value

	Table A-2 Initial Calibration Verification Non-Conformance								
Method	Analyte	ICVID	%R	Limit	Associated Samples	Qualifier			
8260C	TETRACHLOROETHENE	P3840.D	150.8	80-120	TI0116-1, -2, -3, -4, -5, -6, -7, - 8, -9, -11, -10RA, -2DL, -4DL, - 9DL, -3DL, -11DL, TI0214-1, -4, -5, -6, -7, -9, -8, -10, 12, -8DL, and - 7DL	Detects: J Non-detects: UJ			
8260C	TRANS-1,3-DICHLOROPROPENE	C6396A.D	123.47	80-120	TI0330-1, -2, -5, -6, -11, -4, -7, - 3DL,-10RA, -12, -13, -14, -17, - 4DL, -6DL, and 7DL	Detects: J Non-detects: UJ			
8260C	ACETONE	P4114A.D	72.62	80-120	TI0428-3RA, -1DL, -2RA, -4, -1, and TI0330-17DL	Detects: J Non-detects: UJ			

Initial calibration verification identification Identification Percent recovery Non-detect estimated value Estimated value

Notes: ICV ID ID %R UJ J

=

	Table A-3 Continuing Calibration Verification Non-Conformance								
Lab ID /Calibration ID	Analyte	%D	%D Limit	Associated Samples	Qualifier				
WG176285-4 / P3940.D	BROMOMETHANE	23.79988	+/- 20	TI0116-1, -2, -3, -4, -5, -6, -7, -8, -9, and -11	Detects: J Non-detects: UJ				
WG176319-4 / P3964.D	BROMOMETHANE	24.57249	+/- 20	TI0116-10RA, TI0214-1, -4, -5, -6, -7, and -9	Detects: J Non-detects: UJ				
WG176319-4 / P3964.D	4-METHYL-2-PENTANONE	20.68251	+/- 20	TI0116-10RA, TI0214-1, -4, -5, -6, -7, and -9	Detects: J Non-detects: UJ				
WG176436-4 / P3989.D	BROMOMETHANE	28.03265	+/- 20	TI0214-8, -10, and -12	Detects: J Non-detects: UJ				
WG176436-4 / P3989.D	CHLOROETHANE	27.72207	+/- 20	TI0214-8, -10, and -12	Detects: J Non-detects: UJ				
WG176436-4 / P3989.D	4-METHYL-2-PENTANONE	21.50811	+/- 20	TI0214-8, -10, and -12	Detects: J Non-detects: UJ				
WG176832-4 / P4138.D	ACETONE	-34.27373	+/- 20	TI0428-4, and -1	Detects: J Non-detects: UJ				
WG176788-4 / C6418.D	CHLOROMETHANE	-20.54679	+/- 20	TI0330-10RA, -12, -13, -14, -17, -4DL, -6DL, -and -7DL	Detects: J Non-detects: UJ				
WG176788-4 / C6418.D	ACETONE	-27.76004	+/- 20	TI0330-10RA, -12, -13, -14, -17, -4DL, -6DL, -and -7DL	Detects: J Non-detects: UJ				
WG176788-4 / C6418.D	TETRACHLOROETHENE	-21.38037	+/- 20	TI0330-10RA, -12, -13, -14, -17, -4DL, -6DL, -and -7DL	Detects: J Non-detects: UJ				
WG176788-4 / C6418.D	METHYL ACETATE	-25.56044	+/- 20	TI0330-10RA, -12, -13, -14, -17, -4DL, -6DL, -and -7DL	Detects: J Non-detects: UJ				
WG176732-4 / C6394.D	ACETONE	57.49765	+/- 20	TI0330-1, -2, -5, -6, -11, -4, -7, and -3DL	Detects: J Non-detects: UJ				
WG176732-4 / C6394.D	2-BUTANONE	25.58556	+/- 20	TI0330-1, -2, -5, -6, -11, -4, -7, and -3DL	Detects: J Non-detects: UJ				
WG176732-4 / C6394.D	2-HEXANONE	22.15201	+/- 20	TI0330-1, -2, -5, -6, -11, -4, -7, and -3DL	Detects: J Non-detects: UJ				

