2014-2015 OU2 GROUNDWATER INVESTIGATION BPOW6-1, BPOW6-2, BPOW6-3, BPOW6-4, BPOW6-5, BPOW6-6 (VPB 145, VPB 146, VPB 147) Installation Report

BETHPAGE, NY

Prepared for:



Department of the Navy Naval Facilities Engineering Command, Mid-Atlantic 9742 Maryland Ave. Norfolk, VA 23511-3095

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Prepared by:



Resolution Consultants

A Joint Venture of AECOM & EnSafe

1500 Wells Fargo Building

440 Monticello Avenue

Norfolk, VA 23510

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List of Acronyms and Abbreviations

AOC Area of Concern
bgs below ground surface
COC Contaminants of Concern

COR Continuously Operating Reference

EPA Environmental Protection Agency, United States

ft feet

GOCO Government-Owned Contractor-Operated

GPS Global Positioning System
IDW Investigation Derived Waste
IR Installation Restoration

Katahdin Katahdin Analytical Services, Inc MWD Massapequa Water District NAD North American Datum

NAVD North American Vertical Datum

NAVFAC Naval Facilities Engineering Command

NG Northrop Grumman

NTU nephlometric turbidity units

NWIRP Naval Weapons Industrial Reserve Plant

NYSDEC New York State Department of Environmental Conservation

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

TCE Trichloroethene

TCL Target Compound List

TCLP Toxicity Characteristic Leaching Procedure

TOC Total Organic Carbon
UFP United Federal Programs

US United States

VOC Volatile Organic Compounds

VPB Vertical Profile Boring

1.0 PROJECT BACKGROUND

Resolution Consultants has prepared this Data Summary Report for the Naval Facilities Engineering Command (NAVFAC), Mid-Atlantic under contract task order WE15 Contract N62470-11-D-8013. This report describes the installation of six monitoring wells and two quarterly groundwater monitoring events (specifically at the Vertical Profile Boring [VPB] locations VPB 145, VPB 146, VPB 147) in 2014 and 2015 for the Naval Weapons Industrial Reserve Plant (NWIRP) Bethpage Operable Unit (OU) 2 Site 1 offsite plume. NWIRP Bethpage is located in east-central Nassau County, Long Island, New York, approximately 30 miles east of New York City (Figure 1).

1.1 Scope and Objectives

This report provides information on the installation of BPOW6-1, BPOW6-2, BPOW6-3, BPOW6-4, BPOW6-5 and BPOW6-6. The purpose of these outpost wells is to ascertain subsurface conditions and contaminant levels upgradient of the Massapequa Water District (MWD) wells and to provide early warning of plume migration toward the MWD wellfield. The locations of BPOW6-1, BPOW6-2, BPOW6-3, BPOW6-4, BPOW6-5 and BPOW6-6, VPBs and monitoring well locations are shown in Figure 2.

The field investigation included completing six monitoring wells, well development, soil/groundwater analysis, groundwater grab samples, and surveying. The tasks also included the decommissioning and relocation of wells BPOW6-5 and BPOW6-6 when it was determined while drilling the borehole for BPOW6-6 in February 2015 that it had damaged BPOW6-5. These activities are described and documented in this report and in Appendix A.

Field tasks were conducted in 2014 and 2015 in accordance with the *United Federal Programs Sampling and Analysis Plan (UFP SAP)*, Bethpage, New York. In addition, the work adhered to the following UFP SAP Addendums: Groundwater Sampling Using Low Stress (Low Flow) Purging and Sampling Protocol (Resolution Consultants, 2013) and Installation of Vertical Profile Borings and Monitoring Wells (Resolution Consultants, 2013).

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 Properties; however, a small portion near Sites 2 and 3 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 observed onsite. The Magothy is characterized by fine to medium sands and silts interbedded with zones of clays, silty sands and sandy clays. Sand and gravel lenses are found in some areas between depths of 600 and 880 ft below ground surface (bgs); these deposits form the Magothy Aquifer.

Investigations performed by the Navy since 2012 indicate that the bottom of the Magothy (top of the Raritan Clay) can extend to depths of 700 to greater than 1,000 ft bgs. The top of the Raritan Clay deepens to the south southeast, as evidenced by clay depths of 1,000 ft bgs (or more) in borings installed offsite. The Raritan Clay Unit is of continental origin and consists of clay, silty clay, clayey silt, and fine silty sand. This member acts as a confining layer over the Lloyd Sand Unit. The

Lloyd Sand Unit is also of continental origin, having been deposited in a large fresh water lacustrine environment. The material consists of fine to coarse-grained sands, gravel, inter-bedded clay, and silty sand. These deposits form the Lloyd Aquifer.

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

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

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

2.0 FIELD PROGRAM

Six monitoring wells were installed in the vicinity of VPB 145, VPB 146 and VPB 147 between July 2014 and June 2015. Field investigation activities consisted of drilling, well installation, well development, sampling, soil/groundwater analysis, surveying and decommissioning and relocation of two of the monitoring wells. 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 BPOW6-1, BPOW6-2, BPOW6-3, BPOW6-4, BPOW6-5 and BPOW6-6 were installed using mud rotary drilling techniques. Well locations are shown on Figure 2 and construction details are summarized in Table 1. Boring logs with lithologic descriptions limited to the well screen interval are included in the Appendix A. Data Summary Reports for VPB 145 (Resolution Consultants, 2014a), VPB 146 (Resolution Consultants, 2014b) and VPB 147 (Resolution Consultants, 2015) document the installation of these VPBs including detailed lithologic descriptions, continuous gamma plots and multiple VOC samples 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 and also to coincide with MWD wells MWD 6442 and MWD 6443. 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 Decommissioning and Relocation of BPOW6-5 and BPOW6-6

Monitoring well BPOW6-5 was initially installed on February 11, 2015 on Crocus Street, Seaford, New York. The total depth was 795 feet bgs. On Friday, February 27, 2015, pieces of Schedule 80 PVC well casing were observed mixed with the cuttings in the drilling mud. It was determined that

borehole BPOW6-6 had intersected the casing of well BPOW6-5, located 13 feet to the west, at a depth of approximately 460 to 480 feet. Resolution Consultants and the Navy assessed the situation and concluded that well BPOW6-5 could not be salvaged. Therefore, the decision was made to decommission well BPOW6-5 and borehole BPOW6-6.

The borehole decommissioning procedure was to tremie grout both the well and the boring from the total depth of each borehole to the ground surface to prevent potential contaminants of concern (COCs) from entering the borehole or the well and migrating vertically between subsurface zones. Decommissioning of both well BPOW6-5 and borehole BPOW6-6 was successfully completed on April 9, 2015. Documentation of the implementation is provided in Appendix A. Monitoring wells BPOW6-5 and BPOW6-6 were relocated and completed on North Hickory Street, Massapequa on May 8, 2015 and May 27, 2015, respectively.

2.3 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 nephlometric turbidity units (NTUs) if possible. Table 2 summarizes total pumped volume from air and pump development and final turbidity.

2.4 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 Volatile Organic Compounds (VOC)s 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 are included in Appendix A.

Monitoring wells BPOW6-1, BPOW6-2, BPOW6-3, BPOW6-4, BPOW6-5 and BPOW6-6 are sampled quarterly as part of the Navy's ongoing Environmental Restoration Program. Resolution Consultants sampled BPOW6-1, BPOW6-2, BPOW6-3, and BPOW6-4 during the March 2015 quarterly monitoring event and sampled BPOW6-5 and BPOW6-6 during the June quarterly monitoring event. Analytical results and stabilized field parameters for these data are summarized in Tables 3 and 4, respectively. Data validation is documented in Appendix A.

Per an agreement between the Navy and Northrop Grumman, Northrop Grumman has taken over the quarterly sampling of BPOW6-1, BPOW6-2, BPOW6-3, and BPOW6-4 since the initial March 2015 quarterly monitoring event. Data validation and results of the June 2015 quarterly monitoring of these wells by Northrop Grumman's consultant is documented in Appendix A. Northrop Grumman will take over the quarterly sampling of BPOW6-5 and BPOW6-6 for the 2015 third quarter sampling event.

2.5 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, 2013).

As part of the IDW management practices and in accordance with the SAP, the investigation waste (consisting of 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.

2.6 Surveying

A survey of the 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 location was 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 (ground, latitude/longitude and northing/easting) and a survey map is included in Appendix A.

3.0 REFERENCES

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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.

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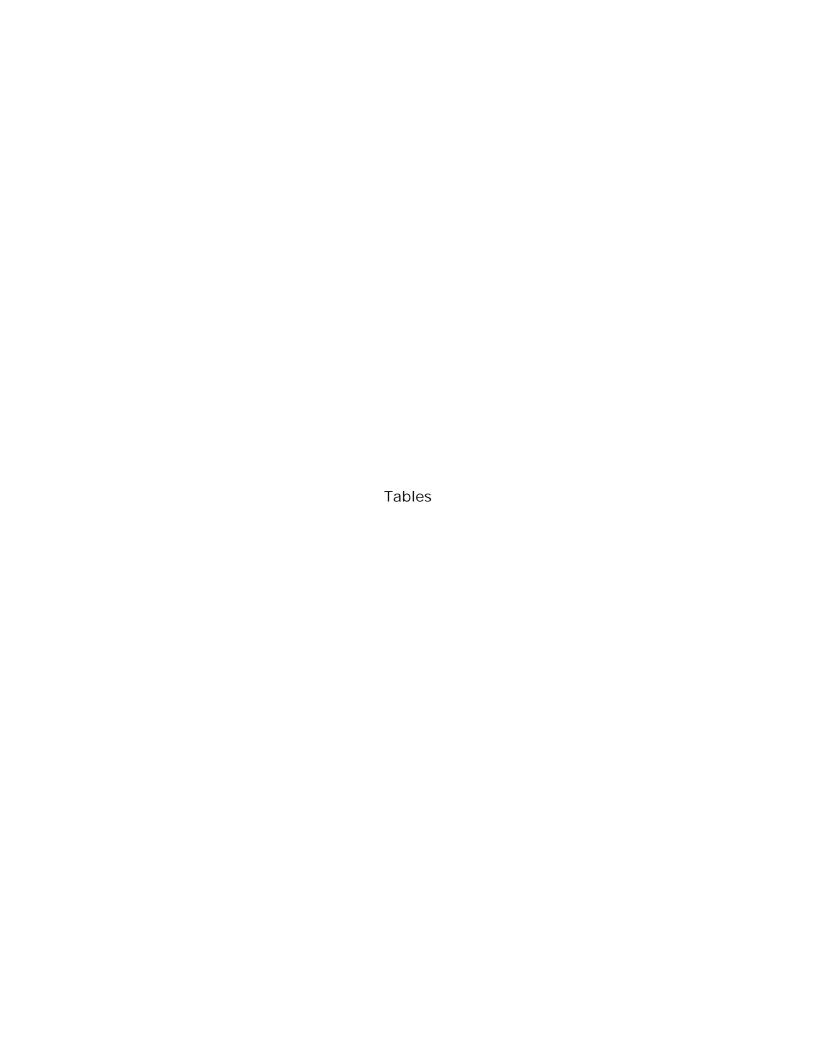


TABLE 1 MONITORING WELL CONSTRUCTION SUMMARY 2014-2015 OU2 GROUNDWATER INVESTIGATION NWIRP BETHPAGE, NY

MONITORING WELL	WELL COMPLETION DATE	GROUND ELEVATION (MSL)	PVC (INNER CASING) (MSL)	WELL DEPTH (ft bgs)	CASING DEPTH (ft bgs)	SCREEN INTERVAL (ft bgs)	SUMP DEPTH INTERVAL (ft bgs)	BORING DEPTH (ft bgs)
BPOW6-1	9/15/2014	43.61	42.93	580	52.5	550-575	5	595
BPOW6-2	8/18/2014	43.58	43.08	785	51.7	755-780	5	798
BPOW6-3	11/25/2014	40.34	39.96	780	52	750-775	5	795
BPOW6-4	12/16/2014	40.40	40.02	575	52.5	545-570	5	590
BPOW6-5	5/8/2015	43.27	42.58	555	54	525-550	5	567
BPOW6-6	5/27/2015	43.17	42.34	800	54	770-795	5	812

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

	AIR DEVEL	OPMENT	PUM	IP DEVELOPME	APPROX. TOTAL	FINAL		
MONITORING WELL	DATE	APPROX. VOLUME (GAL)	DATE	FINAL PUMP DEPTH (FT BGS)	APPROX. VOLUME (GAL)	DEVELOPMENT VOLUME (GAL)	TURBIDITY (NTUs)	
BPOW6-1	10/7/2014	4,800	10/13/2014	550-575	10,000	14,800	4.61	
BPOW6-2	10/6/2014	5,000	10/8/2014- 10/10/14	755-780	10,000	15,000	60	
BPOW6-3	1/5/2015	10,000	1/7/15- 1/8/2015	750-775	5,000	15,000	30.03	
BPOW6-4	1/6/2015	5,000	1/6/2015	545-570	4,000	9,000	0.68	
BPOW6-5	6/3/2015	7,400	6/5/2015	525-550	4,700	12,100	7.76	
BPOW6-6	6/4/2015	8,300	6/8/2015	770-795	4,000	12,300	30.27	

Location Sample Date Sample ID	NYSDEC Groundwater Guidance or Standard Value	BPOW6-1 3/26/2015 BPOW6-1-GW- 032615	BPOW6-2 3/26/2015 BPOW6-2-GW- 032615	BPOW6-3 3/26/2015 BPOW6-3-GW- 032615	BPOW6-4 3/26/2015 BPOW6-4-GW- 032615
Sample type code	(Note 1)	N	N	N	N
VOC 8260C (ug/L)					
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	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
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	< 0.75 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	< 0.50 U
1,2-DICHLOROETHENE, TOTAL	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
1,2-DICHLOROPROPANE	1	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,3-DICHLOROBENZENE	3	< 0.50 U	< 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
1,4-DIOXANE (Method 8270D_SIM)	NL	< 0.17 U	< 0.19 U	< 0.17 U	< 0.17 U
2-BUTANONE	50	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ
2-HEXANONE	50	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ
4-METHYL-2-PENTANONE	NL	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 U
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 UJ	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ
CARBON DISULFIDE	60	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
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 UJ	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ
CHLOROFORM	7	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROMETHANE	5	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ
CIS-1,2-DICHLOROETHENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 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 UJ	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
DIBROMOCHLOROMETHANE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
DICHLORODIFLUOROMETHANE	5	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ
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 U	< 2.5 U	< 2.5 U	< 2.5 U
O-XYLENE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
STYRENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TETRACHLOROETHENE	5	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
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	< 0.50 U	< 0.50 U	< 0.50 U	< 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	< 1.5 U	< 1.5 U	< 1.5 U	< 1.5 U

Location Sample Date	NYSDEC Groundwater	BPOW6-5 6/25/2015	BPOW6-6 6/25/2015
Sample ID	Guidance or Standard Value	BPOW6-5-GW-	PPOW/ / CW 0/3515
Sample type code	(Note 1)	062515 N	BPOW6-6-GW-062515 N
VOC 8260C (ug/L)			
1,1,1-TRICHLOROETHANE	5	< 0.50 U	< 0.50 U
1,1,2,2-TETRACHLOROETHANE	5	< 0.50 U	< 0.50 U
1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	5	< 0.50 U	< 0.50 U
1,1,2-TRICHLOROETHANE	1	< 0.50 U	< 0.50 U
1,1-DICHLOROETHANE	5	< 0.50 U	< 0.50 U
1,1-DICHLOROETHENE	5	< 0.50 U	< 0.50 U
1,2,4-TRICHLOROBENZENE	5	< 0.50 U	< 0.50 U
1,2-DIBROMO-3-CHLOROPROPANE	0.04	< 0.75 U	< 0.75 U
1,2-DIBROMOETHANE	NL	< 0.50 U	< 0.50 U
1,2-DICHLOROBENZENE	3	< 0.50 U	< 0.50 U
1,2-DICHLOROETHANE	5	< 0.50 U	< 0.50 U
1,2-DICHLOROETHENE, TOTAL	5	< 1.0 U	< 1.0 U
1,2-DICHLOROPROPANE	1	< 0.50 U	< 0.50 U
1,3-DICHLOROBENZENE	3	< 0.50 U	< 0.50 U
1,4-DICHLOROBENZENE	3	< 0.50 U	< 0.50 U
1,4-DIOXANE (Method 8270D_SIM)	NL	< 0.17 U	< 0.17 U
2-BUTANONE	50	< 2.5 U	< 2.5 U
2-HEXANONE	50	< 2.5 U	< 2.5 U
4-METHYL-2-PENTANONE	NL	< 2.5 U	< 2.5 U
ACETONE	50	< 2.5 U	< 2.5 U
BENZENE	1	< 0.50 U	< 0.50 U
BROMODICHLOROMETHANE	50	< 0.50 U	< 0.50 U
BROMOFORM	50	< 0.50 U	< 0.50 U
BROMOMETHANE	5	< 1.0 U	< 1.0 U
CARBON DISULFIDE	60	< 0.50 U	< 0.50 U
CARBON TETRACHLORIDE	5	< 0.50 U	< 0.50 U
CHLOROBENZENE	5	< 0.50 U	< 0.50 U
CHLOROETHANE	5	< 1.0 UJ	< 1.0 UJ
CHLOROFORM	7	< 0.50 U	< 0.50 U
CHLOROMETHANE	5	< 1.0 U	< 1.0 U
CIS-1,2-DICHLOROETHENE	5	< 0.50 U	< 0.50 U
CIS-1,3-DICHLOROPROPENE	0.4	< 0.50 U	< 0.50 U
CYCLOHEXANE	NL	< 0.50 U	< 0.50 U
DIBROMOCHLOROMETHANE	5	< 0.50 U	< 0.50 U
DICHLORODIFLUOROMETHANE	5	< 1.0 U	< 1.0 U
ETHYLBENZENE	5	< 0.50 U	< 0.50 U
ISOPROPYLBENZENE	5	< 0.50 U	< 0.50 U
M- AND P-XYLENE	NL	< 1.0 U	< 1.0 U
METHYL ACETATE	NL	< 0.75 U	< 0.75 U
METHYL CYCLOHEXANE	NL	< 0.50 U	< 0.50 U
METHYL TERT-BUTYL ETHER	10	< 0.50 U	< 0.50 U
METHYLENE CHLORIDE	5	< 2.5 U	< 2.5 U
O-XYLENE	NL	< 0.50 U	< 0.50 U
STYRENE	5	< 0.50 U	< 0.50 U
TETRACHLOROETHENE	5	< 0.50 U	< 0.50 U
TOLUENE	5	0.76 J	1.0
TRANS-1,2-DICHLOROETHENE	5	< 0.50 U	< 0.50 U
TRANS-1,3-DICHLOROPROPENE	0.4	< 0.50 U	< 0.50 U
TRICHLOROETHENE	5	< 0.50 U	< 0.50 U
TRICHLOROFLUOROMETHANE	5	< 1.0 U	< 1.0 U
VINYL CHLORIDE	2	< 1.0 U	< 1.0 U
XYLENES, TOTAL	5	< 1.5 U	< 1.5 U

Notes:

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

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

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

Yellow highlighted values exceed Groundwater Standards or guidance value

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

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

UJ = The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is

approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte.

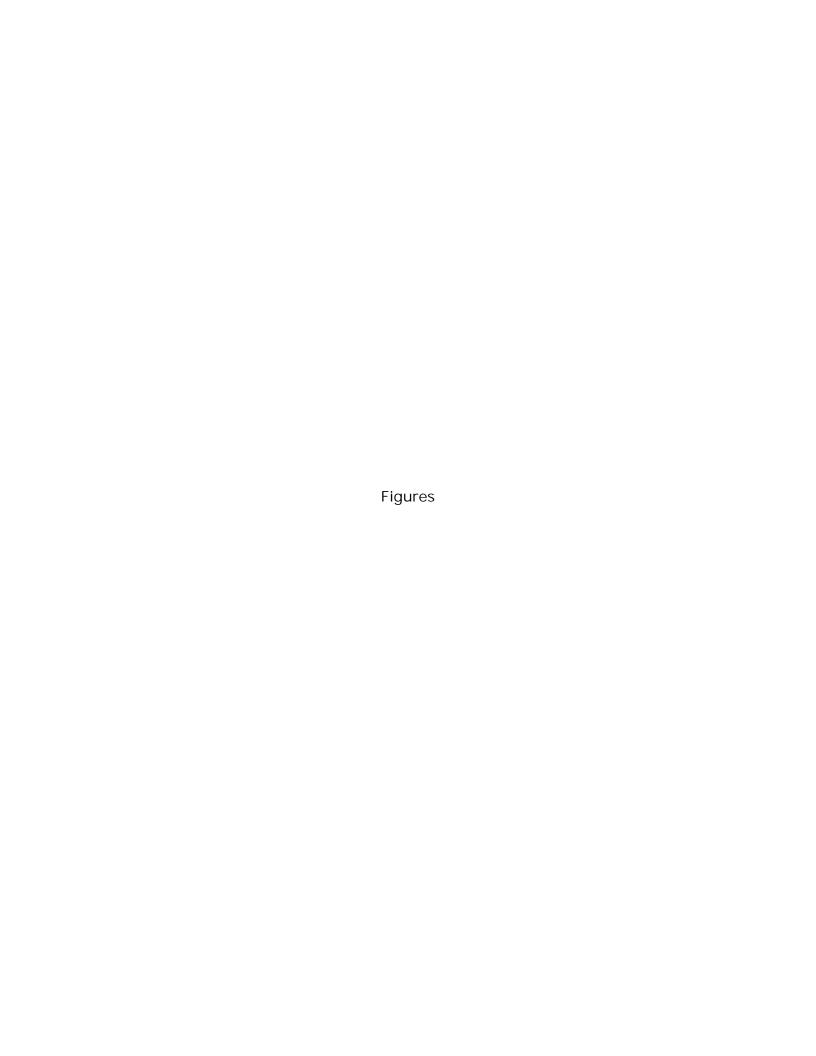
J = The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.

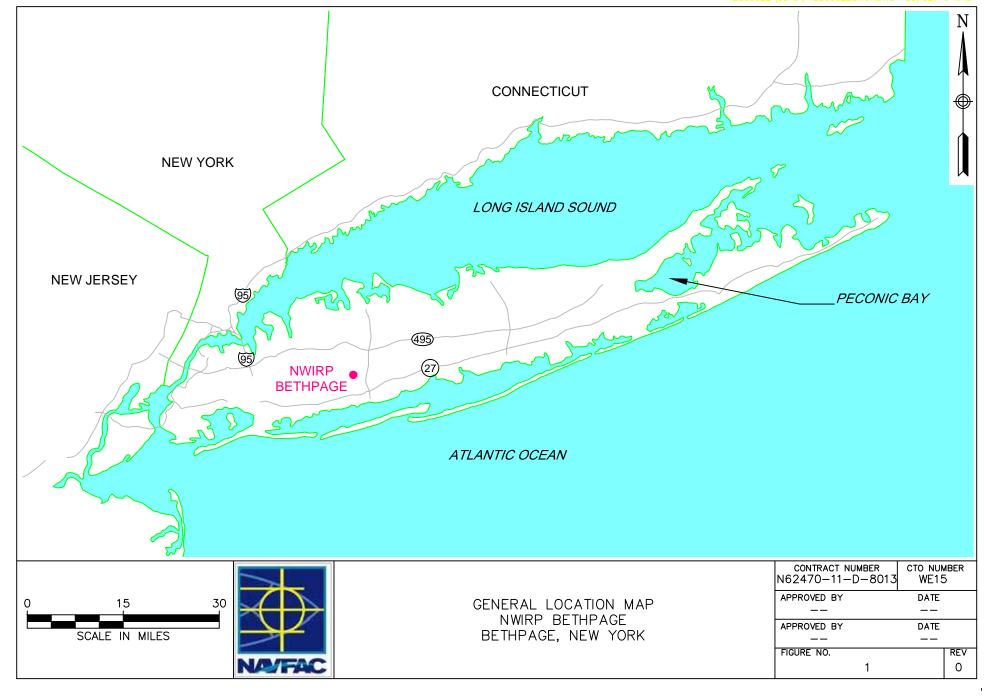
M = the matrix spike or matrix spike duplicate did not meet recovery or precision requirements.

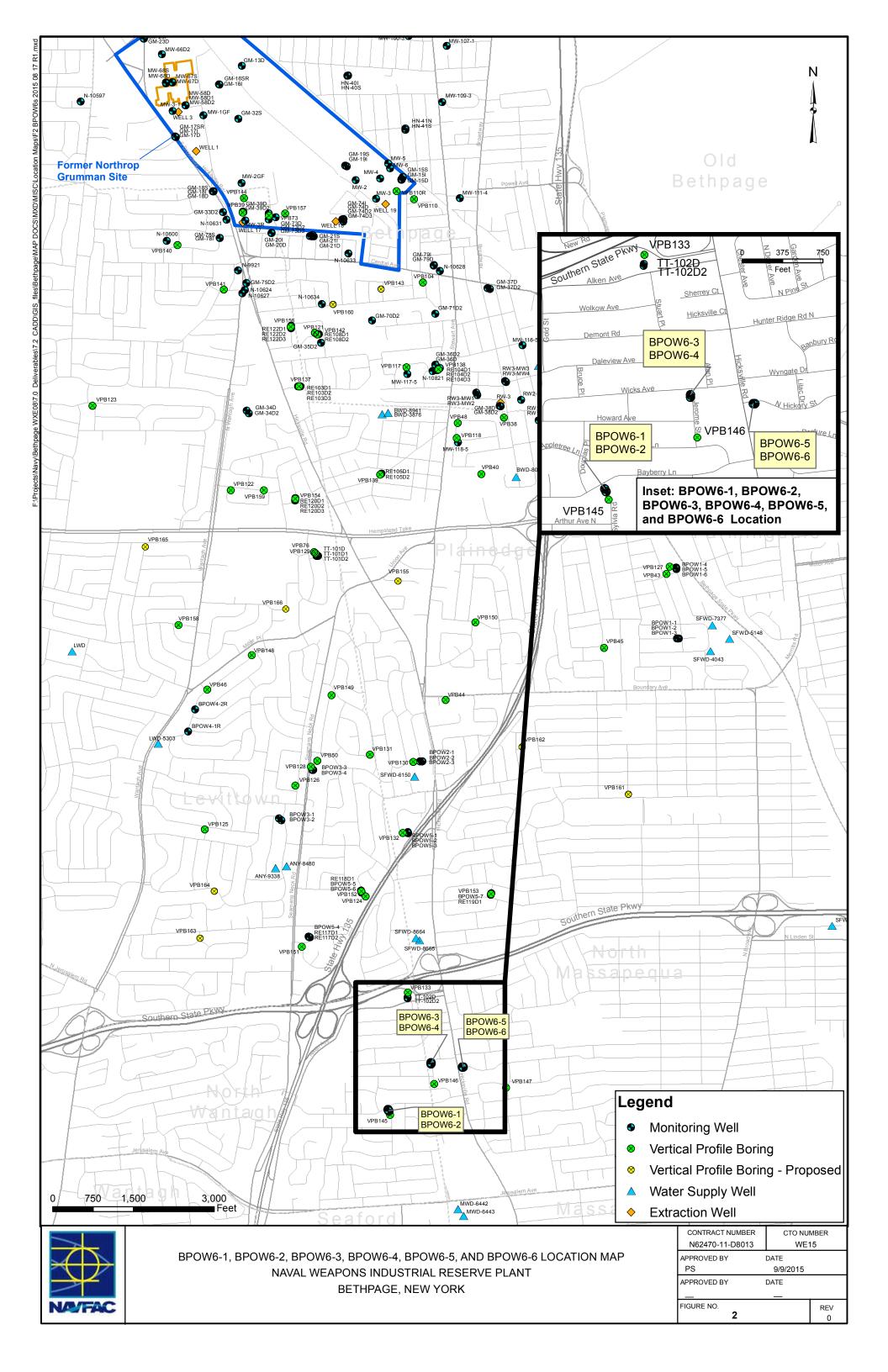
Table 4 Stabilized Field Parameters

Well	Date	Temperature (°C)	рН	Specific Conductance (μS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Depth to water (ft bgs)	Flow rate (ml/min)
BPOW6-1	3/26/2015	12.94	5.10	0.088	1.18	60.8	163	13.6	500
BPOW6-2	3/26/2015	12.98	5.33	0.032	1.18	-22.5	30	13.97	600
BPOW6-3	3/26/2015	12.88	4.94	0.027	0.88	218.4	4.47	10.78	550
BPOW6-4	3/26/2015	12.91	4.59	0.093	0.91	-34.1	3.99	10.11	750
BPOW6-5	6/25/2015	20.80	4.81	0.054	0.58	221.1	1.57	18.00	200
BPOW6-6	6/25/2015	16.46	4.10	0.028	0.61	330.7	31.6	18.51	700

^{*} Initial water level not equilibrated due to pump installation; drawdown during sampling not determined.







Appendix A

BPOW6-1, BPOW6-2, BPOW6-3, BPOW6-4, BPOW6-5, BPOW6-6

Section 1

Boring Logs

Boring Log

BORING #: BPOW6-1
Sheet 1 of 2

Client: Department of the Navy, Naval Facilit	Logged By: V. Thayer	
Location: Bayberry Ln and Sylvia Rd, Seaford	Drilling Company: DELTA WELL AND PUMP COMPANY	
Project #: 60266526	Well Screen Interval (ft): 550-575	
Start Date: 8/21/2014	Drilling Method: Auger (0-50' bgs) Mud Rotary (>50' bgs)	Water Level (ft):
Finish Date: 9/15/2014	Northing: 193228.16 Easting: 1126796.92	Total Depth (ft): 595.0

		ı					T
ДЕРТН (ft)	PID (ppm)	Formation	SOSN	GRAPHIC LOG	MATERIAL DESCRIPTION	Well	Well Construction
0					0-553 ft bgs: See VPB145 for Descriptions		10" Diameter Steel Casing
50							Gasing
100							
150							
200							
250						-	Bentonite Grout
300							
350							
400							
450						-	4" Diameter Schedule 80 PVC Riser
500						0000000	

Boring Log

BORING #: BPOW6-1 Sheet 2 of 2

Client: Department of the Navy, Naval Facilit	Logged By: V. Thayer	
Location: Bayberry Ln and Sylvia Rd, Seaford	Drilling Company: DELTA WELL AND PUMP COMPANY	
Project #: 60266526	Well Screen Interval (ft): 550-575	
Start Date: 8/21/2014	Drilling Method: Auger (0-50' bgs) Mud Rotary (>50' bgs)	Water Level (ft):
Finish Date: 9/15/2014	Northing: 193228.16 Easting: 1126796.92	Total Depth (ft): 595.0

DЕРТН (ft)	PID (ppm)	Formation	SSO	GRAPHIC LOG	MATERIAL DESCRIPTION	Well	Well Construction
506 508 510 512 514 516					0-553 ft bgs: See VPB145 for Descriptions (continued)		4" Diameter Schedule 80 PVC Riser (continued)
518 520 522 524 526 528 530						-	#0 Filter Sand
532 534 536 538 540 542 544 546 548 550 552							#1 Filter Sand
554 556 558 560	0 0		SP-SM SP-SM		Light gray (Gley 1 7/1) poorly graded SAND with Silt, angular medium sand, trace to few silt Dark gray (Gley 1 4/1) poorly graded SAND with Silt, angular medium sand, bottom 2 in. sand interbedded with black laminated lignite		
562 564 566	0		SP		Gray (Gley 15/1) poorly graded SAND, angular medium Sand, one 1/4" band of lignite		4" Diameter Schedule 80 PVC, 10 Slot Well Screen (550-575 ft bgs)
568 570 572	0		SP		Gray (Gley 1 5/1) poorly graded SAND, angular medium Sand, muscovite flakes, six 1/4" layers of lignite		(555 5.5 // 1295)
574 576 578 580 582	0		SP		Gray (Gley 1 6/1) poorly graded SAND, subangular medium to coarse Sand, two black layers of 1/4" lignite		Sump
584 586 588 590 592							#1 Sand to bottom
594					End of boring at 595.0 ft. bgs.		

Boring Log

BORING #: BPOW6-2 Sheet 1 of 2

Client: Department of the Navy, Naval Facilit	Logged By: V. Thayer			
Location: Bayberry Ln and Sylvia Rd, Seaford	Drilling Company: DELTA WELL AND PUMP COMPANY			
Project #: 60266526	Project #: 60266526			
Start Date: 7/25/2014	Date: 7/25/2014 Drilling Method: Auger (0-50' bgs) Mud Rotary (>50' bgs)			
Finish Date: 8/18/2014	Northing: 193253.75 Easting: 1126785.34	Total Depth (ft): 798.0		

					T	I	I
DEPTH (ft)	PID (ppm)	Formation	nscs	GRAPHIC LOG	MATERIAL DESCRIPTION	Well	Well Construction
0					0-758 ft bgs: See VPB145 for Descriptions		4011 Diversity Object
50							10" Diameter Steel Casing
100							
150							
200							
250							
300						-	Bentonite Grout
300							
350							
400							
450							
500							
550							
600						-	4" Diameter Schedule 80 PVC Riser
							80 PVC Riser
650							
700							

Boring Log

BORING #: BPOW6-2 Sheet 2 of 2

Client: Department of the Navy, Naval Facilit	Logged By: V. Thayer		
Location: Bayberry Ln and Sylvia Rd, Seaford	Drilling Company: DELTA WELL AND PUMP COMPANY		
Project #: 60266526	Well Screen Interval (ft): 755-780		
Start Date: 7/25/2014	Drilling Method: Auger (0-50' bgs) Mud Rotary (>50' bgs)	Water Level (ft):	
Finish Date: 8/18/2014	Northing: 193253.75 Easting: 1126785.34	Total Depth (ft): 798.0	

702 708 709 709 710 711 711 713 714 715 716 718 720 722 724 724 725 728 730 732 733 734 736 737 738 739 740 741 744 744 745 745 756 758 758 758 758 758 758 758 758 758 758	Well Construction	Completion	Well	MATERIAL DESCRIPTION	GRAPHIC LOG	nscs	Formation	PID (ppm)	DEPTH (ft)
714 716 718 720 722 722 724 724 726 728 739 730 732 734 736 738 738 740 741 742 744 746 748 759 750 752 754 756 758 0 SW-SC SW	4" Diameter Schedule 80 PVC Riser (continued)		60000	0-758 ft bgs: See VPB145 for Descriptions (continued)					702 - 704 - 706 - 708 - 710
Total	#0 Filter Sand								714 716 718 720 722 724 726 728 730 732 734 736 738 740 742 744 746
gravel, few clay Light gray (Gley 1) widely graded SAND with Clay, medium to coarse sand, few subrounded fine gravel, few clay Light gray (Gley 1) widely graded SAND with Clay, medium to coarse sand, trace subrounded fine gravel, few clay Light gray (Gley 1) widely graded SAND with Clay, medium to coarse sand, trace subrounded fine gravel, few clay Light gray (Gley 1) widely graded GRAVEL, subangular fine to coarse Gravel, few white clay Light gray (Gley 1) widely graded GRAVEL with Clay, subrounded fine to coarse gravel, some fine to coarse sand, few clay	#1 Filter Sand			Light gray (Gley 1) widely graded SAND with Clay, angular	•••••			0	754 756
Coarse sand, few subrounded fine gravel, few clay Light gray (Gley 1) widely graded SAND with Clay, medium to coarse sand, trace subrounded fine gravel, few clay SW-SC Light gray (Gley 1) widely graded GRAVEL, subangular fine to coarse Gravel, few white clay Light gray (Gley 1) widely graded GRAVEL with Clay, subrounded fine to coarse gravel, some fine to coarse sand, few clay				medium to coarse sand, little fine sand, trace subrounded fine gravel, few clay		SW-SC			
Total Course sand, trace subrounded fine gravel, few clay SW-SC SW-SC Light gray (Gley 1) widely graded SAND with Clay, medium to coarse sand, trace subrounded fine gravel, few clay Light gray (Gley 1) widely graded GRAVEL, subangular fine to coarse Gravel, few white clay Light gray (Gley 1) widely graded GRAVEL with Clay, subrounded fine to coarse gravel, some fine to coarse sand, few clay				Light gray (Gley 1) widely graded SAND with Clay, medium to coarse sand, few subrounded fine gravel, few clay		SW-SC		0	
Light gray (Gley 1) widely graded GRAVEL, subangular fine to coarse Gravel, few white clay CW-GC Light gray (Gley 1) widely graded GRAVEL, subangular fine to coarse Gravel, few white clay Light gray (Gley 1) widely graded GRAVEL with Clay, subrounded fine to coarse gravel, some fine to coarse sand, few clay	4" Diameter Schedule 80 PVC, 10 Slot Well Screen					SW-SC		0	770
T78 0 Using the control of the contr	(755-780 ft bgs)			Light gray (Gley 1) widely graded GRAVEL, subangular fine to coarse Gravel, few white clay		GW-GC		0	774
GW-GC GW-GC few clay				Light gray (Gley 1) widely graded GRAVEL with Clay,				0	778
784 786 788 790	Sump #1 Sand to bottom			subrounded fine to coarse gravel, some fine to coarse sand,		GW-GC			782 784 786 788 790 792 794

Boring Log

BORING #: BPOW6-3
Sheet 1 of 2

Client: Department of the Navy, Naval Facilit	Logged By: V. Thayer		
Location: Jerome St & Wicks Ave, Seaford, N	Drilling Company: DELTA WELL AND PUMP COMPANY		
Project #: 60266526	Well Screen Interval (ft): 750-775		
Start Date: 10/28/2014	Drilling Method: Auger (0-50' bgs) Mud Rotary (>50' bgs)	Water Level (ft):	
Finish Date: 11/25/2014	Northing: 194106.07 Easting: 1127582.36	Total Depth (ft): 795.0	

ОЕРТН (ft)	PID (ppm)	Formation	nscs	GRAPHIC LOG	MATERIAL DESCRIPTION	Well	Well Construction
0					0-753 ft bgs: See VPB146 for Descriptions		10" Diameter Steel Casing
50						6,636,535 anist 644	J
100							
150							
200							
250							
300						-	Bentonite Grout
350						00000000000000000000000000000000000000	
400							
450							
500							
550							
600						-	
							4" Diameter Schedule 80 PVC Riser
650							
700							

Boring Log

BORING #: BPOW6-3 Sheet 2 of 2

Client: Department of the Navy, Naval Facilit	Logged By: V. Thayer			
Location: Jerome St & Wicks Ave, Seaford, N	Drilling Company: DELTA WELL AND PUMP COMPANY			
Project #: 60266526	Ground Elevation (msl): 40.34	Well Screen Interval (ft): 750-775		
Start Date: 10/28/2014	Drilling Method: Auger (0-50' bgs) Mud Rotary (>50' bgs)	Water Level (ft):		
Finish Date: 11/25/2014	Northing: 194106.07 Easting: 1127582.36	Total Depth (ft): 795.0		

DEPTH (ft)	PID (ppm)	Formation	SSSN	GRAPHIC LOG	MATERIAL DESCRIPTION	Well	Well Construction
716 718 720					0-753 ft bgs: See VPB146 for Descriptions (continued)		4" Diameter Schedule 80 PVC Riser (continued)
722 724 726 728						-	#0 Filter Sand
730 732 734 736 738							
740 742 744 746							#1 Filter Sand
748 750 752 754 756	0.1		SM		Gray (7.5 YR 5/1) SILTY SAND, micaceous fine Sand, some silt, clay		
758 760 762					No Recovery		
764 766	0.3		SP GC		Light gray (7.5 YR 5/1) SAND with one orange band, fine Sand Light gray (7.5 YR 7/1) CLAYEY GRAVEL, subrounded, fine to coarse Gravel, little fine to coarse sand, clay		4" Diameter Schedule 80 PVC, 10 Slot Well Screen (750-775 ft bgs)
768 770 772	0.1		GC		Light gray (7.5 YR 7/1) CLAYEY GRAVEL, laminated clay (768-768.1) overlying gravel with minor orange staining, fine gravel, little fine to coarse sand, few clay		
774 776 778	0.1		GP-GC		Light gray (7.5 YR 7/1) GRAVEL with Clay; subrounded to subangular fine gravel, little fine to coarse sand, silt, clay		Sump
780 782 784							·
786 788 790							#1 Sand to bottom
792 794					End of boring at 795.0 ft. bgs.		