ID = Identification %D = Percent difference

UJ = Non-detect estimated value
J = Detected estimated value

	Table A-4 Surrogate Non-Conformance							
Method Surrogate %R Limits Associated Sample Qualifier								
8260C	1,2-DICHLOROETHANE-D4	121	70-120	RE114D1-GW-122115	Detects: J			
8260C	DIBROMOFLUOROMETHANE	117	85-115	RE114D1-GW-122115	Detects: J			
8260C	DIBROMOFLUOROMETHANE	116	85-115	RE121D2-GW-122115	Detects: J			

=

Notes: %R UJ Percent recovery Non-detect estimated value J Detected estimated value

	Table						
	Matrix Spike/Matrix Spike Du (Micrograms		n-Conform	ance			
	(Which ograms	per inter)		1			
		Sample	Spike	MS	MSD	%R	
Spiked Sample	Analyte	Result	Added	%R	%R	Limits	Qualifier
TT101D2-GW-122115	METHYLENE CHLORIDE	<2.5	50.0	53.4	60.2	55-140	UJ
TT101D2-GW-122115	CIS-1,2-DICHLOROETHENE	1.7	50.0	54.6	59.8	70-125	J
TT101D2-GW-122115	1,2,4-TRICHLOROBENZENE	< 0.50	50.0	50.8	56.2	65-135	UJ
TT101D2-GW-122115	CHLOROBENZENE	< 0.50	50.0	55.8	60.2	80-120	UJ
TT101D2-GW-122115	1,1-DICHLOROETHANE	< 0.50	50.0	64.4	70.8	70-135	UJ
TT101D2-GW-122115	CIS-1,3-DICHLOROPROPENE	< 0.50	50.0	60.8	64.4	70-130	UJ
TT101D2-GW-122115	1,2-DIBROMO-3-CHLOROPROPANE	< 0.75	50.0	46.4	57	50-130	UJ
TT101D2-GW-122115	ISOPROPYLBENZENE	< 0.50	50.0	56.2	61	75-125	UJ
TT101D2-GW-122115	TRANS-1,2-DICHLOROETHENE	< 0.50	50.0	58.4	65.2	60-140	UJ
TT101D2-GW-122115	BENZENE	< 0.50	50.0	62.4	67.6	80-120	UJ
TT101D2-GW-122115	1,2-DICHLOROPROPANE	< 0.50	50.0	63.8	68.8	75-125	UJ
TT101D2-GW-122115	O-XYLENE	< 0.50	50.0	58.8	62.2	80-120	UJ
TT101D2-GW-122115	1,3-DICHLOROBENZENE	< 0.50	50.0	51.4	56.8	75-125	UJ
TT101D2-GW-122115	1,1-DICHLOROETHENE	3.6	50.0	56.8	63.4	70-130	J
TT101D2-GW-122115	1,1,2-TRICHLOROETHANE	0.50	50.0	63.4	65.4	75-125	J
TT101D2-GW-122115	CYCLOHEXANE	< 0.50	50.0	63.8	69	71-133	UJ
TT101D2-GW-122115	TOLUENE	< 0.50	50.0	65	69.6	75-120	UJ
TT101D2-GW-122115	CARBON TETRACHLORIDE	1.3	50.0	63.6	66.6	65-140	J
TT101D2-GW-122115	1,2-DICHLOROETHANE	< 0.50	50.0	61.6	66	70-130	UJ
TT101D2-GW-122115	1,2-DICHLOROETHENE, TOTAL	1.7	100	56.5	62.5	84-121	J
TT101D2-GW-122115	XYLENES, TOTAL	<1.5	150	59.8	63.2	89-116	UJ
TT101D2-GW-122115	STYRENE	< 0.50	50.0	60.4	63.4	65-135	UJ
TT101D2-GW-122115	DIBROMOCHLOROMETHANE	< 0.50	50.0	59.4	63.6	60-135	UJ
TT101D2-GW-122115	1,4-DICHLOROBENZENE	< 0.50	50.0	50.8	54.6	75-125	UJ
TT101D2-GW-122115	ETHYLBENZENE	< 0.50	50.0	56.2	60.8	75-125	UJ
TT101D2-GW-122115	M- AND P-XYLENE	<1.0	100	60.2	63.8	75-130	UJ
TT101D2-GW-122115	BROMOFORM	< 0.50	50.0	55.6	57	70-130	UJ
TT101D2-GW-122115	BROMODICHLOROMETHANE	< 0.50	50.0	67.6	70	75-120	UJ
TT101D2-GW-122115	CHLOROFORM	0.90	50.0	57	62.4	65-135	J
TT101D2-GW-122115	1,1,1-TRICHLOROETHANE	0.34	50.0	58.9	64.3	65-130	J
TT101D2-GW-122115	1,2-DIBROMOETHANE	< 0.50	50.0	61.2	66.6	80-120	UJ
TT101D2-GW-122115	1,1,2,2-TETRACHLOROETHANE	< 0.50	50.0	51.6	59.2	65-130	UJ
	•		•	1		•	