Boring Log

BORING #: BPOW6-4
Sheet 1 of 2

Client: Department of the Navy, Naval Facilit	Logged By: V. Thayer			
Location: Jerome St & Wicks Ave, Seaford, N	Drilling Company: DELTA WELL AND PUMP COMPANY			
Project #: 60266526	60266526 Ground Elevation (msl): 40.40			
Start Date: 12/3/2014	Drilling Method: Auger (0-50' bgs) Mud Rotary (>50' bgs)	Water Level (ft):		
Finish Date: 12/16/2014	Northing: 194127.00 Easting: 1127580.66	Total Depth (ft): 590.0		

DEPTH (ft)	PID (ppm)	Formation	nscs	GRAPHIC LOG	MATERIAL DESCRIPTION	Well	Completion	Well Construction
0					0-548 ft bgs: See VPB146 for Descriptions			10" Diameter Steel Casing
50								
100								
150								
200							-4	Bentonite Grout
250								
300								
350								
400								
450							-	4" Diameter Schedule 80 PVC Riser

Boring Log

BORING #: BPOW6-4 Sheet 2 of 2

Client: Department of the Navy, Naval Facilit	Logged By: V. Thayer			
Location: Jerome St & Wicks Ave, Seaford, N	Drilling Company: DELTA WELL AND PUMP COMPANY			
Project #: 60266526	Ground Elevation (msl): 40.40	Well Screen Interval (ft): 545-570		
Start Date: 12/3/2014	Drilling Method: Auger (0-50' bgs) Mud Rotary (>50' bgs)	Water Level (ft):		
Finish Date: 12/16/2014	Northing: 194127.00 Easting: 1127580.66	Total Depth (ft): 590.0		

DEPTH (ft)	PID (ppm)	Formation	nscs	GRAPHIC LOG	MATERIAL DESCRIPTION	Well	Well Construction
500 502 504 506 508 510 512					0-548 ft bgs: See VPB146 for Descriptions (continued)		4" Diameter Schedule 80 PVC Riser <i>(continued)</i>
514 516 518 520 522 524 526							#0 Filter Sand
528 530 532 534 536 536 538 540							#1 Filter Sand
544 546 548 550 552	0.1		SP-SM		Gray (5/1) poorly graded SAND with Silt, angular fine to medium sand, few silt		
554 556 558 560	0.1		SP-SM SP-SM		Dark gray (4/1) poorly graded Sand with SILT, medium Sand, few coarse sand, few silt Gray (5/1) poorly graded SAND with Silt, angular, medium sand, little fine sand, few silt		4" Diameter Schedule 80 PVC, 10 Slot Well Screen
562 564 566	0.5		SP		Gray (5/1) poorly graded SAND, fine to medium Sand, muscovite flakes, trace silt, trace small gravel, one 0.25" band of lignite		(545-570 ft bgs)
568 570 572	0.1		SP		Gray (5/1) poorly graded SAND, medium Sand, trace silt, several 0.25" lignite seams		Sump
574 576 578 580							
582 584 586							#1 Sand to bottom
588					End of boring at 590.0 ft. bgs.		

Boring Log

BORING #: BPOW6-5
Sheet 1 of 2

Client: Department of the Navy, Naval Facilit	Logged By: V. Varricchio			
Location: North Hickory, Massapequa, NY	Drilling Company: DELTA WELL AND PUMP COMPANY			
Project #: 60266526	Ground Elevation (msl): 43.27	Well Screen Interval (ft): 525-550		
Start Date: 5/1/2015	Drilling Method: Auger (0-50' bgs) Mud Rotary (>50' bgs)	Water Level (ft):		
Finish Date: 5/8/2015	Northing: 194045.49 Easting: 1128178.28	Total Depth (ft): 567.0		

					,		
DЕРТН (ft)	PID (ppm)	Formation	nscs	GRAPHIC LOG	MATERIAL DESCRIPTION	Well	Well Construction
o					0-528 ft bgs: see VPB147 for Descriptions		— 10" Diameter Steel Casing
50							
100							
150							
200							
250						•	4" Diameter Schedule 80 PVC Riser
300							
350							
400							
450							

Boring Log

BORING #: BPOW6-5 Sheet 2 of 2

Client: Department of the Navy, Naval Facilit	Logged By: V. Varricchio		
Location: North Hickory, Massapequa, NY	Drilling Company: DELTA WELL AND PUMP COMPANY		
Project #: 60266526	Well Screen Interval (ft): 525-550		
Start Date: 5/1/2015	Drilling Method: Auger (0-50' bgs) Mud Rotary (>50' bgs)	Water Level (ft):	
Finish Date: 5/8/2015	Northing: 194045.49 Easting: 1128178.28	Total Depth (ft): 567.0	

DEPTH (ft)	PID (ppm)	Formation	nscs	GRAPHIC LOG	MATERIAL DESCRIPTION	Well	Well Construction
490					0-528 ft bgs: see VPB147 for Descriptions (continued)		Bentonite Grout (continued)
494							
496							
498							
500						620000000000000000000000000000000000000	
502							
504							
508							
510							
512							#1 Filter Sand
514							#11 liter Sand
516							
518							
520							
524							
526							
528	0				Gray (10 YR 5/1) poorly graded medium SAND		
530	0		SP				
534	0		SP		Gray (10 YR 5/1) poorly graded medium SAND		
538	0				Gray (10 YR 5/1) silty fine poorly graded SAND, Lignite banding throughout		4" Diameter schedule
540	0		SM		banding throughout		80 PVC, 10 Slot Well Screen (525-550 ft
542		-			Gray (10 YR 6/1) poorly graded medium SAND		bgs)
544	0		SP		Gray (10 TK 0/1) poorly graded medicin SAND		
546 548		_					
550	0		SP		Gray (10 YR 5/1) poorly graded fine to medium SAND Gray (10 YR 5/1) SANDY lean CLAY		
552			CL				
554				<i>\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\</i>			Sump
556							
558							
560							#1 Sand to bottom of
562							boring
564							
566					End of boring at 567.0 ft. bgs.	1	

Boring Log

BORING #: BPOW6-6 Sheet 1 of 2

Client: Department of the Navy, Naval Facilit	Logged By: S. Wright / P. Kareth		
Location: North Hickory, Massapequa, NY	Drilling Company: DELTA WELL AND PUMP COMPANY		
Project #: 60266526	Ground Elevation (msl): 43.17	Well Screen Interval (ft): 770-795	
Start Date: 5/15/2015	Drilling Method: Auger (0-50' bgs) Mud Rotary (>50' bgs)	Water Level (ft):	
Finish Date: 5/27/2015	Northing: 194044.59 Easting: 1128163.72	Total Depth (ft): 812.0	

O-773 ft bgs: See VPB147 for Descriptions	Diameter Steel
10" E Casir	Diameter Steel ing
Casi	ing
100	
150	
250	
300	
350	
4" Di 80 Pi	iameter Schedule VC Riser
450	
500	
550	
600	
650	
700	

Boring Log

BORING #: BPOW6-6 Sheet 2 of 2

Client: Department of the Navy, Naval Facilit	Logged By: S. Wright / P. Kareth		
Location: North Hickory, Massapequa, NY	Drilling Company: DELTA WELL AND PUMP COMPANY		
Project #: 60266526	Well Screen Interval (ft): 770-795		
Start Date: 5/15/2015	Drilling Method: Auger (0-50' bgs) Mud Rotary (>50' bgs)	Water Level (ft):	
Finish Date: 5/27/2015	Northing: 194044.59 Easting: 1128163.72	Total Depth (ft): 812.0	

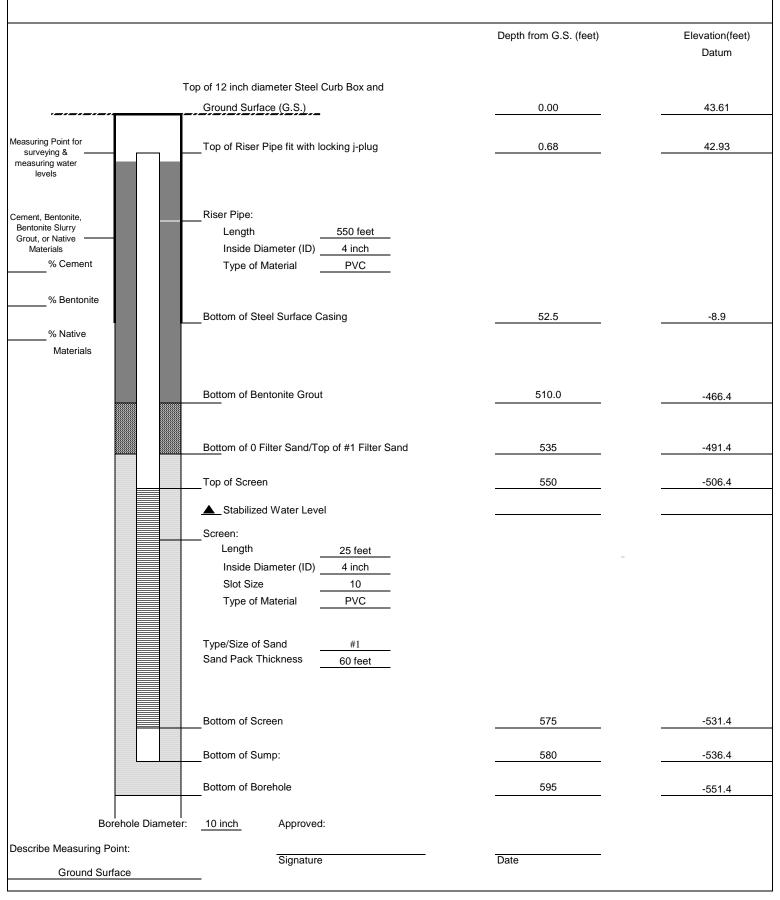
	ı		ı	T		T	Т
DEРТН (ft)	PID (ppm)	Formation	SOSN	GRAPHIC LOG	MATERIAL DESCRIPTION	Well	Well Construction
730 - 732 - 734 - 736 - 738 - 740 - 742					0-773 ft bgs: See VPB147 for Descriptions (continued)	•	Bentonite Grout
744 746 748 750 752 754 756 758 760 762 764 766 768 770							#1 Filter Sand
774 776 778 780	0 ,		SW		Light gray (10 YR 7/1) fine to coarse SAND, 20% fine to coarse Gravel, trace silt Light gray (10 YR 7/1) fine to coarse SAND, 20% fine to coarse Gravel, 5% silt		
782 784 786	0		SM		Light gray with grayish mottling (N7+10 YR 7/4) Silty fine SAND, spoontip is light gray (N7) Clayey fine Sand		4" Diameter schedule 80 PVC, 10 Slot Well Screen (770-795 ft bgs)
790	0		CL		Light Gray (5 YK 6/1) fine sandy CLAY. 7" layer of Silty fine Sand mid spoon, orange mottling in tip		-3-7
794 796 798 798 800	<u> </u>		SP-SM		Light gray (N7) medium to fine SAND, some Silt, few orange mottling, bottom 4" layer clayey fine Sand		Sump
804 806 808 810							#1 Sand to bottom of boring
812					End of boring at 812.0 ft. bgs.		

Section 2

Monitoring Well Construction Log

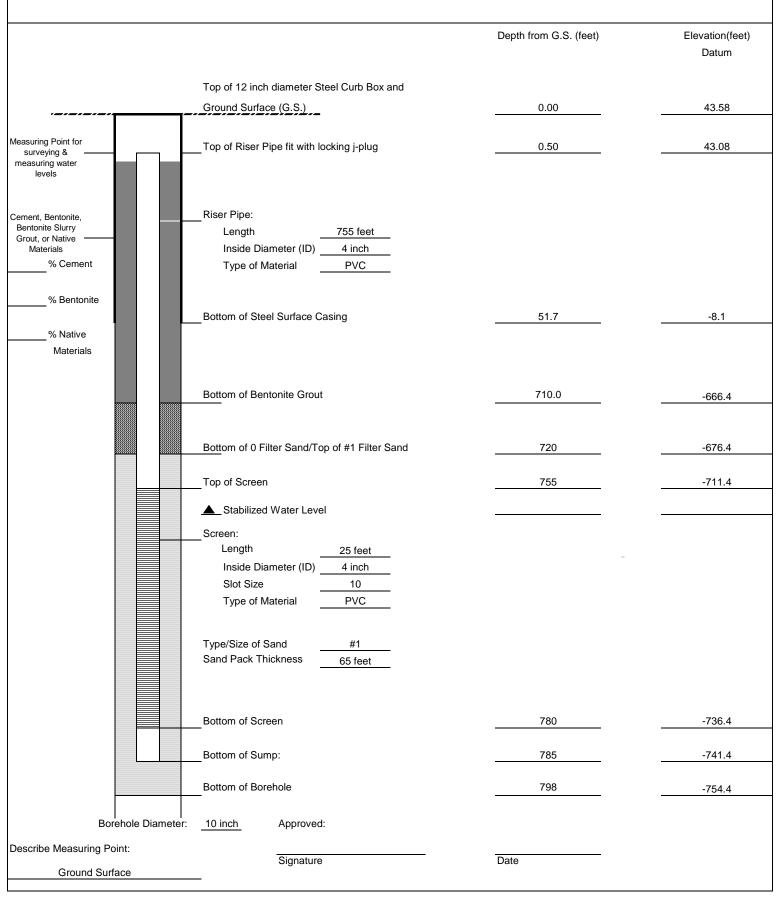


Client: Site Locati		Project Number:	60266526	WELI	LID: BPOW6-1
	ion: Bayberry Lane and Sy	, ,	NY	Date Installed:	8/21/2014 - 9/15/14
Method:	MUD ROTARY			Inspector:	V. THAYER
Coords:	Northing: 193228 16	Fasting: 1126796.9	2	Contractor	DELTA WELL & PLIMP



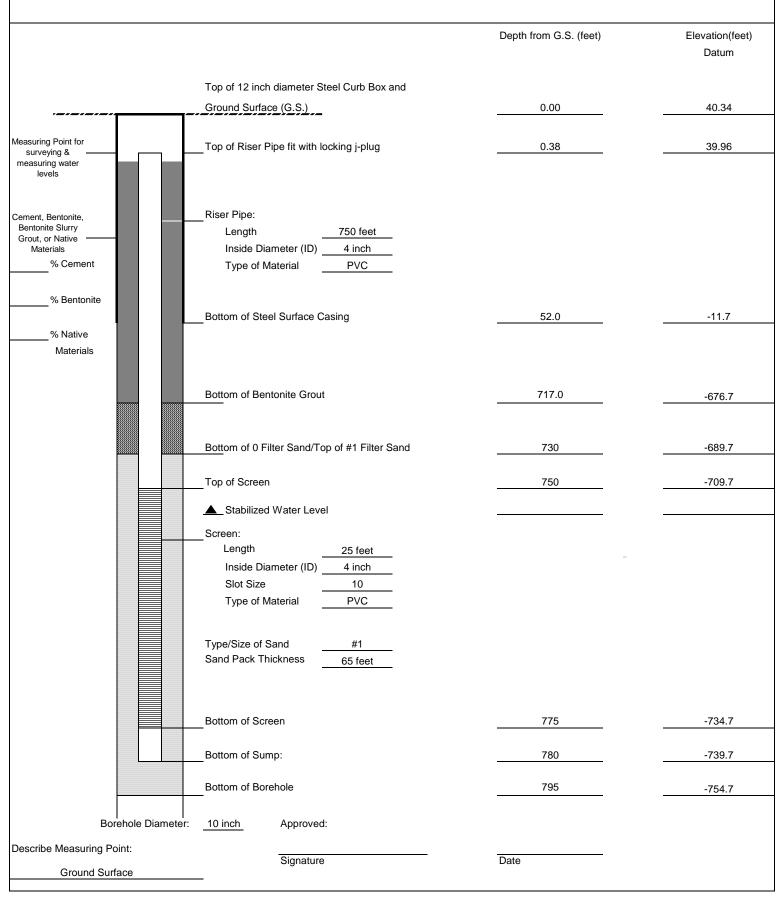


Client: Site Location	NAVFAC n: NWIRP BETHPAG	Project Number: 602665 E, NY	526	WELL ID: BPOW6-2
Well Locati	on: Bayberry Lane and Sy	lvia Road, Seaford, NY		Date Installed: 7/25/2014 - 8/18/14
Method:	MUD ROTARY			Inspector: V. THAYER
Coords:	Northing: 193253.75	Easting: 1126785.34		Contractor: DELTA WELL & PUMP



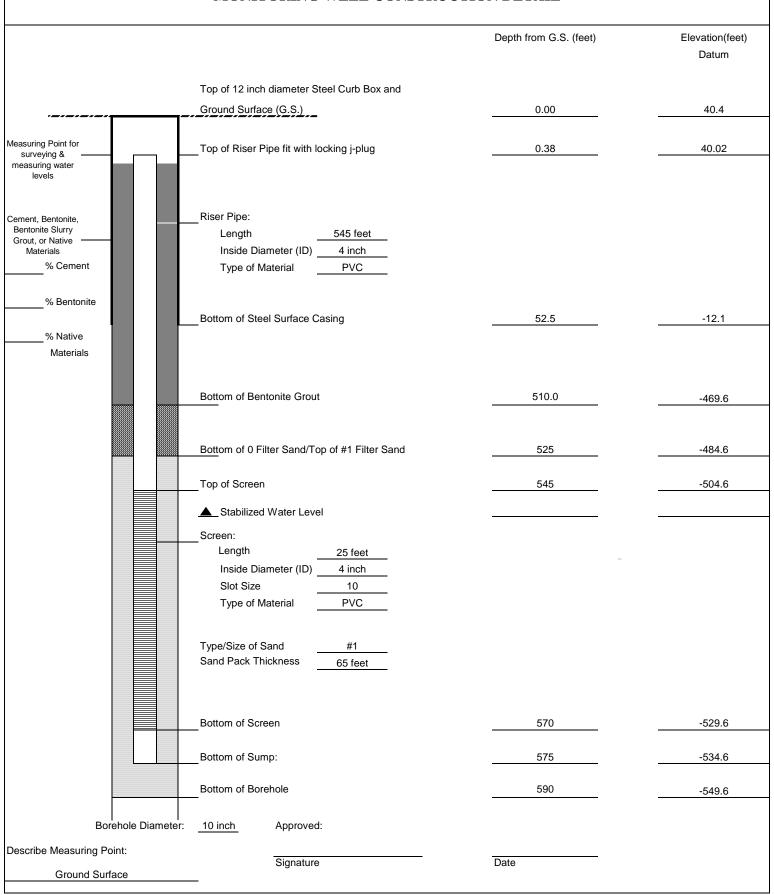


Client:	NAVFAC on: NWIRP BETHPAC	Project Number:	60266526	WELL	LID: BPOW6-3
	ion: Jerome St and Wicks			Date Installed:	10/28/2014 - 11/25/14
Method:	MUD ROTARY			Inspector:	V. THAYER
Coords:	Northing: 19/106.07	Easting: 1127582 3/	5	Contractor	DELTA WELL & DUMP



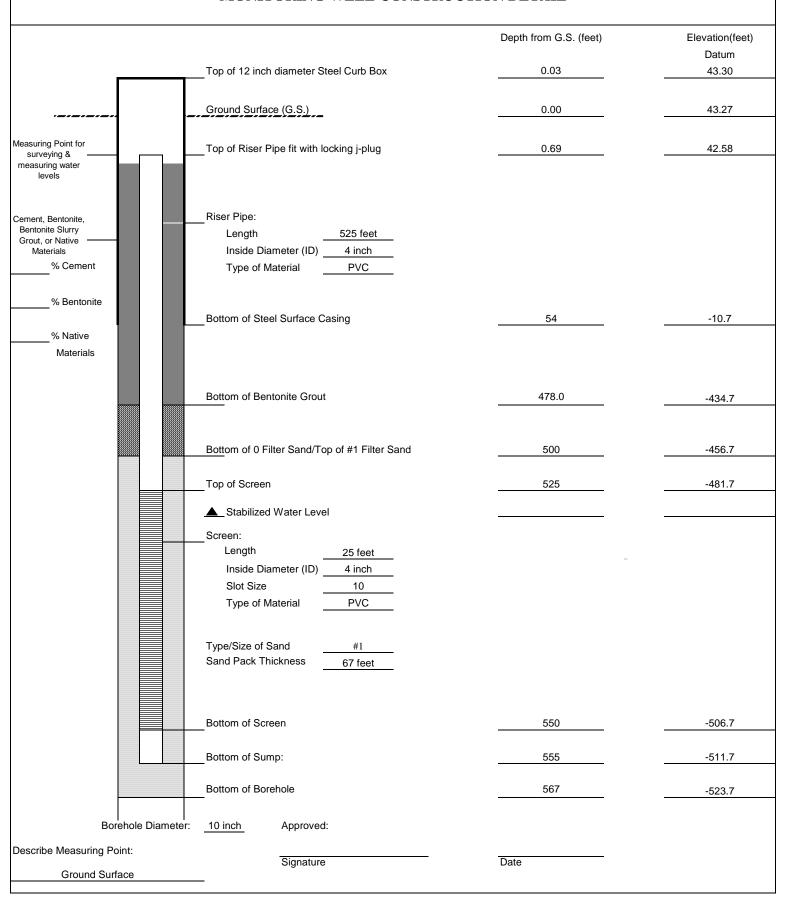


Client: NAVFAC	Project Number: 602665	WELL ID: BPOW6-4
Site Location: NWIRP BET	THPAGE, NY	
Well Location: Jerome St and	Wicks Ave, Seaford, NY	Date Installed: 12/3/2014 - 12/16/14
Method: MUD ROTARY		Inspector: V. THAYER
Coords: Northing: 19412	7.00 Fasting: 1127580.66	Contractor: DELTA WELL & PLIMP





Client:	NAVFAC	Project Number:	60266526	WELL	ID: BPOW6-5
Site Locati	on: NWIRP BETHPAC	SE, NY		,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,	2. 21 0 110 0
Well Locat	ion: North Hickory, Mass	apequa, NY		Date Installed:	5/1/2015 - 5/8/2015
Method:	MUD ROTARY			Inspector: V	/. VARRICCHIO
Coords:	Northing: 194045.49	Easting: 1128178.2	28	Contractor:	DELTA WELL & PUMP





Client:	NAVFAC	Project Number:	60266526	WELL ID: BPOW6-6
Site Locati	on: NWIRP BETHPAC	SE, NY		
Well Loca	tion: North Hickory, Mass	apequa, NY		Date Installed: 5/15/2015 - 5/27/2015
Method:	MUD ROTARY			Inspector: S. WRIGHT/P. KARETH
Coords:	Northing: 194044 59	Fasting: 1128163.7	22	Contractor: DELTA WELL & PLIMP

			Depth from G.S. (feet)	Elevation(feet) Datum
		Ground Surface (G.S.)	0.00	43.17
		Top of 12 inch diameter Steel Curb Box	0.01	43.16
easuring Point for surveying & ———————————————————————————————————		Top of Riser Pipe fit with locking j-plug	0.83	42.34
ement, Bentonite,		Riser Pipe:		
dentonite Slurry Grout, or Native Materials Cement		Length 770 feet Inside Diameter (ID) 4 inch Type of Material PVC		
% Bentonite % Native		Bottom of Steel Surface Casing	54	-10.8
Materials		Bottom of Bentonite Grout	725.0	-681.8
		Bottom of 0 Filter Sand/Top of #1 Filter Sand	745	-701.8
		Top of Screen	770	-726.8
		Stabilized Water Level		
		Screen:		
		Length 25 feet	_	
		Inside Diameter (ID) 4 inch		
		Slot Size 10 Type of Material PVC		
		Type/Size of Sand #1 Sand Pack Thickness 67 feet		
		Bottom of Screen	795	-751.8
		Bottom of Sump:	800	-756.8
		Bottom of Borehole	812	-768.8
Boreh	ole Diameter:	10 inch Approved:		
escribe Measuring Point	:	Signatura	Data	
Ground Surfac		Signature	Date	

Section 3

BPOW6-5 and BPOW6-6 Abandonment



To: Navy CLEAN Bethpage Project File

From: Bill Spronz and Valerie Thayer, Resolution Consultants

Subject: Bethpage Well BPOW 6-5 and Boring BPOW 6-6 Decommissioning

Procedure

Date: 10 April 2015

Introduction

On Friday, February 27, 2015, while drilling the borehole for planned observation well BPOW 6-6, pieces of Schedule 80 PVC well casing were observed mixed with the cuttings in the drilling mud. Investigation by the Delta Well and Pump Co. Inc. ("Delta") crew and the Resolution Consultants ("Resolution") geologist determined that borehole BPOW 6-6 had intersected the borehole and casing of nearby observation well BPOW 6-5 located about 13 feet away. Well BPOW 6-5 and borehole BPOW 6-6 were being installed at vertical profile boring location VPB 147 on Crocus Street in Seaford, New York. The purpose of this memorandum is to present the events that occurred and how the situation was resolved by the grouting and decommissioning of both boreholes.

Background:

Monitoring well BPOW 6-5 was completed to a depth of 795 feet below ground surface ("bgs") on February 11, 2015. The well was constructed of 4-inch diameter Schedule 80 PVC well screen installed from 770 to 795 feet bgs and 4-inch diameter Schedule 80 PVC casing installed from 770 feet bgs to the ground surface. The screen was packed with well sand, the casing tremie grouted to the ground surface, and a temporary PVC cap was installed at the top of the casing.

The BPOW 6-6 borehole was started on February 25, 2015. The planned depth for proposed well BPOW 6-6 was 550 feet bgs. Two days later, on February 27, 2015, Schedule 80 PVC casing was observed in the mud tub while drilling at about 460 to 480 feet bgs. When the crew opened up nearby well BPOW 6-5, water was flowing from the well casing suggesting that the well casing had been breached. The difference in hydrostatic head between the drilling mud in borehole BPOW 6-6 and the water in well BPOW 6-5 was causing water to flow from the well as the fluids equalized.

The crew measured the depth of borehole BPOW 6-6 at 450 feet bgs, and then reamed the borehole to 478 feet bgs to clean it out. Delta then set temporary plugs inside of the 10-inch surface casing and the mud return line welded to the top of the surface casing at boring BPOW 6-6 to prevent fluids from flowing from the borehole while investigating the condition of nearby well BPOW 6-5.

Resolution Consultants and the Navy assessed the situation and concluded that well BPOW 6-5 could not be salvaged. Therefore, the decision was made to decommission well BPOW 6-5 and borehole BPOW 6-6.

Planning:

On Monday, March 2, 2015, a conference call was held with Resolution and Delta personnel to discuss an approach to properly decommission well BPOW 6-6 and borehole BPOW 6-5. The goal for the well and borehole abandonment procedure was to tremie grout both the well and the boring from the total depth of each borehole to the ground surface to prevent potential contaminants of concern (COCs) from entering the borehole or the well and migrating vertically between subsurface zones. The following procedures were proposed:

- Set a temporary packer or plug in the surface casing of boring BPOW 6-5 to prevent fluids from being pushed up into this boring and from spilling into the street.
- Move the drilling rig off boring 6-5 and onto well BPOW 6-6.
- Confirm the depth of the breach and/or sediment blockage in the BPOW 6-6 well casing by dropping a weighted line into the well to determine whether debris or sediments blocked the casing.
- Install small diameter drill rods and drill bit into well BPOW 6-6 to drill or jet through potential obstructions in order to open the well to the total depth of 795 feet if possible.
- If the drill pipe or a tremie pipe could be advanced the total 795 foot depth of well BPOW 6-6, tremie grout the well from the bottom to the ground surface.
- If the drill pipe could not bypass the casing breach or debris fallen into the casing, tremie grout the well from the breach or blocked casing to the ground surface. At a minimum, the grout would be placed to form a seal from the breach in the casing to the ground surface.
- The 4-inch PVC well casing in BPOW 6-5 was already grouted in-place when the well was installed. Therefore, the formation outside of the casing is isolated from the inside of the well preventing vertical migration of potential COCs. Grouting the casing interior or at least from the casing breach to the ground surface in well BPOW 6-5 and fully grouting borehole BPOW 6-6 would complete the integrity of this seal.
- Once the level of grout was equal to the casing breach, the temporary plug in boring BPOW 6-5 would be removed and both borings would be tremie grouted the casing breach to the ground surface through the tremie pipe installed in well BPOW 6-5.
- Confirm that grout is present in both boreholes to the ground surface by monitoring returns at the mud pits and at the top of each borehole.
- Allow the grout to settle overnight and then top out both boreholes.
- Finish decommissioning each borehole by cutting off the casings below ground surface, filling the holes to grade, and restoring the grass in the tree lawn where the well and boring were located.

Implementation:

A drilling rig was set up over well BPOW 6-5 and on March 6, 2015, the depth of the breach in the casing was confirmed at 451 feet bgs using a weighted wire-line lowered into the casing. The drill crew was eventually able to by-pass this area and lowered the weighted wire-line to a depth of 470' within the PVC well casing where sediments and drilling fluid had fallen into well BPOW 6-5 preventing the weighted line from going any further.

A drill bit was attached to 1.25-inch drill pipe and lowered into well BPOW 6-5 to clean the debris out of the well casing to the extent possible so that the grout seal could be tremied from inside of the well casing below the breach to the ground surface. Delta was able to advance the drill bit past the casing breach at 451 feet bgs and felt that they were inside of the BPOW 6-5 PVC well casing based upon drilling conditions. The crew had difficulty maintaining circulation of the drilling mud once they passed the casing breach at 451 feet bgs and had to curtail cleaning the pipe at about 480 feet bgs because drill cuttings and mud were no longer returning to the surface and increased pressure in borehole BPOW 6-6 was causing the temporary plugs sealing the borehole to leak.

Delta was able to remove the drill pipe and install a tremie pipe to a depth of 470 feet bgs in well BPOW 6-5 on the morning of March 7, 2015. The temporary plugs in boring BPOW 6-6 were removed and Cetco high solids bentonite grout was pumped through the tremie pipe from 470 feet bgs to seal both boreholes. Grout returns were identified at ground surface in boring BPOW 6-6 at 12:50 PM and in well BPOW 6-5 at 1:15 PM on March 7, 2015.

On March 9, 2015, the grout seal was measured at 7 feet bgs in well BPOW 6-5 and at 55 feet bgs in boring BPOW 6-6. The Delta crew then tremied additional grout into the well and the boring to seal each to the ground surface. The temporary PVC well cap was removed from well BPOW 6-5 and the casings for well BPOW 6-5 and boring BPOW 6-6 were cut off below the ground surface. The grass strip where the two borings are located was restored to pre drilling condition on April 9, 2015.

Section 4

Groundwater Sample Log Sheets

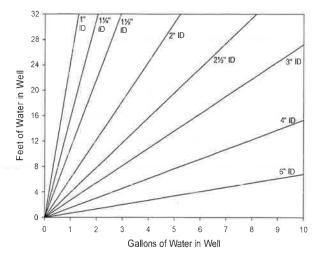


Well ID: BPOW6-1

Low Flow Ground Water Sample Collection Record

Client: Project No	Project No: 60266526						126 11	5 Tin	ne: Start <u> </u>	
Site Locat Weather (tion: Conds:	Cloud	ia by 40	o overaight	vaile (Collector(s)	Pau	1 Kant	H	35.
1. WATE a. Tota b. Wat 2. WELL a. Pur	R LEVEL al Well Len ter Table D PURGE D ge Method	DATA: ngth <u>ら</u> Depth <u>/</u> DATA : Ge	(measu 80 3.53 otech b	c. Length of \d. Calculated	of Casing Water Colo System V	g) umn /olume (see	(a-b)		Casing Diam	neter/Material th PVC
- Tem _l - pH - Sp _{.,} 0	perature	± 3 ± 0 ± 3	% 0.1 unit 9%	(see workplan) - D.O ORP - Drawdown	± 10% ± 10m < 0.3'		Re	Turbidi move a mir	nimum 1 scre	en volume
c. Field	d Testing E	=quipm	ent use -	у.	ake 57		Model 536		55	Number 474
Time	Volume Removed	Temp.	- р <u>Н</u>	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)		Flow Rate (ml/min)		Color/Odor
(24hr)	(Liters)			(ms/cm)	(mg/L)	(1117)	(1410)	(IIII/IIIII)	water (it)	2 · 1
200	- 1	in it is	5.06	0.073	6.61	196,9	13.62	3 30	1362	
and		12.41	4.79	0.075	4.16	720.2	19.5	(41)	3500	
115		12.68	5.62	0.105	2.60	1376				
920		12.72	5.74	1.113	1.82	103.4	A	16.1	13.58	
925		12.70	5,65	0106	1.97	86.0	(*) j.k	- fr 15	•	gray, cloudy
930	5acc	12.68		0.125	2.19	87.1	7/000	tilk hi X		Wisible silt
Ha: Ha:	ceptance of serequired serequired ve parame If no or Note that the control of the	volume turbidit ters sta /A - Exp	e been r y been i abilized	emoved reached low.	Yes No C	urge (k	it bold	an Aux	vest up se	(continued on back)
3. SAMP	LE COLLE	ECTION	N:	Method: Geo	tech blade	der pump v	vith drop tu	be assemb	ly	
Sample II) 			Container Type 40-mL vial 1-L amber	e No. of	Containers 3 2	s Prese HC no		Analysis Red VOCs ,4-Dioxane	Time
Comment	ts <u>// // // // // // // // // // // // //</u>	7 60	Hom	, tubry t	oo long	1				
Signature		Pau	l Ka	ee H				Date	3/26/	4

Purge Volume Calculation



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

1 screen volume 15 ft = 56.8 L / 9.8 G 20 ft = 75.7 L / 13.1 G 25 ft = 94.6 L / 16.3 G

Well ID:	800005-1	3	850 =	Zhrs

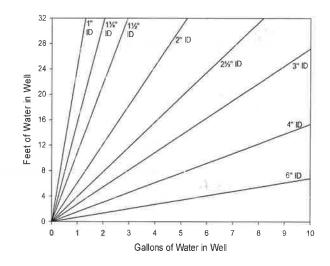
(continued										
	Volume				士10%					
Time (24 hr)	Removed (Liters)	Temp (°C)	ρН	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
935	(Encra)	12.61	5.60	0.102	(mg/L)	69.6	>1000	500	water (it)	
940		12.71	5.81	0.100	2.25	62.7	21/10	320		
645		12.75	3.01 453	0.098	2.00	581	7/000		13.60	
950		12.72	4150	0.097	1.41	57.5	71000		73.5	clear flowcell.
955		12.68	5.45	0.094	1.58	244	7/100	500		TIPLOCETT
1000		12.62	538	0.092	150	33.0	905		13.58	clear flowcell
1005	1000	12.72	5.76	0.090	1.20	57.0	850		30100	110000011
1010	0	12.74	5.81	0.091	1.16	548	786		A	744
1015		H.69	577	0.089	1.29	JJ .0	576	177	fy. I	
1020		12.74	7.74	0.089	1.31	55.5	511		13.6/	
1025		12.76	15,24	5.088	1.34	63.6	443			
1030		12.81	5.19	0.089	1.00	58.8	37.7			111
1035	1590	12.90	5.16	0.08%	1.16	59.7	265			
1040	-	12.88	5.14	0.088	1.15	60.0	243		13,60	
1015		12.97	5.12	0.088	1.14	60.6	162			
1050	17.5gal	1294	5.10	0.088	1.17.	60.8	163		13,60	
	/							Soci		
1100										Suple
							2.42	epre.		