Matrix spike Matrix spike duplicate MS MSD %R Percent recovery

TT101D2-GW-122115

Bold Percent recovery not within control limit

Nondetect analyte in associated sample qualified estimated "UJ" because the %R is lower than the control limit. UJ J Detected analyte in associated sample qualified estimated "J" because %R is lower than the control limit.

< 0.50

50.0

52

57.6

70-120

UJ

1,2-DICHLOROBENZENE

Table A-6 Field Duplicate (Micrograms per liter)						
			Sample	Duplicate		
Sample ID	Duplicate ID	Analyte	Result	Result	RPD	Qualifiers
TT101D2-GW-122115	DUPLICATE-GW-122115	1,1-DICHLOROETHENE	3.6	5	32.6	J - both results

Notes: RPD J Relative percent difference Estimated value =

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
bm	Missing blank information
bt	Trip blank contamination
С	Calibration issue
cr	Chromatographic resolution
d	Reporting limit raised due to chromatographic interference
dt	Dissolved result > total over limit
е	Ether interference
ej	Above calibration range; result estimated.
f	Presumed contamination from FB or ER.
fd	Field duplicate RPDs
h	Holding times
hs	Headspace greater than 6mm in all sample vials
i	Internal standard areas
ii	Injection internal standard area or retention time exceedance
it	Instrument tune
k	Estimated maximum possible concentrations (EMPC)
l	LCS recoveries
Ic	Labeled compound recovery
ld	Laboratory duplicate RPDs
lp	Laboratory control sample/laboratory control sample duplicate RPDs
m	Matrix spike recovery
mc	Deviation from the method
md	MS/MSD RPDs
nb	Negative laboratory blank contamination
р	Chemical preservation issue
p-h	Uncertainty near detection limit (< Reporting Limit), historical reason code applied.
pe	Post Extraction Spike
q	Quantitation issue
r	Dual column RPD
rt	SIM ions not within + 2 seconds
S	Surrogate recovery
sp	Sample preparation issue
SU	Evidence of ion suppression
t	Temperature Preservation Issue
X	Low % solids
У	Serial dilution results
Z	ICS results