Low Flow Ground Water Sample Collection Record

Client:	Navy NW	IRP Re	thnage		Г	ate: 3	126 11	5 Tir	me: Start _	25 Á pm
Client: Navy NWIRP Bethpage Project No: 60266526							1 1	- '"	Finish	
Site Locat	ion:	54/4	ia							
Weather (Conds:	385	,	eloxdy		Collector(s)		<u></u>		
1. WATE	R LEVEL I	DATA:	(measu	red from Top	of Casin	g)				
a. Total Well Length 785 c. Length of Water Column 13.96 (a-b) Casing Diameter/Material										
										ch PVC
b. Wat	er Table D	epth_		d. Calculated	System \	/olume (see	back)	5 ft scr	een , 16.	3 gal
	PURGE D		-4b bl		ith drop to	iba aaaamh	als e			
a. Purç	ge Method	: <u>Ge</u>	otech bi	adder pump w	ith drop it	ibe assemi	лу			
	•			see workplan)		(values >0) E ma/l)	Tuebid	: 100	/
- I emp - pH	perature		% 1 unit	- D.O. - ORP	± 10%	տ (values >0 ւ∨	J.5 mg/L)	i urbia	ity ± 10%	o .
	Cond.			- Drawdown	< 0.3'	•	Re	move a mi	nimum 1 scre	en volume
c Field	d Testing E	- -auinm	ent used	d: Ma	ıke		Model		Seria	l Number
0.1101	u 100mg	-qa.p	-	V51			5560	PS	711	
	\		=	1anna			H1 987	03		
<u>Time</u>	Volume Removed	Temn	<u>рН</u> -	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)	
845		10.80	7.85	0.054	7.70	-42.3		500	13.79	cloudy
855		12.71	5.56	0.044	2.30	-72.3	26.4	600	13.99	clusely
900		12.84	5.51	0.041	2.07	-26.2		600	13.99	"lundy
905	Syal	12.87	5.69	0.045	1.75	-27.5	1000	600	13.99	cloudy
910		12.82	5,66	0.042	1.62	-37:0		600	13.99	cloudy
915		12.77	5,54	0.037	1.51	-38,4	770	600	13.99	cloudy
920		12.87	5.50	0.036	1.45	-38.6		600	17.98	elovely
	ceptance c				Yes N	_	1	fic		(continued on back)
	s required s required									
	ve parame		•	odonod		i i				
	If no or N/	A - Exp	lain bel	ow.						
	! 									
3. SAMP	LE COLLE	CTION	l : I	Method: Geo	tech blad	der pump w	ith drop tu	be assemb	oly	
Sample IE)		(Container Type	e No of	· Containers	s Prese	rvation	Analysis Re	q. Time
	2 - GIW	0326		40-mL vial	7 140. 01	3	HC		VOCs	1025
BPOW 6-	Z-62W -	0326	15	1-L amber		2	no	ne	1,4-Dioxane	1025
	4.4	.1 1	4.1							
Comment	s H	+ >0	Hom							- 1
Cianatura	(1					Date	3/76	12015
Signature			1			0			vFlow-GWa - rev	- 1

Purge Volume Calculation



	Volume /	Linear F	t. of Pipe
	ID (in)	Gallon	Liter
ì	0.25	0.0025	0.0097
ļ	0.375	0.0057	0.0217
	0.5	0.0102	0.0386
ľ	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
þ	4	0.6528	2.4711
	6	1.4688	5.5600

1 screen volume 15 ft = 56.8 L / 9.8 G 20 ft = 75.7 L / 13.1 G 25 ft = 94.6 L / 16.3 G

Well ID:

(continued	-	,									
Ti	Volume	T		0 0 1	50	000	-				
Time (24 hr)	Removed (Liters)	Temp (°C)	pН	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/O	dor
925	(Enters)	12.86	5.44	0.036	1.41	-39.1	265	600	13.99	Cloudy	
930		12.52	5.44	0.035	1.69	-32.6	244	600	17.99	Cloudy Cloudy Cloudy	
935		12.92	5.44	0.034	1.77	-75.0		600	17.99	Cloudy	
940	[Vegal	12.90		0.034	1.34	-75.1	134	600	17.77	Cloudy	
945		12.41	5.79	0.033	1.32	-378	82.1	600	13.97	Cloudy	clegrin
950		17.89	5.38	0.037	1.70	-32.5	69.0	600	13.47	-1	7
955		12.88	5.37	0.077	1.27	-30.7	56.4	640	13.97	& e	
1000		12.85	5.36	0.032	1.24	-28.5	52.8	600	13.97	11	
1005		12.88	5.36	0.032	1.23	-27.5	38.4	600	17.97	e,	
1010	15 341	12.89	5.35	DIOJZ	1.23	-76.2	33.1	600	13.97	4,0	
1015		12.93		0.032	1.70	-24.3	29.9	600	17.97	f ₄	
1020		12.48	\$.37	0.072	1,18	-22.5	30,4	600	13.47	Eq.	
						ie.					
F.1											
	a										
										1 -8	

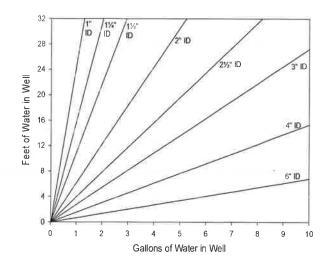




Low Flow Ground Water Sample Collection Record

Client: Project No Site Locat			6026	66526 1 Seromo		-	126 11		Finish	am/pm 3/5 am/pm
Weather 0	Conds:	40 "		rle	c	collector(s)	_ /-	al Have	and a second	
a. Tota b. Wat 2. WELL a. Puro b. Acco	al Well Lenter Table Description Descripti	gth_7 Pepth_ ATA : Ge	%, 78 otech bl	c. Length of \ d. Calculated adder pump w see workplan)	of Casing Water Colu System V	olume (see	(a-b)	=5 ft su	Casing Diamo 4-incl CECN 16.	1 gal
- Temp - pH	perature		5% 0.1 unit	- D.O. - ORP	± 10% ± 10m).5 mg/L)	Turbidit	ty ± 10%	
-	Cond.		%				Re	move a mir	nimum 1 scree	en volume
c. Field	d Testing E	Equipm	ent used	d: Ma YS Has		-4 ×	Model 556	S (0)	71	Number 477 777
	Volume		-			7/	V	5 (8
<u>Time</u> (24hr)	Removed (Liters)	Temp.	<u>PH</u>	Spec. Cond. (mS/cm)	<u>DO</u> (mg/L)	<u>ORP</u> (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
1300	6 N				7-	- 1	A		3. %	
1325		1258	5.24	0.032	2.01	177.8		550	10.65	Can ly
1350		1243	5.02	0.028	1.78	1920	(3)	7	* = 2 t	32/
1335	5 gal	12.13	5.00	00028	461	195,2	27,3			
1340		12,13	4.96	0.027	1.42	149.6	181	y V	10.62	
1345		/Z35	4.96	0.027	1.20	202,9	46.5	-		
/350			4.95	0.027	1.15	204.5	17.5			
Ha: Ha:	ceptance c s required s required ve parame If no or N/	volume turbidit ters sta	e been re y been r abilized	emoved eached	Yes No					(continued on back)
3. SAMP	LE COLLE	ECTION	N: 1	Method: Geo	otech blado	ler pump w	vith drop tu	be assemb	ly	_
Sample II	D 16-3-44	V-033		Container Type 40-mL vial	No. of	Containers	НС		Analysis Req	Time
				1-L amber		2	no	ne 1	,4-Dioxane	
Commen	ts _ 45 	ho	Homa .	stirredy	y 34/1-,	pra ci	to teching	g is to l	ng	
Signature)	/	Paul	Kart				Date	3/26/ Flow-GWa - rev	March 2015.xls

Purge Volume Calculation



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

1 screen volume

15 ft = 56.8 L / 9.8 G 20 ft = 75.7 L / 13.1 G 25 ft = 94.6 L / 16.3 G

Well ID:

(continued	from front)									
	Volume									
Time	Removed	Temp	рН	Spec. Cond.	DO	ORP	Turbidity	Flow Rate	Depth to	Color/Odor
(24 hr)	(Liters)	(°C)		(mS/cm)	(mg/L)	(mV)	(NTU)	(ml/min)	water (ft)	
1355		12.78	4.95	0.027	1.10	207.4	30.6	530	10.68	
1400		1233	4.95	0.027	1.07	208.5	17.0			
1415	10001	(250)	11-921	0.027	1.07	210,2	14.0			
1410		12.47	4.94	0.027	1.04	211.3	11.1			
1415		12.44	4.94	0.077	1.01	212.4	658			
1420		12.31	4.94	0.027	098	215.5	5.86			
1425		12.32	4.94	0.027	0.2%	2144	5.85			
1930	1500	12.25	4.94	0.027	0.95	2153	6.02	55%)	10.78	
1935	7	12.60	494	0.027	092	2/7.3	4,57	A-1		
1440		1267	4.94	0.027	092	217.9	4:28	200	- 3 PeV	
444-5	170al	12,86	494	0.027	0.98	2/8.4	4.47	756 ×	UN.	2
	,						0.08	V 5 X 3	Ac 4.	
1450										ale Dil
141										
										ř
				+						



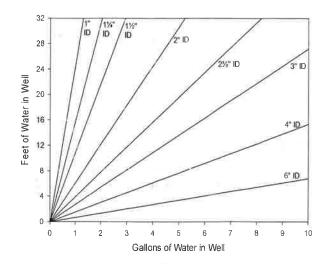
Well ID: 8 POW6-4

LowFlow-GWa - rev March 2015.xls

Low Flow Ground Water Sample Collection Record

Client:	Navy NW	IRP Be			D:	ate: 3	126 11	5 Tir		320 am/pm
Project No	_			66526					Finish	am/pm
Site Locat			- 1	1 eron		N 11 (/ X		JC		
Weather (Conds:		dri-	رياح.		collector(s)	!	<u> </u>		
1. WATE	R LEVEL	DATA:	(meası	red from Top	of Casing	1) 565	70°			
a. Tota	al Well Len	<u>ک</u> igth	75	c. Length of \	Water Colu	ımn_ **	3 (a-b)		Casing Diame	eter/Material n PVC
b. Wat	ter Table D	epth_	9.92	d. Calculated	System V	olume (see	back) 2	set sur	eau 16.	laal
	PURGE D								-	
			otech b	adder pump w	ith drop tu	be assemb	oly			
h Acc	entance C	ritoria d	lefined (see workplan)						
	perature	± 3		- D.O.		(values >0).5 mg/L)	Turbidi	ty ± 10%	
- pH	p 0. 0. 0.		.1 unit	- ORP	± 10m	•	3 – ,			
- Sp. C	Cond.	± 3	%	- Drawdown	< 0.3'		Re	move a mi	nimum 1 scree	en volume
c Field	d Testing E	-auipm	ent use	d: Ma	ake		Model		Serial I	Number
		-4		V5"			56mm		711	
				11-10114		l t	1198703	Į.	691	77
т	Volume	-		0		000	T	Flam Data	D 41- 4-	0-1/0-1
<u>Time</u> (24hr)	Removed (Liters)	(°C)	<u>pH</u>	Spec. Cond. (mS/cm)	<u>DO</u> (mg/L)	<u>ORP</u> (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
1339	(Eitoro)		4.12	0.105	4.31	15-9	(,,,,,,	750	9.98	(buy)
1375		1281		0.02	4.14	10.5		750	10.03	cley
1340			4.14	0.101	7.01	70.0		750	10.07	claus
1345		12.71		0.093	1.36	5,t	76:2	758	10.09	clear
1350		12.64		0.093	1.27	-9.4	20	750	10.09	chier
1355		12.59		0.097	1.14	-15.8	6.71	750	10.09	cher
1400		12,62	4.57	0.092	1.10	18.2		750	10.09	chin
d. Acc	ceptance c				Yes No	N/A	4			(continued on back)
	s required									
	s required		•	eached						
па	ve parame If no or N/			ΟW		L				
			,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,	· · ·						
		-07101				•	101 1			
3. SAMP	LE COLLE	CHON	N: I	Method: Geo	tech bladd	er pump v	vith drop tu	be assemb	oly	-
Sample II)		(Container Type	No. of	Containers	s Preser	rvation	Analysis Req.	Time
BPOW	6-4-GW	-032	615	40-mL vial		3	HC		VOCs	1500
Brow	6-4-GW	-032	615	1-L amber		2	noi	ne ´	1,4-Dioxane	1500
		_								
Comment	ts	_							- 1:	
		1								
		1	9						3/2/	
Signature		1						Date	2/26/	2915

Purge Volume Calculation



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

1 screen volume

15 ft = 50.8 L / 9.8 G 20 ft = 75.7 L / 13.1 G 25 ft = 94.6 L / 16.3 G

Well ID:

(continued f	,									
	Volume									
Time	Removed	Temp	рН	Spec. Cond.	DO	ORP		Flow Rate	Depth to	Color/Odor
(24 hr)	(Liters)	(°C)		(mS/cm)	(mg/L)	(mV)	(NTU)	(ml/min)	water (ft)	
1405		17.80	4.57	6.077	1.0.3	-21.0	3.15	750	10.11	ched
1410		1277	4.57	O. VH	1,04			750	10,11	Clear
1417		12.83	4.77	0.042	1.01	-25.8	254	750	10.11	cher
1420		12.77	4.57	0.042	1.01	-76.4		750	10.11	clear
1425		12.71	4,57	0,04	1.00	-78.3	7.24	750	10.11	Elvar
1430		12.61	4.57	0.092	1.00	-24.8		750	10.11	cleyr
1475		12.64	4.58	0.092	0.98	-70:7	7.63	750	10.11	Clear cher Clear clear clear clear chear chear chear
6PPi		12.95	4.54	Ovaz	0.45	-72.1		750	10.11	Chus
1445		12.90	4.59	0.093	0.41	-3219	7.49	750	10.11	chear
1450		12.91	4,57	0.097	0.91	-34.1		756	10.11	elm
•										

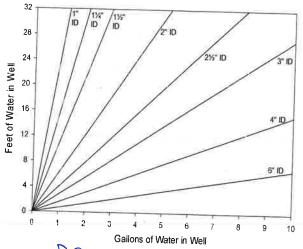


DOWE S					
Well ID: Row 6 5	Well ID:	Rpow	6-	5	

Low Flow Ground Water Sample Collection Record

	avy NWIF	RP Bethpa				Date:	61 25	/15	Time: Sta	O-X	m/pm m/pm
roject No:		0266526							Liii	ISII_IVZ	im biii
ite Locatio	100					0-11	to=(o):				
/eather C	onds:					Coll	ector(s):_				
. WATER	R LEVEL I	DATA: (m ell Length:	easured 1	from Top	of Casing) c. Length of	f Water Colu	umn: 53	7. 60 ft	Casir	ng Diameter/M 4-inch PV	aterial
. t	o. Water T	able Dept	h: <u>[8 00</u>	_ft (d. Calculat	ed System	Volume (se	ee back)			
. WELL	PURGE D a. Purge M	ATA Method: <u>C</u>	Geotech b	ladder pun	np with dro	p tube asse	mbly			· ·	
	o. Accepta - Temper - pH ± 0. - Sp. Con	ature ± 3° 1 unit	%	d (see work - D.O. ± 1 - ORP ± 1 - Drawdow	0% (values 0mV	s >0.5 mg/L))	- Turbidity - Remove	a minimum	n 1 screen volu	
	c. Field Te	esting Equ	ipment us	sed:	Make	5	Model SS6 m	25		Serial Number	
	22			0.	HANN	/A		98703		1240 24X	
	Volume						Turbidity	Flow Rate	Depth to		
Time	Removed	Temp		Sp. Cond.	DO (ma/l)	ORP (mV)	(NTU)	(ml/min)	water (ft)	Color / Od	or
(24hr)	(liters)	(°C)	pH	(mS/cm) (0.67)	(mg/L)	134.6	(4.6	700	18.04	Claruli	non(
1835		6.29	10.99		3.84		11.0	800	(8.05	i,	
1745		15.83	6.31	0-085		150.6 275.2		700	18.02	53	
1855	19	15.76	4.25	0-052	3.18		,		18.06	150	
1905		15.90	4.26	0-053	1.73	279.6	11.5			46	+
1915		15.95	4,25	6-053	1.55	279.>	10.2	700	18.06	377	
0925	38	16.15	4.43	0.053	1-08	975.3	3.49	700	18.02	73.	
0935	70	16.09	4.23	0.053	1.48	265.7	2,69	100	18.00		
W J Z	Has r Has r Have	otance crit	olume bee Irbidity be Irs stabiliz	en removed en reached ed	3	Yes A T T	№	N/A		(continued on back)	
3. SAM	IPLE COL	LECTION	:	Method	Geotech	bladder pur	mp with dro	op tube as	sembly		
Sample	: 1D 5w 6-5-6	m -18622;	3015	<u>Cc</u>	ntainer typ 40-mL via	ls	container	HC	CI	Analysis Req. VOCs 1,4-Dioxane	102
	d				1-L amb	er	2	non	e	1,T DIOXAIIO	1000
Comme	eņts										
		1//	1								

Purge Volume Calculation



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

1 screen volume

15 ft = 37.1 L / 9.8 G 20 ft = 49.6 L / 13.1 G 25 ft = 61.7 L / 16.3 G

Well ID:

Brow 6-5

continuea t	from front)									
	Volume									
Time	Removed	Temp	pН	Spec. Cond.	DO	ORP	Turbidity	Flow Rate	D" ·	
(24 hr)	(Liters)	(°C)		(mS/cm)	(mg/L)	(mV)	(NTU)	(ml/min)	Depth to water (ft)	Color/Odor
0945	(A)	16.00	4.49	0.053	03.0	251.9	2.30	750	18,00	1,
1955	62	16.03	4.46	6.053	0.75		2.14	700		a a
1005		16.25	4.40	0-653	0.69	233.1	2.24	750	18.00	
1010		(6.98		6-053	0.65	239.6	1.65	200		
1015		20-19	4.78	0.054	0.57	220.7	1.80		18.00	
620		2080		0.05-1	0-58	221.1	1.57	260	18.00	G.
				0.03 -1		(0)	1.21	dou	18,00	Clearloone Simpled
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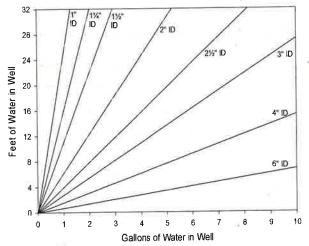


Well ID: BPOW 6-6

Low Flow Ground Water Sample Collection Record

						. .	Pio oc	, nooth	III IXGC	,or a	
Client:		WIRP Bet				Date:	6/ 7	? 5 /15	Time:	Start 810	_ám/pm
Project		602665				=: =:				Finish 1075	
Site Loc			ckory			_					=
vveaine	r Conds:	40	3 F	dias		C	collector(s)	:			
1. WAT	ER LEVE	L DATA:	(measure	ed from To	p of Casin	g)					
			gth:&oc			of Water C	olumn: 🗾	281-47 f	t C	asing Diameter/	
	b. Water	r Table De	epth: 18.	53 ft	d. Calcul	ated Syster	m Volume	(see back)	16.3	4-inch P	<u>VC</u>
2. WEL	L PURGE		-								
	a. Purge	Method:	Geotech	bladder pu	ımp with dr	op tube ass	sembly				
	- Tempe - pH ± (erature ±	3%	ed (see wo - D.O. ± - ORP ± - Drawdo	10% (value 10mV	es >0.5 mg/	L)		ty ± 10% e a minimu	um 1 screen vol	ume
	c. Field 1	resting Ed	quipment ι	used:	Make		Model			Serial Number	r
	Volume										
Time	Removed	Temp		Sp. Cond.	DO	ORP	Turbidity	Flow Rate	Depth to		
(24hr)	(liters)	(°C)	pН	(mS/cm)	(mg/L)	(mV)	(NTU)	(ml/min)	water (ft)	Color / Ode	or
225		16.04	5.05	0.048	1.59	265.6	627	700	18.55	stable ilid,	Inem
838	Saullon		4.63	0.035	1.45	283.6	-	700	18.56	7.	14
845		15,99	4.14	0.079	1.16	315.3	104	700	18.56	Cleanny	- 5
855		16.08	4.27	0.028	0.89	311.9	75.7	700	18.56	5/194Hs do	sede Ino
905	(Ogellons	16.04	4.34	0.028	0.85	311, 9	69.0	700	18.56	Clearlhony	-
915		16,13	4.18	0 028	0.76	320.2	60.7	700	(8.56	U	
925		16.29	4.21	0-628		312	46.3	700	18.56	tc .	
			eria pass/f			Yes	No	N/A	10 0	(continued on back)	
				n removed		<u>*</u>					
	Have	parameter	s stabilize	n reached		[]	片				
			- Explain b			3	ш	Ш			
			_								
B. SAMPI	LE COLL	ECTION:		Method:	Geotech bl	adder pum	p with drop	tube asse	embly		
Sample ID)			Cont	ainer type	No of c	ontainers	Preserv	vation A	nalysis Bos	
	GW-662				0-mL vials	3	Ontainers	HCI	<u>valion A</u>	nalysis Req. VOCs	Time 1015
POW 6-6	-GW-06.	25 2015			1-L amber	2		none		1,4-Dioxane	1015
comments											
	-										
						·					
ignature	p.:						_				
				(

Purge Volume Calculation



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

1 screen volume

15 ft = 37.1 L / 9.8 G 20 ft = 49.6 L / 13.1 G 25 ft = 61.7 L / 16.3 G

Well ID:

BPOW 6-6

(continued f	from front)	000								
Time (24 hr)	Volume Removed (Liters)	Temp (°C)	рН	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
135	15 galle	16-13	4.21	0-005	0-66	324-6	36.0	700	18.56	If
940	16.59	16.78	4.45	0.078	0.56	3/3.9	39-1	700	18.56	fe -
945	100	16.43	4.57	0.078	0.58	313.1	32.9	700	18.56	K
950		16.39	4.52	0.078	0.58	316.0	31.5	700	18.56	4
9.55		16.40	4.26	0.028	6.59	330.5	31.5	700	18.52	/ (
1000		16.46	4.10	4.029	0.61	3307	31.6	700	18-51	Ny
1015		10 10								Sant
1010										
-						Κ.				
										-

Section 5

Analytical Data Validation – Resolution Consultants

BPOW6-1, BPOW6-2, BPOW6-3, BPOW6-4, BPOW6-5, BPOW6-6 (VPB 145, VPB 146, VPB 147) Installation Report BETHPAGE, NY

Documentation of well name change

Well BPOW6-1 was originally named RE111D1

TOC data reported in data validation report package SH7523 contains sample RE111D1-SOIL-090514-573-575 and thus pertains to well BPOW6-1.

Well BPOW6-2 was originally named RE111D2

TOC data reported in data validation report package SH6618 contains sample RE111D2081114-768-770 and thus pertains to well BPOW6-2.



Data Validation Report

Project:	Regional Groundwater Investigation - N	IWIRP Bethpage
Laboratory:	Katahdin Analytical	
Service Request:	SH6618	
Analyses/Method:	EPA SW-846 Method 9060A for TOC	
Validation Level:	3	
AECOM Project Number:	60266526.SA.DV	
Prepared by:	Dawn Brule/RESCON	Completed on: 01/06/2015
Reviewed by:	Lori Herberich/RESCON	File Name: SH6618_9060A

SUMMARY

The samples listed below were collected by Resolution Consultants from the Regional Groundwater Investigation - NWIRP Bethpage site on August 11, 2014.

Sample ID	Matrix/Sample Type
RE111D2081114-768-770	Soil

Data validation activities were conducted with reference to these methods, *Test Methods for Evaluating Solid Waste, Physical/Chemical Methods SW-846*, Method 9060A, *Total Organic Carbon* (USEPA, 1996), *USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Superfund Data Review* (January 2010), and Quality Systems Manual (QSM) for Environmental Laboratories, Version 4.2 (DoD, October 2010) where applicable. 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 [COC])/sample integrity
- ✓ Holding times and sample preservation
- ✓ Initial calibration/continuing calibration verification
- Laboratory blanks/equipment blanks
- NA Matrix spike (MS)/Matrix duplicate (MD) and/or matrix spike duplicate (MSD) results
- ✓ Laboratory control sample (LCS) results
- NA Field duplicates
- ✓ 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. The symbol (X) indicates that a QC nonconformance resulted in the qualification of data. Any QC nonconformance that resulted in the qualification of

2

data is discussed below. In addition, nonconformances or other issues that were noted during validation, but did not result in qualification of data, may be discussed for informational purposes only.

The data appear valid as reported and may be used for decision making purposes. There were no data points qualified or rejected on the basis of this data review.

RESULTS

Data Completeness (COC)/Sample Integrity

The data package was reviewed and found to meet acceptance criteria for completeness:

- The COCs were reviewed for completeness of information relevant to the samples and requested analyses, and for signatures indicating transfer of sample custody.
- The laboratory sample login sheet(s) were reviewed for issues potentially affecting sample integrity, including the condition of sample containers upon receipt at the laboratory.
- Completeness of analyses was verified by comparing the reported results to the COC requests.

Due to limitations in the reporting system, the laboratory omitted the "RE-" prefix from the sample ID in the report. The submitted EDD file reflects the full sample ID.

Holding Times/Sample Preservation

Sample preservation and preparation/analysis holding times were reviewed for conformance with the QC acceptance criteria. The QC acceptance criteria were met.

Initial Calibration/Continuing Calibration Verification

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

- all criteria were met for the calibration curves
- the initial calibration verification (ICV) percent recovery (%R) criteria were met; and
- the continuing calibration verification standard (CCV) method %Rs were met

The QC acceptance criteria were met.

Laboratory Blanks/Equipment Blanks

Laboratory method blanks and equipment rinsate blanks were evaluated as to whether there were contaminants detected above the detection limit (DL). An equipment blank was not submitted with the samples in this data set.

Data validation qualifications for individual samples are based on the maximum contaminant concentration detected in all associated blanks.

Method blank results were reviewed for conformance with the QC acceptance criteria. Detected results in blanks are not discussed in this data validation report if the associated results were nondetect or if qualification of sample results was not required.

The QC acceptance criteria were met and/or qualification of the sample results was not required.

MS/MD and/or MSD Results

MS/MD/MSD analyses were not performed on samples reported in this SDG. There were no validation actions taken on this basis.

LCS Results

The LCS %Rs were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

Field Duplicate Results

There were no field duplicate samples submitted with this data set. No validation actions were taken on this basis.

Sample Results/Reporting Issues

Compounds that were not detected in the sample are reported as not detected (U) at the Limit of Detection (LOD).

Compounds detected at concentrations less than the LOQ but greater than the detection limit (DL) were qualified by the laboratory as estimated (J). This "J" qualifier was retained during data validation.

Any sample that was analyzed at a dilution due to high concentrations of target or non-target compounds or matrix interferences was checked to ensure that the results and/or sample specific LODs and LOQs were adjusted accordingly by the laboratory.

QUALIFICATION ACTIONS

No sample results were qualified as a result of this data review.



600 Technology Way Scarborough, ME 04074 Tel: (207) 874-2400 Fax: (207) 775-4029

CHAIN of CUSTODY

PLEASE BEAR DOWN AND PRINT LEGIBLY IN PEN

Page / A

Resolution	V Consultan	its	Contac	it Læno:	a ViV	audou	Phone #)		. F	ax #		
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Purchase Order #	Proj.	. Name / N	10. Bel	Yh pu	ne 6	02.66	521	1	2	in Quote		//	
Bill (if different than above)				dress				<i>a</i> '	,		~		-
Sampler (Print / Sign)	erie Thay	13	V	aleu	e The	alex		Cop	ies To:	Value	ix Tha	LUN	
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Report of Analytical Results

Client: Rick Purdy AECOM

701 Edgewater Drive

Wakefield,MA 01880

Lab Sample ID: SH6618-2

Report Date: 04-SEP-14 Client PO: 16518

Project: Navy Clean WE15-03-0 SDG: SH6618

Sample Description 111D2081114-768-770	<u>1</u> 077	*	000			Matrix SL	Date Sampled 11-AUG-14		Date Received 9-AUG-14		
rameter	Result Adj	Adj LOQ	Adj MDL	Adj LOD	LOQ Adj MDL Adj LOD Anal. Method QC, Batch	QC.Batch	Anal, Date	Prep. Method Prep. Date Footnotes	Prep. Date	Footnotes	
AC In Soil	920 ug/gdrywt	440	45.	330	SW846 9060A Mod.	WG148851	SW846 9060A WG148851 22-AUG-14 10:39:30 Mod.	N/A	N/A		1
tal Solids	% 06	_		N/A	SM2540G		WG148631 21-AUG-14 10:56:13	SM2540G	20-AUG-14		

TOC In Soil Parameter

Total Solids



Data Validation Report

Project:	Regional Groundwater Investigation - N	IWIRP Bethpage
Laboratory:	Katahdin Analytical	
Service Request:	SH7523	
Analyses/Method:	EPA SW-846 Method 9060A for TOC	
Validation Level:	3	
AECOM Project Number:	60266526.SA.DV	- 7
Prepared by:	Dawn Brule/RESCON	Completed on: 01/06/2015
Reviewed by:	Lori Herberich/RESCON	File Name: SH7523_9060A

SUMMARY

The samples listed below were collected by Resolution Consultants from the Regional Groundwater Investigation - NWIRP Bethpage site on September 5, 2014.

Sample ID	Matrix/Sample Type
RE111D1-SOIL-090514-573-575	Soil

Data validation activities were conducted with reference to these methods, *Test Methods for Evaluating Solid Waste, Physical/Chemical Methods SW-846*, Method 9060A, *Total Organic Carbon* (USEPA, 1996), *USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Superfund Data Review* (January 2010), and Quality Systems Manual (QSM) for Environmental Laboratories, Version 4.2 (DoD, October 2010) where applicable. 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 review elements (where applicable to the method):

- ✓ Data completeness (chain-of-custody [COC])/sample integrity
- ✓ Holding times and sample preservation
- ✓ Initial calibration/continuing calibration verification
- ✓ Laboratory blanks/equipment blanks
- NA Matrix spike (MS)/Matrix duplicate (MD) and/or matrix spike duplicate (MSD) results
- ✓ Laboratory control sample (LCS)/laboratory control sample duplicate (LCSD) results
- NA Field duplicates
- ✓ 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. The symbol (X) indicates that a QC nonconformance resulted in the qualification of data. Any QC nonconformance that resulted in the qualification of

Resolution Consultants 2

data is discussed below. In addition, nonconformances or other issues that were noted during validation, but did not result in qualification of data, may be discussed for informational purposes only.

The data appear valid as reported and may be used for decision making purposes. There were no data points qualified or rejected on the basis of this data review.

RESULTS

Data Completeness (COC)/Sample Integrity

The data package was reviewed and found to meet acceptance criteria for completeness:

- The COCs were reviewed for completeness of information relevant to the samples and requested analyses, and for signatures indicating transfer of sample custody.
- The laboratory sample login sheet(s) were reviewed for issues potentially affecting sample integrity, including the condition of sample containers upon receipt at the laboratory.
- Completeness of analyses was verified by comparing the reported results to the COC requests.

Due to limitations in the reporting system, the laboratory omitted the "RE111D1-" prefix from the sample ID in the report. The submitted EDD file reflects the full sample ID.

Holding Times/Sample Preservation

Sample preservation and preparation/analysis holding times were reviewed for conformance with the QC acceptance criteria. The QC acceptance criteria were met.

Initial Calibration/Continuing Calibration Verification

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

- all criteria were met for the calibration curves
- the initial calibration verification (ICV) percent recovery (%R) criteria were met; and
- the continuing calibration verification standard (CCV) method %Rs were met

The QC acceptance criteria were met.

Laboratory Blanks/Equipment Blanks

Laboratory method blanks and equipment rinsate blanks were evaluated as to whether there were contaminants detected above the detection limit (DL). An equipment blank was not submitted with the samples in this data set.

Data validation qualifications for individual samples are based on the maximum contaminant concentration detected in all associated blanks.

Method blank results were reviewed for conformance with the QC acceptance criteria. Detected results in blanks are not discussed in this data validation report if the associated results were nondetect or if qualification of sample results was not required.

The QC acceptance criteria were met and/or qualification of the sample results was not required.

MS/MD and/or MSD Results

MS/MD/MSD analyses were not performed on samples reported in this SDG. There were no validation actions taken on this basis.

LCS/LCSD Results

The LCS/LCSD %Rs and/or relative percent recoveries (RPDs) were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

Field Duplicate Results

There were no field duplicate samples submitted with this data set. No validation actions were taken on this basis.

Sample Results/Reporting Issues

Compounds that were not detected in the sample are reported as not detected (U) at the Limit of Detection (LOD).

Compounds detected at concentrations less than the LOQ but greater than the detection limit (DL) were qualified by the laboratory as estimated (J). This "J" qualifier was retained during data validation.

Any sample that was analyzed at a dilution due to high concentrations of target or non-target compounds or matrix interferences was checked to ensure that the results and/or sample specific LODs and LOQs were adjusted accordingly by the laboratory.

QUALIFICATION ACTIONS

No sample results were qualified as a result of this data review.

ATTACHMENTS

Attachment A: Qualifier Codes and Explanations

Attachment A

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 limit of quantitation 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.								
R	The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified.								



600 Technology Way Scarborough, ME 04074 Tel: (207) 874-2400 Fax: (207) 775-4029

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Report of Analytical Results

Client: Rick Purdy

AECOM

701 Edgewater Drive Wakefield,MA 01880

Lab Sample ID: SH7523-1

Report Date: 25-SEP-14 Client PO: 16518

Project: Navy Clean WE15-03-0 SDG: SH7523

Matrix 7 SOIL-090514-573-575 Sample Description

Date Received 05-SEP-14 12:30:00 Date Sampled

SOIL-090514-573-57;	3-575					ST	05-SEP-14 12:30:00 11-SEP-14	:00 11-SEP	2.14	
Parameter	Result	Adj LOQ	Adj MDL	Adj LOD	Adj LOQ Adj MDL Adj LOD Angl. Method QC.Batch	QC.Batch	Anal. Date	Prep. Method Prep. Date Footnotes	Prep. Date	Footnotes
TOC In Soil	2200 ug/gdrywt	470	100	N/A	SW846 9060A	WG150664	N/A SW846 9060A WG150664 22-SEP-14 13:08:25 Mod	N/A	N/A	
Total Solids	84.%	-		N/A	SM2540G	WG150535	WG150535 22-SEP-14 09:16:37	SM2540G 19-SEP-14	19-SEP-14	



DATA VALIDATION REPORT

Project:	Regional Groundwater Inves	tigation — NWIRP Bethpage
Laboratory:	Katahdin Analytical	
Sample Delivery Groups:	SI0596	
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: 02/27/2015 Revised on: 10/27/2015
Reviewed by:	Tina Cantwell/Resolution Consultants	File Name: SI0596_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 22 January 2015 in accordance with the following Sampling and Analysis Plans:

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

Sample ID	Lab ID	Matrix/Sample Type	Analysis
BPOW6-5-SOIL-012215-778-780		Soil	9060A
BPOW6-5-SOIL-D-012215		Field Duplicate	9060A
BPOW6-5-Equipment Blank-012215		Equipment Blank	5310B

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



(DoD) Quality Systems Manual (QSM) for Environmental Laboratories, Version 4.2 (October 2010). In the absence of method-specific information, laboratory quality control (QC) limits, project-specific requirements and/or professional judgment were used as appropriate.

REVIEW ELEMENTS

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

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

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

Qualifications Actions

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

No results were qualified during this review. Analytical completeness was calculated to be 100% and the data are usable for their intended purpose, according to U.S. Environmental Protection Agency and Department of Defense guidelines. Katahdin Analytical discovered a QC error in Total Organic



Carbon calculations, informed us, and has implemented corrective action provided in Attachment A. Attachment B provides final results after data review.

ATTACHMENTS

Attachment A: Katahdin Analytical Corrective Action Report

Attachment B: Final Results after Data Review



Attachment A Katahdin Analytical Corrective Action Report

KATAHDIN ANALYTICAL SERVICES, INC. - CORRECTIVE ACTION REPORT

	,			
Problem Identification	on (Person initiating CAR) N	Name:	Leslie Dimond	Date: 08/28/2015
Discovered by Lab	oratory X	Discov	ered by Client (Compl	aint) Other
Details of Problem:				
by a factor of ten. Ger Calculations are set us method name, LOQ, process chains). This sample results, adjus was not adjusting LOT The lab had felt that the the lab decided to cha LOQ/LOD/MDL adjust	orge Brewer looked into the up to be performed automati LOD, MDL, limits and a spe is process chain contains all ted LOQs, LODs and MDLs Qs, LODs and/or MDLs for here wasn't really a standariange this and use 500 mg a	e matter ically in cific pro of the control of the control of the same of the street the street weight is the street weight evise the street weight is the street weight in the street weight is the street weight in the street weight weight in the street weight weight in the street weight weight in the street weight w	and found that their sa KIMS. Each product of cess chain for reporting calculations associated esults for QC samples. The amount used, but of the amount, so this was a candard and to incorportis TOC calculation in o	nt the results for their samples seemed to be off mple results were calculated incorrectly. ode for a test contains information such as the g to the either the LOQ, LOD or MDL (i.e. three with a test including separate calculations for In October of 2014, a client asked why the laborally for an instrument dilution and total solids. not incorporated. After numerous discussions, ate any different amount into the cur Laboratory Information System (KIMS). The tion was wrong.
The formula was: AD)J. LOQ/LOD/MDL = LOQ/L	.OD/MD	L * DF * (Sample amou	unt / 1000) * (100/TS).
calculations. Sample 2014, a data reviewe	results are checked in this	manner d limits v	, but adjusted limits are were not correct. This	are checked at a frequency of 10% of all e not necessarily checked. In December of issue was reviewed with MIS, and it was changed to:
ADJ. LOQ/LOD/MDL	= LOQ/LOD/MDL * DF * (50	00 / Sar	mple Amount) * (100/T	6) (where 500 is the standard sample amount).
At the time this chang needed to be corrected		er the in	npression that the sam	ple calculation was also incorrect and also
The formula for the sa	ample result was: Results (u		= <u>Total carbon</u> * sample amount / 1000)	(100/TS) (where 1000 is a conversion factor)
This was incorrectly of	changed to: Results (ug/g)		<u>l carbon</u> * TS Sample Amount)	
problem was not add during data review ar	ressed with the MIS departr	ment. H	lowever, there were se	ing data review and manually corrected. The veral cases where this error was not caught s issue was brought to the attention of MIS and
	otal carbon * (100/Ts ble amount / 1000)	S) (whe	ere 1000 is a conversio	n factor)
solids is not used in t		and MS	S/MSD samples are als	ated using different calculations since total o not affected because they use different
Associated Non-Co	nformances: List logbool	k and pa	age numbers	
There are no non-cor	nformances associated with	this cor	rective action.	
Root Cause Investig	gation & Determination (To	o be cor	npleted by Department	Manager, Operations Manager and/or QA
	elow and investigate to dete	rmine w	hether one of them, or	more than one, could be the cause of the
Possible Causes		· · · · · · · · · · · · · · · · · · ·	Details	

KATAHDIN ANALYTICAL SERVICES, INC. - CORRECTIVE ACTION REPORT

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

Corrective Action Plan: Name: Leslie Dimond

Date: 08/28/2015

Details of Corrective Action Plan -

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

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

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

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

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

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

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

KATAHDIN ANALYTICAL SERVICES, INC. – CORRECTIVE ACTION REPORT

senior management or the Quality Assurance Officer must also check the new calculations.							
Review & Approval of Corrective Action	n Plan						
Supervisor Approval:	SIGN,		Date: 99.11.15				
Operations Manager Approval:	total J. had	eau	Date: 9:1/15				
Quality Assurance Officer: Luseic T	_U		Date: 09-11-15				
Monitoring of Corrective Action (To be	completed by QA Officer	and/or Operations	Manager): List details of follow-up				
Corrective Action Effective	Return to Control –	Yes No	Further Monitoring Needed/Additional Corrective Action				
QA Approval:			Date:				
Additional Information:							
Additional information.							