Attachment D Final Results after Data Review

		Sample	Delivery Group Lab ID Sample ID	TIO	THPAGE- 330-10RA 01-GW-12	١.
			Sample Date Sample Type	12	2/18/2015 oundwater	
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	0.95	J	
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	0.21	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	U	
8260C	2-HEXANONE	591-78-6	UG L	2.5	U	
8260C	4-METHYL-2-PENTANONE	108-10-1	UG L	2.5	U	
8260C	ACETONE	67-64-1	UG L	2.5	UJ	С
8260C	BENZENE	71-43-2	UG L	0.5	U	
8260C	BROMODICHLOROMETHANE	75-27-4	UG L	0.5	Ū	
8260C	BROMOFORM	75-25-2	UG L	0.5	Ū	
8260C	BROMOMETHANE	74-83-9	UG L	1	Ū	
8260C	CARBON DISULFIDE	75-15-0	UG L	0.5	Ū	
8260C	CARBON TETRACHLORIDE	56-23-5	UG L	0.5	Ū	
8260C	CHLOROBENZENE	108-90-7	UG L	0.5	Ū	
8260C	CHLOROETHANE	75-00-3	UG L	1	UJ	С
8260C	CHLOROFORM	67-66-3	UG L	0.5	U	
8260C	CHLOROMETHANE	74-87-3	UG L	1	UJ	С
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG L	0.21	J	
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG L	0.5	Ü	
8260C	CYCLOHEXANE	110-82-7	UG L	0.5	Ü	
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG L	0.5	Ū	
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG L	1	Ü	
8260C	ETHYLBENZENE	100-41-4	UG L	0.5	Ū	
8260C	ISOPROPYLBENZENE	98-82-8	UG L	0.5	Ü	
8260C	M- AND P-XYLENE	108-38-3/106-42	UG L	1	Ü	
8260C	METHYL ACETATE	79-20-9	UG L	0.75	UJ	С
8260C	METHYL CYCLOHEXANE	108-87-2	UG L	0.5	U	Ť
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	Ü	<u> </u>
8260C	METHYLENE CHLORIDE	75-09-2	UG L	2.5	Ü	<u> </u>
8260C	O-XYLENE	95-47-6	UG L	0.5	Ü	1
8260C	STYRENE	100-42-5	UG L	0.5	Ü	
8260C	TETRACHLOROETHENE	127-18-4	UG L	1.6	J	С
8260C	TOLUENE	108-88-3	UG L	0.5	Ŭ	Ť
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG L	0.5	Ü	1
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG L	0.5	UJ	С
8260C	TRICHLOROETHENE	79-01-6	UG_L	17		Ť
8260C	TRICHLOROFLUOROMETHANE	75-69-4	UG L	1	U	
8260C	VINYL CHLORIDE	75-01-4	UG L		Ü	
8260C	XYLENES, TOTAL	1330-20-7	UG L	1.5	Ü	
3270D_SIM	1,4-DIOXANE	123-91-1	UG L	6.9		1

		Sample	E Delivery Group Lab ID Sample ID Sample Date Sample Type	T RE107[12	THPAGE-3 10330-11 D2-GW-12 2/18/2015 oundwater	21815
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	Ü	-
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	79-34-3 76-13-1	UG L	15	0	
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG L	0.5	U	
8260C	1,1-DICHLOROETHANE	75-34-3	UG L	0.5	Ü	-
8260C	1,1-DICHLOROETHANE 1,1-DICHLOROETHENE	75-34-3 75-35-4	UG L	0.5	U	
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG L	0.5	U	
			UG L		U	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8		0.75	_	
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	U	ļ
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	-
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	U	ļ
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	2.7		ļ
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	С
8260C	2-HEXANONE	591-78-6	UG_L	2.5	UJ	С
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	U	
8260C	ACETONE	67-64-1	UG_L	2.5	UJ	С
8260C	BENZENE	71-43-2	UG L	0.5	U	
8260C	BROMODICHLOROMETHANE	75-27-4	UG L	0.5	U	1
8260C	BROMOFORM	75-25-2	UG L	0.5	Ū	
8260C	BROMOMETHANE	74-83-9	UG L	1	Ü	
8260C	CARBON DISULFIDE	75-15-0	UG L	0.5	Ü	1
8260C	CARBON TETRACHLORIDE	56-23-5	UG L	0.5	Ü	1
8260C	CHLOROBENZENE	108-90-7	UG L	0.5	Ü	
8260C	CHLOROETHANE	75-00-3	UG L	1	UJ	С
8260C	CHLOROFORM	67-66-3	UG L	0.5	U	
8260C	CHLOROMETHANE	74-87-3	UG L	1	Ü	
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG L	2.7		1
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG L	0.5	U	-
8260C	CYCLOHEXANE	110-82-7	UG L	0.5	Ü	-
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG L	0.5	Ü	-
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG L	1	Ü	-
8260C	ETHYLBENZENE	100-41-4	UG L	0.5	Ü	+
8260C	ISOPROPYLBENZENE	98-82-8	UG L	0.5	U	+
8260C 8260C		108-38-3/106-42	UG_L	0.5 1	U	
	M- AND P-XYLENE				U	
8260C	METHYL ACETATE	79-20-9	UG_L	0.75		
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	U	ļ
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	<u> </u>
8260C	STYRENE	100-42-5	UG_L	0.5	U	<u> </u>
8260C	TETRACHLOROETHENE	127-18-4	UG_L	6.4		<u> </u>
8260C	TOLUENE	108-88-3	UG_L	0.5	U	1
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	U	ļ
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	UJ	С
8260C	TRICHLOROETHENE	79-01-6	UG_L	140		
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	
270D SIM	1,4-DIOXANE	123-91-1	UG L	9.3	I '	