Attachment B Final Results after Data Review

	Sample [eliver	y Group	SIOS	596	SIO	596	S105	596
			Lab I D	S105 ^o	96-1	S105	96-2	S1059	96-3
		Sa	mple I D	BP0W6-5-EQUIPMENT R	RINSATE BLANK-012215	BPOW6-5-SOIL-	012215-778-780	BPOW6-5-SO	IL-D-012215
		Samp	ole Date	1/22/	2015	1/22	/2015	1/22/2	2015
Sample Type		Equipme	nt Blank	S	oil	Field Du	plicate		
CAS									
Method Analyte No Units		Result	Qual	Result	Qual	Result	Qual		
5310B	TOTAL ORGANIC CARBON	-28	MG_L	0.36	J	NA		NA	
9060A	TOTAL ORGANIC CARBON	-28	UG_G	NA		390	J	290	J

Notes:

Milligrams per liter Micrograms per gram Not analyzed Final qualifier MG_L = UG_G =

NA Qual =

Final Qualifier: J

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



DATA VALIDATION REPORT

Project:	Regional Groundwater Investigation —	NWIRP Bethpage
Laboratory:	Katahdin Analytical	
Sample Delivery Group:	SI1908	
Analyses/Method:	Volatile Organic Compounds by U.S. EF 1,4-Dioxane by U.S. EPA SW-846 Meth (SIM)	
Validation Level:	3	
Project Number:	0888812477.SA.DV	
Prepared by:	Dana Miller/Resolution Consultants	Completed on: 06/08/2015
Reviewed by:	Tina Cantwell/Resolution Consultants	File Name: SI1908_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 26 March 2015 in accordance with the following Sampling and Analysis Plans:

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

Sample ID	Matrix/Sample Type	Analysis
BPOW6-1-GW-032615	Groundwater	8260C / 8270D_SIM
BPOW6-2-GW-032615	Groundwater	8260C / 8270D_SIM
BPOW6-3-GW-032615	Groundwater	8260C / 8270D_SIM
BPOW6-4-GW-032615	Field Duplicate	8260C / 8270D_SIM
TRIPBLANK032615	Groundwater	8260C

Data validation activities were conducted using the following guidance documents: *Test Methods for Evaluating Solid Waste, Physical/Chemical Methods SW-846, specifically Method 8260C, Volatile Organic Compounds by Gas Chromatography/Mass Spectrometry* (U.S. EPA, 2006), *SW-846 Method 8270D, Semivolatile Organic Compounds by Gas Chromatograph/Mass Spectrometry*



(U.S. EPA, 2007), *U.S. Environmental Protection Agency* (U.S. EPA) *Contract Laboratory Program National Functional Guidelines for Superfund Organic Methods Data Review* (June 2008), and Department of Defense Quality Systems Manual for Environmental Laboratories, Version 4.2 (October 2010). In the absence of method-specific information, laboratory quality control (QC) limits, project-specific requirements and/or professional judgment were used as appropriate.

REVIEW ELEMENTS

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

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

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

RESULTS

Initial Calibration/Continuing Calibration Verification
Calibration data were reviewed for conformance with the QC acceptance criteria to ensure that:

• the initial calibration percent relative standard deviation, correlation coefficient/coefficient of determination, and/or response factor method acceptance criteria were met;



- the initial calibration verification (ICV) standard percent recovery acceptance criteria were met;
- the continuing calibration verification standard (CCV) method percent difference or percent drift (%Ds) and response factor acceptance criteria were met; and
- the retention time method acceptance criteria were met.

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

ICV Recovery Non-conformance:

Cuitouio	Ac	tions
Criteria	Detected Results	Non-detected Results
Recovery >120%	J	UJ
Recovery < 80%	J	UJ

Notes:

J = Estimated

UJ = Undetected and estimated

CCV Linearity Non-conformance:

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

Notes:

J = Estimated

UJ = Undetected and estimated

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

Qualification 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 rejected; therefore, analytical completeness was calculated to be 100 percent. Data not qualified during data review are considered usable by the project. The remaining results qualified as estimated may be high or low, but the data are usable for their intended purpose, according to U.S. Environmental Protection Agency and Department of Defense guidelines. Final



data review qualifiers used to describe results and how they should be interpreted by the end data user are provided in Attachment B and Attachment C. Attachment D provides final results after data review.

ATTACHMENTS

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

Attachment A Non-Conformance Summary Table

		Initial Calibratio	Table A		onformance	
Method	Analyte	ICVID	%R	Limit	Associated Samples	Qualifier
8260C	Dichlorodifluoromethane	WG160458-7	132.53	80-120	BPOW6-1-GW-032615	UJ
8260C	Dichlorodifluoromethane	WG160458-7	132.53	80-120	BPOW6-2-GW-032615	UJ
8260C	Dichlorodifluoromethane	WG160458-7	132.53	80-120	BPOW6-3-GW-032615	UJ
8260C	Dichlorodifluoromethane	WG160458-7	132.53	80-120	BPOW6-4-GW-032615	UJ
8260C	Dichlorodifluoromethane	WG160458-7	132.53	80-120	TRIPBLANK032615	UJ
8260C	Chloromethane	WG160458-7	121.77	80-120	BPOW6-1-GW-032615	UJ
8260C	Chloromethane	WG160458-7	121.77	80-120	BPOW6-2-GW-032615	UJ
8260C	Chloromethane	WG160458-7	121.77	80-120	BPOW6-3-GW-032615	UJ
8260C	Chloromethane	WG160458-7	121.77	80-120	BPOW6-4-GW-032615	UJ
8260C	Chloromethane	WG160458-7	121.77	80-120	TRIPBLANK032615	UJ
8260C	Bromomethane	WG160458-7	130.01	80-120	BPOW6-1-GW-032615	UJ
8260C	Bromomethane	WG160458-7	130.01	80-120	BPOW6-2-GW-032615	UJ
8260C	Bromomethane	WG160458-7	130.01	80-120	BPOW6-3-GW-032615	UJ
8260C	Bromomethane	WG160458-7	130.01	80-120	BPOW6-4-GW-032615	UJ
8260C	Bromomethane	WG160458-7	130.01	80-120	TRIPBLANK032615	UJ
8260C	Carbon Disulfide	WG160458-7	544.89	80-120	BPOW6-1-GW-032615	UJ
8260C	Carbon Disulfide	WG160458-7	544.89	80-120	BPOW6-2-GW-032615	UJ
8260C	Carbon Disulfide	WG160458-7	544.89	80-120	BPOW6-3-GW-032615	UJ
8260C	Carbon Disulfide	WG160458-7	544.89	80-120	BPOW6-4-GW-032615	UJ
8260C	Carbon Disulfide	WG160458-7	544.89	80-120	TRIPBLANK032615	UJ
8260C	Acetone	WG160458-7	137.34	80-120	BPOW6-1-GW-032615	UJ
8260C	Acetone	WG160458-7	137.34	80-120	BPOW6-2-GW-032615	UJ
8260C	Acetone	WG160458-7	137.34	80-120	BPOW6-3-GW-032615	UJ
8260C	Acetone	WG160458-7	137.34	80-120	BPOW6-4-GW-032615	UJ
8260C	Acetone	WG160458-7	137.34	80-120	TRIPBLANK032615	UJ
8260C	2-Butanone	WG160458-7	134.65	80-120	BPOW6-1-GW-032615	UJ
8260C	2-Butanone	WG160458-7	134.65	80-120	BPOW6-2-GW-032615	UJ
8260C	2-Butanone	WG160458-7	134.65	80-120	BPOW6-3-GW-032615	UJ
8260C	2-Butanone	WG160458-7	134.65	80-120	BPOW6-4-GW-032615	UJ
8260C	2-Butanone	WG160458-7	134.65	80-120	TRIPBLANK032615	UJ
8260C	Cyclohexane	WG160458-7	193.37	80-120	BPOW6-1-GW-032615	UJ
8260C	Cyclohexane	WG160458-7	193.37	80-120	BPOW6-2-GW-032615	UJ
8260C	Cyclohexane	WG160458-7	193.37	80-120	BPOW6-3-GW-032615	UJ
8260C	Cyclohexane	WG160458-7	193.37	80-120	BPOW6-4-GW-032615	UJ
8260C	Cyclohexane	WG160458-7	193.37	80-120	TRIPBLANK032615	Ŋ
8260C	Tetrachloroethene	WG160458-7	124.99	80-120	BPOW6-1-GW-032615	UJ
8260C	Tetrachloroethene	WG160458-7	124.99	80-120	BPOW6-2-GW-032615	UJ
8260C	Tetrachloroethene	WG160458-7	124.99	80-120	BPOW6-3-GW-032615	UJ
8260C	Tetrachloroethene	WG160458-7	124.99	80-120	BPOW6-4-GW-032615	UJ
8260C	Tetrachloroethene	WG160458-7	124.99	80-120	TRIPBLANK032615	UJ
8260C	2-Hexanone	WG160458-7	130.94	80-120	BPOW6-1-GW-032615	UJ
8260C	2-Hexanone	WG160458-7	130.94	80-120	BPOW6-2-GW-032615	UJ
8260C	2-Hexanone	WG160458-7	130.94	80-120	BPOW6-3-GW-032615	UJ
8260C	2-Hexanone	WG160458-7	130.94	80-120	BPOW6-4-GW-032615	UJ
8260C	2-Hexanone	WG160458-7	130.94	80-120	TRIPBLANK032615	UJ 03

Notes: ICV Initial calibration verification

%R

Percent recovery Non-detected analyte in associated sample qualified estimated "UJ" due to potential bias. UJ

	Table A-2 Continuing Calibration Verification Non-Conformance					
Method	Analyte	CCVID	%D	Limit	Associated Samples	Qualifier
8260C	Chloroethane	C2041.D	21.59	20	BPOW6-1-GW-032615	UJ
8260C	Chloroethane	C2041.D	21.59	20	BPOW6-2-GW-032615	ΟΊ
8260C	Chloroethane	C2041.D	21.59	20	BPOW6-3-GW-032615	UJ
8260C	Chloroethane	C0241.D	21.59	20	BPOW6-4-GW-032615	UJ
8260C	Chloroethane	C0241.D	21.59	20	TRIPBLANK032615	UJ

Notes: CCV %D UJ =

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

Attachment B Qualifier Codes and Explanations

Qualifier	Explanation
UJ	The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual quantitation limit necessary to accurately and precisely measure the analyte in the sample.
U	The analyte was analyzed for, but was not detected above the reported sample quantitation limit.

Attachment C Reason Codes and Explanations

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

Attachment D Final Results after Data Review

		Sa	Lab ID Sample ID mple Date	SI1908 SI1908-1 BPOW6-1-GW-032615 3/26/2015		
	1		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	0.5	U	
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.5	U	
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	0.5	U	
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	0.5	U	
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	U	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	U	
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	U	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	U	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	1	U	
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	U	
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	U	
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	U	
8260C	2-BUTANONE	78-93-3	UG_L	2.5	UJ	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	Ü	
8260C	BROMOFORM	75-25-2	UG L	0.5	U	
8260C	BROMOMETHANE	74-83-9	UG L	1	UJ	С
8260C	CARBON DISULFIDE	75-15-0	UG L	0.5	UJ	С
8260C	CARBON TETRACHLORIDE	56-23-5	UG L	0.5	U	
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	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	UJ	С
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG L	0.5	U	
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG L	1	UJ	С
8260C	ETHYLBENZENE ETHYLBENZENE	100-41-4	UG L	0.5	U	
8260C	ISOPROPYLBENZENE	98-82-8	UG L	0.5	Ü	
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	1
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG L	0.5	U	1
8260C 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		_		UJ	-
8260C 8260C	TOLUENE	127-18-4 108-88-3	UG_L UG L	0.5 0.5	U	С
	TRANS-1,2-DICHLOROETHENE		_			1
8260C		156-60-5	UG_L	0.5	U	-
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	U	1
8260C	TRICHLOROETHENE	79-01-6	UG_L	0.5	U	
8260C	TRICHLOROFLUOROMETHANE	75-69-4	UG_L	1	U	
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	U	
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	U	-
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	0.17	U	

			very Group Lab ID Sample ID ample Date ample Type	SI BPOW6-3/2	611908 1908-2 2-GW-03261 26/2015 undwater	15
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	U	
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	U	
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	0.5	U	
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.5	U	
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	0.5	U	
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	0.5	U	
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	U	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	U	
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	U	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	U	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	1	U	
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	U	
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	U	
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	U	
8260C	2-BUTANONE	78-93-3	UG_L	2.5	UJ	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	
8260C	BROMOFORM	75-25-2	UG L	0.5	U	
8260C	BROMOMETHANE	74-83-9	UG L	1	UJ	С
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	UJ	С
8260C	CARBON TETRACHLORIDE	56-23-5	UG L	0.5	U	
8260C	CHLOROBENZENE	108-90-7	UG L	0.5	U	
8260C	CHLOROETHANE	75-00-3	UG L	1	UJ	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	0.5	U	
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG L	0.5	U	
8260C	CYCLOHEXANE	110-82-7	UG L	0.5	UJ	С
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG L	0.5	U	
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	1	UJ	С
8260C	ETHYLBENZENE	100-41-4	UG_L	0.5	U	
8260C	ISOPROPYLBENZENE	98-82-8	UG_L	0.5	U	
8260C	M- AND P-XYLENE	108-38-3/106-42	UG_L	1	U	
8260C	METHYL ACETATE	79-20-9	UG_L	0.75	U	
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	U	
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	U	
8260C	METHYLENE CHLORIDE	75-09-2	UG_L	2.5	U	
8260C	O-XYLENE	95-47-6	UG_L	0.5	U	
8260C	STYRENE	100-42-5	UG_L	0.5	U	
8260C	TETRACHLOROETHENE	127-18-4	UG_L	0.5	UJ	С
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	Ü	
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	0.19	Ü	

		Sa	Lab ID Sample ID mple Date mple Type	SI1908 SI1908-3 BPOW6-3-GW-032615 3/26/2015 Groundwater		
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	U	
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	U	
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	0.5	U	
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.5	U	
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	0.5	U	
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	0.5	U	
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	U	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	U	
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	U	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	U	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	1	U	
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	U	
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	U	
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	U	
8260C	2-BUTANONE	78-93-3	UG_L	2.5	UJ	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	
8260C	BROMOFORM	75-25-2	UG_L	0.5	U	
8260C	BROMOMETHANE	74-83-9	UG_L	1	UJ	С
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	UJ	С
8260C	CARBON TETRACHLORIDE	56-23-5	UG_L	0.5	U	
8260C	CHLOROBENZENE	108-90-7	UG_L	0.5	U	
8260C	CHLOROETHANE	75-00-3	UG_L	1	UJ	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	0.5	U	
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG_L	0.5	U	
8260C	CYCLOHEXANE	110-82-7	UG_L	0.5	UJ	С
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG_L	0.5	U	
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	1	UJ	С
8260C	ETHYLBENZENE	100-41-4	UG_L	0.5	U	
8260C	ISOPROPYLBENZENE	98-82-8	UG_L	0.5	U	
8260C	M- AND P-XYLENE	108-38-3/106-42	UG_L	1	U	
8260C	METHYL ACETATE	79-20-9	UG_L	0.75	U	
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	U	
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	U	
8260C	METHYLENE CHLORIDE	75-09-2	UG_L	2.5	U	
8260C	O-XYLENE	95-47-6	UG_L	0.5	U	
8260C	STYRENE	100-42-5	UG_L	0.5	U	
8260C	TETRACHLOROETHENE	127-18-4	UG_L	0.5	UJ	С
8260C	TOLUENE	108-88-3	UG_L	0.5	U	
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	U	
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	U	
8260C	TRICHLOROETHENE	79-01-6	UG_L	0.5	U	
8260C	TRICHLOROFLUOROMETHANE	75-69-4	UG_L	1	U	
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	U	
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	U	
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	0.17	U	