		Sample	Delivery Group Lab ID		THPAGE-3 0428-3RA	3
			Sample ID Sample Date Sample Type	12	03-GW-12 2/29/2015 oundwater	
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	4.9		
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	Ü	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG L	0.5	Ü	
8260C	1,2-DICHLOROETHANE	107-06-2	UG L	0.5	Ü	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG L	1	Ü	
8260C	1,2-DICHLOROPROPANE	78-87-5	UG L	0.5	Ü	
8260C	1.3-DICHLOROBENZENE	541-73-1	UG L	0.5	Ü	
8260C	1,4-DICHLOROBENZENE	106-46-7	UG L	0.5	Ü	
8260C	2-BUTANONE	78-93-3	UG L	2.5	Ü	
8260C	2-HEXANONE	591-78-6	UG L	2.5	Ü	
8260C	4-METHYL-2-PENTANONE	108-10-1	UG L	2.5	Ü	
8260C	ACETONE	67-64-1	UG L	2.5	UJ	С
8260C	BENZENE	71-43-2	UG L	0.5	Ü	•
8260C	BROMODICHLOROMETHANE	75-27-4	UG L	0.5	Ü	
8260C	BROMOFORM	75-25-2	UG L	0.5	Ü	
8260C	BROMOMETHANE	74-83-9	UG L	1	U	
8260C	CARBON DISULFIDE	75-15-0	UG L	0.5	Ü	
8260C	CARBON TETRACHLORIDE	56-23-5	UG L	0.5	Ü	
8260C	CHLOROBENZENE	108-90-7	UG L	0.5	Ü	
8260C	CHLOROETHANE	75-00-3	UG L	1	Ü	
8260C	CHLOROFORM	67-66-3	UG L	0.5	Ü	
8260C	CHLOROMETHANE	74-87-3	UG L	1	Ü	
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	0.5	Ü	
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG L	0.5	U	
8260C	CYCLOHEXANE	110-82-7	UG L	0.5	Ü	
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG L	0.5	Ü	
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG L	1	Ü	
8260C	ETHYLBENZENE	100-41-4	UG L	0.5	Ü	
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	0.75	Ü	
8260C	METHYL CYCLONEXANE 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	•
8260C	TOLUENE	108-88-3	UG_L	0.5	U	С
8260C 8260C	TRANS-1.2-DICHLOROETHENE				U	
8260C 8260C	TRANS-1,2-DICHLOROETHENE TRANS-1,3-DICHLOROPROPENE	156-60-5	UG_L	0.5	U	
8260C 8260C	- ,	10061-02-6	UG_L	0.5		
	TRICHLOROETHENE	79-01-6	UG_L	0.36	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	0.17	U	