			Lab ID Sample ID ample Date ample Type	SI* BPOW6-4 3/2	11908 1908-4 I-GW-03261 6/2015 Indwater	5
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	U	
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	U	
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	0.5	U	
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.5	U	
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	0.5	U	
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	0.5	U	
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	U	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	U	
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	U	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	U	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	1	U	
8260C	1,2-DICHLOROPROPANE	78-87-5	UG L	0.5	U	
8260C	1,3-DICHLOROBENZENE	541-73-1	UG L	0.5	U	
8260C	1,4-DICHLOROBENZENE	106-46-7	UG L	0.5	U	
8260C	2-BUTANONE	78-93-3	UG L	2.5	UJ	С
8260C	2-HEXANONE	591-78-6	UG L	2.5	U.J	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	Ů
8260C	BROMODICHLOROMETHANE	75-27-4	UG L	0.5	U	
8260C	BROMOFORM	75-25-2	UG L	0.5	U	
8260C	BROMOMETHANE	74-83-9	UG L	1	UJ	С
8260C	CARBON DISULFIDE	75-15-0	UG L	0.5	UJ	С
8260C	CARBON TETRACHLORIDE	56-23-5	UG L	0.5	U	
8260C	CHLOROBENZENE	108-90-7	UG L	0.5	U	1
8260C	CHLOROETHANE	75-00-3	UG L	1	UJ UJ	С
8260C	CHLOROFORM	67-66-3	UG L	0.5	U	-
8260C	CHLOROMETHANE	74-87-3	UG L	1	UJ UJ	С
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG L	0.5	U	C
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG_L	0.5	U	
8260C	CYCLOHEXANE				UJ	-
8260C 8260C	DIBROMOCHLOROMETHANE	110-82-7 124-48-1	UG_L UG L	0.5 0.5	U	С
8260C 8260C		75-71-8	UG_L	1	UJ	-
	DICHLORODIFLUOROMETHANE			·	UJ UJ	С
8260C 8260C	ETHYLBENZENE ISOPROPYLBENZENE	100-41-4 98-82-8	UG_L	0.5 0.5	U	
8260C 8260C		108-38-3/106-42	UG_L UG L	1	U	1
	M- AND P-XYLENE					
8260C	METHYL ACETATE	79-20-9	UG_L	0.75	U	
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	U	-
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	U	-
8260C	METHYLENE CHLORIDE	75-09-2	UG_L	2.5	U	
8260C	O-XYLENE	95-47-6	UG_L	0.5	U	1
8260C	STYRENE	100-42-5	UG_L	0.5	U	1
8260C	TETRACHLOROETHENE	127-18-4	UG_L	0.5	UJ	С
8260C	TOLUENE	108-88-3	UG_L	0.5	U	1
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	U	<u> </u>
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	U	ļ
8260C	TRICHLOROETHENE	79-01-6	UG_L	0.5	U	
8260C	TRICHLOROFLUOROMETHANE	75-69-4	UG_L	1	U	<u> </u>
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	U	<u> </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	

Samp Samp			Livery Group Lab ID Sample ID Sample Date Sample Type	Lab ID SI1908-5 Sample ID TRIPBLANK03 ample Date 3/26/201 ample Type Trip Blant		
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	U	
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	U	
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	0.5	U	
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.5	U	
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	0.5	U	
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	0.5	U	
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	U	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	U	
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	U	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	U	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	1	U	
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	U	
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	U	
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	U	
8260C	2-BUTANONE	78-93-3	UG_L	2.5	UJ	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	
8260C	BROMOFORM	75-25-2	UG_L	0.5	U	
8260C	BROMOMETHANE	74-83-9	UG_L	1	UJ	С
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	UJ	С
8260C	CARBON TETRACHLORIDE	56-23-5	UG_L	0.5	U	
8260C	CHLOROBENZENE	108-90-7	UG_L	0.5	U	
8260C	CHLOROETHANE	75-00-3	UG_L	1	UJ	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	0.5	U	
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG_L	0.5	U	
8260C	CYCLOHEXANE	110-82-7	UG_L	0.5	UJ	С
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG_L	0.5	U	
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	1	UJ	С
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	1
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	С
8260C	TOLUENE	108-88-3	UG_L	0.5	U	
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	U	
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	U	
8260C	TRICHLOROETHENE	79-01-6	UG_L	0.5	U	
8260C	TRICHLOROFLUOROMETHANE	75-69-4	UG_L	1	U	
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	U	
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	U	
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	NA		

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



DATA VALIDATION REPORT

Project:	Regional Groundwater Investigation — NWIRP Bethpage				
Laboratory:	Katahdin Analytical				
Sample Delivery Group:	SI2993				
Analyses/Method:	Total Organic Carbon by U.S. EPA SW-846 Method 9060A				
Validation Level:	3				
Project Number:	0888812477.SA.DV				
Prepared by:	Dana Miller/Resolution Consultants Completed on: 05/9/2015				
Reviewed by:	Tina Cantwell/Resolution File Name: S12993_9060A Consultants				

SUMMARY

This report summarizes data review findings for samples listed below, collected by Resolution Consultants from the Regional Groundwater Investigation — NWIRP Bethpage Site on 5 to 6 May 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
BPOW6-5-SOIL-050515-528-530	Soil	9060A
BPOW6-5-EB-050615	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), *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 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
- NA Gas chromatography/Mass spectrometer performance checks
- ✓ Initial calibration verification /continuing calibration verification
- ✓ Laboratory blanks/trip blanks
- NA Surrogate spike recoveries
- ✓ Matrix spike and/or matrix spike duplicate results
- ✓ Laboratory control sample/laboratory control sample duplicate results
- NA Field duplicates
- NA Internal standards
- ✓ Sample results/reporting issues

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

Qualifications Actions

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

No results were qualified during this data review. Attachment A provides final results after data review.

ATTACHMENTS

Attachment A: Final Results after Data Review

Attachment A Final Results after Data Review

Sample Delivery Group			SI299	93	SI299	93	
Lab ID			SI2993	3-1	SI299	3-2	
Sample ID B			BP0W6-5-SOIL-050515-528-530		30 BP0W6-5-EB-050615		
			Sample Date	5/5/20	15	5/6/20)15
			Sample Type	Soil		Equipmen	t Blank
Method	Analyte	CAS No.	Units	Result	Qual	Result	Qual
5310B	TOTAL ORGANIC CARBON	-28	MG_L	NA		0.2	J
9060A	TOTAL ORGANIC CARBON	-28	UG_G	590		NA	_

Notes:

ID = Identification Milligrams per liter
Micrograms per gram
Not analyzed
Final qualifier
Estimated value below the quantitation limit MG_L = UG_G =

NA = Qual =

J



DATA VALIDATION REPORT

Project:	Regional Groundwater Investigation — NWIRP Bethpage				
Laboratory:	Katahdin Analytical				
Sample Delivery Group:	SI3513				
Analyses/Method:	Total Organic Carbon by U.S. EPA SW-846 Method 9060A				
Validation Level:	3				
Project Number:	0888812477.SA.DV				
Prepared by:	Dana Miller/Resolution Consultants Completed on: 05/30/2015				
Reviewed by:	Tina Cantwell/Resolution File Name: SI3513_9060A Consultants				

SUMMARY

This report summarizes data review findings for samples listed below, collected by Resolution Consultants from the Regional Groundwater Investigation — NWIRP Bethpage Site on 21 May 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
BPOW6-6-052115-773-775	Soil	9060A
BPOW6-6-FD-052115	Field Duplicate	9060A
BPOW6-6-052115-EB	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), *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 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
- NA Gas chromatography/Mass spectrometer performance checks
- ✓ Initial calibration verification /continuing calibration verification
- ✓ Laboratory blanks/trip blanks
- NA Surrogate spike recoveries
- ✓ Matrix spike and/or matrix spike duplicate results
- ✓ Laboratory control sample/laboratory control sample duplicate results
- ✓ Field duplicates
- NA Internal standards
- ✓ Sample results/reporting issues

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

Qualifications Actions

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

No results were qualified during this data review. Attachment A provides final results after data review.

ATTACHMENTS

Attachment A: Final Results after Data Review

Attachment A Final Results after Data Review

Sample Delivery Group			SI3513		SI351	3	SI3513	3	
Lab ID			SI3513-1		SI3513	-2	SI3513-	-3	
Sample ID			BPOW6-6-052115	-773-775	BPOW6-6-FD	-052115	BPOW6-6-052	2115-EB	
	Sample Date		5/21/201	5	5/21/20	15	5/21/20	15	
		Sar	mple Type	Soil		Field Dupl	icate	Equipment	Blank
Method	Analyte	CAS No	Units	Result	Qual	Result	Qual	Result	Qual
5310B	TOTAL ORGANIC CARBON	-28	MG_L	NA		NA		0.36	J
9060A	TOTAL ORGANIC CARBON	-28	UG_G	260	J	250	J	NA	

Notes:

ID = Identification Milligrams per liter
Micrograms per gram
Not analyzed
Final qualifier
Estimated value below the quantitation limit $MG_L =$ UG_G =

NA = Qual =



DATA VALIDATION REPORT

Project:	Regional Groundwater Investigation — NWIRP Bethpage		
Laboratory:	Katahdin Analytical		
Sample Delivery Group:	S14556		
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: 07/30/2015	
Reviewed by:	Tina Cantwell/Resolution Consultants	File Name: SI4556_8260C_8270D	

SUMMARY

This report summarizes data review findings for samples listed below, collected by Resolution Consultants from the Regional Groundwater Investigation — NWIRP Bethpage Site on 24 and 25 June 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
BPOW5-6-GW-062415	Groundwater	8260C/8270D_SIM
BPOW5-5-GW-062415	Groundwater	8260C/8270D_SIM
RE118D1-GW-062415	Groundwater	8260C/8270D_SIM
RE108D1-GW-062415	Groundwater	8260C/8270D_SIM
RE108D2-GW-062415	Groundwater	8260C/8270D_SIM
BPOW6-5-GW-062515	Groundwater	8260C/8270D_SIM
BPOW6-6-GW-062515	Field Duplicate	8260C/8270D_SIM
RE117D1-GW-062515	Groundwater	8260C/8270D_SIM
TRIPBLANK-062515	Trip Blank	8260C



Data validation activities were conducted using the following guidance documents: *Test Methods for Evaluating Solid Waste, Physical/Chemical Methods SW-846, specifically Method 8260C, Volatile Organic Compounds by Gas Chromatography/Mass Spectrometry* (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 FLEMENTS

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 verification (ICV)/continuing calibration verification (CCV)
- ✓ Laboratory blanks/trip blanks
- ✓ Surrogate spike recoveries
- NA Matrix spike and/or matrix spike duplicate results
- ✓ Laboratory control sample/laboratory control sample duplicate results
- NA Field duplicates
- ✓ Internal standards
- ✓ Sample results/reporting issues

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



RESULTS

Initial Calibration/Continuing Calibration Verification

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

- The initial calibration percent relative standard deviation, correlation coefficient/coefficient of determination, and/or response factor method acceptance criteria were met
- The 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 calibration verification was as follows:

ICV Recovery Non-conformance:

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

Notes:

J = Estimated

UJ = Undetected and estimated

ICV non-conformances are summarized in Attachment A in Table A-1.

Qualifications Actions

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



No results were rejected; therefore, analytical completeness was calculated to be 100 percent. Data not qualified during data review are considered usable by the project. The remaining results qualified as estimated may be high or low, but the data are usable for their intended purpose, according to U.S. 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 Verification Non-Conformance									
Method	Analyte	ICVID	%R	Limit	Associated Samples	Qualifier				
8260C	Chloroethane	P1539A	127.5	80-120	BPOW5-6-GW-062415	UJ				
8260C	Chloroethane	P1539A	127.5	80-120	BPOW5-5-GW-062415	UJ				
8260C	Chloroethane	P1539A	127.5	80-120	RE118D1-GW-062415	UJ				
8260C	Chloroethane	P1539A	127.5	80-120	RE108D1-GW-062415	UJ				
8260C	Chloroethane	P1539A	127.5	80-120	RE108D2-GW-062415	UJ				
8260C	Chloroethane	P1539A	127.5	80-120	BPOW6-5-GW-062515	UJ				
8260C	Chloroethane	P1539A	127.5	80-120	BPOW6-6-GW-062515	UJ				
8260C	Chloroethane	P1539A	127.5	80-120	RE117D1-GW-062515	UJ				
8260C	Chloroethane	P1539A	127.5	80-120	TRIPBLANK-062515	UJ				

Notes:

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

Attachment B

Qualifier Codes and Explanations

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

Attachment C Reason Codes and Explanations

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

Attachment D Final Results after Data Review

			very Group Lab ID Sample ID mple Date	SI BPOW5-6	I 4556 4556-1 5-GW-0624 24/2015	15
			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	0.5	U	
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.5	U	
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	0.5	U	
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	0.5	U	
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	U	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	U	
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	U	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	U	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	1	U	
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	U	
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	U	
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	U	
8260C	2-BUTANONE	78-93-3	UG_L	12		
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	U	
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	UJ	С
8260C	CHLOROFORM	67-66-3	UG_L	0.5	U	
8260C	CHLOROMETHANE	74-87-3	UG_L	1	U	
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 0.5	U	
8260C	ETHYLBENZENE	100-41-4	UG_L	0.5		
8260C 8260C	ISOPROPYLBENZENE M- AND P-XYLENE	98-82-8 108-38-3/106-42	UG_L UG L	0.5 1	U	
		79-20-9	UG_L UG L	0.75	U	
8260C 8260C	METHYL ACETATE METHYL CYCLOHEXANE	108-87-2	UG_L	0.75	U	
8260C		1634-04-4	UG_L		U	
8260C	METHYL TERT-BUTYL ETHER METHYLENE CHLORIDE	75-09-2	UG_L	0.5 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	U	
8260C	TOLUENE	108-88-3	UG_L	0.74	J	
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG L	0.74	U	
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG L	0.5	U	
8260C	TRICHLOROETHENE	79-01-6	UG_L	0.45	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	

		Sa	very Group Lab ID Sample ID ample Date ample Type	SI- BPOW5-5 6/2	14556 4556-2 5-GW-0624 14/2015 Indwater	115
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG L	0.5	U	
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	U	
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	0.5	U	
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.5	U	
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	0.5	U	
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	0.5	U	
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	U	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	U	
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	U	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	U	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	1	U	
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	U	
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	U	
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	U	
8260C	2-BUTANONE	78-93-3	UG_L	2.5	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	8		
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	UJ	С
8260C	CHLOROFORM	67-66-3	UG_L	0.5	U	-
8260C	CHLOROMETHANE	74-87-3	UG_L	1	U	-
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	0.5	U	
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5 110-82-7	UG_L	0.5	U	
8260C	CYCLOHEXANE		UG_L	0.5	_	
8260C	DIBROMOCHLOROMETHANE DICHLORODIFLUOROMETHANE	124-48-1	UG_L	0.5 1	U	
8260C 8260C	ETHYLBENZENE	75-71-8 100-41-4	UG_L UG L	0.5	U	
8260C 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 8260C	METHYL ACETATE	79-20-9	UG_L	0.75	U	1
8260C	METHYL CYCLOHEXANE	108-87-2	UG L	0.75	U	
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG L	0.5	U	
8260C	METHYLENE CHLORIDE	75-09-2	UG L	2.5	U	
8260C	O-XYLENE	95-47-6	UG L	0.5	U	1
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	
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	0.18	U	

		Sa	very Group Lab ID Sample ID mple Date mple Type	SI- RE118D1 6/2	14556 4556-3 I-GW-0624 24/2015 Indwater	15
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG L	0.5	U	
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	U	
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	0.5	U	
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.5	U	
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	0.5	U	
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	0.5	U	
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	U	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	U	
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	U	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	U	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	1	U	
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	U	
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	U	
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	U	
8260C	2-BUTANONE	78-93-3	UG_L	2.5	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	2.7	J	
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	11	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	UJ	С
8260C	CHLOROFORM	67-66-3	UG_L	0.5	U	
8260C	CHLOROMETHANE	74-87-3	UG_L	1 0.5	U	
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	0.5	U	
8260C 8260C	CIS-1,3-DICHLOROPROPENE CYCLOHEXANE	10061-01-5 110-82-7	UG_L UG_L	0.5 0.5	U	
8260C 8260C	DIBROMOCHLOROMETHANE	124-48-1	UG_L UG L	0.5	U	
8260C 8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L UG L	1	U	
8260C 8260C	ETHYLBENZENE	100-41-4	UG_L	0.5	U	
8260C 8260C	ISOPROPYLBENZENE	98-82-8	UG_L UG L	0.5	U	+
8260C 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	1
8260C	METHYL CYCLOHEXANE	108-87-2	UG L	0.5	U	
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG L	0.5	U	
8260C	METHYLENE CHLORIDE	75-09-2	UG L	2.5	U	
8260C	O-XYLENE	95-47-6	UG L	0.5	U	
8260C	STYRENE	100-42-5	UG L	0.5	U	
8260C	TETRACHLOROETHENE	127-18-4	UG_L	0.5	U	
8260C	TOLUENE	108-88-3	UG_L	0.38	J	
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	U	
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	U	
8260C	TRICHLOROETHENE	79-01-6	UG_L	0.5	U	
8260C	TRICHLOROFLUOROMETHANE	75-69-4	UG_L	1	U	
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	U	
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	U	
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	0.18	U	

		Sample Deli Sa Sa	very Group Lab ID Sample ID ample Date ample Type	SI RE108D 6/	514556 14556-4 1-GW-062 24/2015 undwater	<u>'</u> 415
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		
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	0.34	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	U	
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	Ü	
8260C	BROMOMETHANE	74-83-9	UG L	1	Ü	
8260C	CARBON DISULFIDE	75-15-0	UG L	0.5	U	
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	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	0.34	J	
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 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	
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG L	0.5	U	
8260C	METHYLENE CHLORIDE	75-09-2	UG L	2.5	U	
8260C	O-XYLENE	95-47-6	UG_L	0.5	U	
8260C	STYRENE	100-42-5	UG_L	0.5	U	
8260C 8260C	TETRACHLOROETHENE	127-18-4	UG_L	1.4	U	
8260C 8260C	TOLUENE	108-88-3	UG_L	0.5	U	<u> </u>
8260C 8260C	TRANS-1,2-DICHLOROETHENE	156-60-5		0.5		1
	·		UG_L		U	-
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	U	-
8260C	TRICHLOROETHENE	79-01-6	UG_L	110	11	-
8260C	TRICHLOROFLUOROMETHANE	75-69-4	UG_L	1	U	
8260C	VINYL CHLORIDE	75-01-4	UG_L	1 -	U	1
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	U	1
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	5.2	L	

		Sample Deliv Sa Sa Sa	SI 4556 SI 4556-5 RE108D2-GW-062415 6/24/2015 Groundwater			
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG L	0.98	J	
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG L	0.5	U	
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	6.8	- U	
8260C	1.1.2-TRICHLOROETHANE	79-00-5	UG L	1.8		
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	4.6		
8260C	1.1-DICHLOROETHENE	75-35-4	UG L	6.6		
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG L	0.5	U	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	U	
8260C	1,2-DIBROMOETHANE	106-93-4	UG L	0.5	U	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	U	
					U	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	8.1	11	1
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	U	
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	U	1
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	U	
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	1.5		
8260C	CHLOROBENZENE	108-90-7	UG_L	0.5	U	
8260C	CHLOROETHANE	75-00-3	UG_L	1	UJ	С
8260C	CHLOROFORM	67-66-3	UG L	3.5	0.5	
8260C	CHLOROMETHANE	74-87-3		1	U	
			UG_L		U	-
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	8.1		
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	U	
8260C	O-XYLENE	95-47-6	UG L	0.5	U	
8260C	STYRENE	100-42-5	UG L	0.5	U	
8260C	TETRACHLOROETHENE	127-18-4	UG_L	2.2		
8260C	TOLUENE	108-88-3	UG L	0.5	U	1
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG L	0.5	U	t
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG L	0.5	U	t
8260C	TRICHLOROETHENE	79-01-6	UG_L	3900	U	1
8260C	TRICHLOROFLUOROMETHANE	75-69-4	UG_L	1	U	+
					U	1
8260C	VINYL CHLORIDE	75-01-4	UG_L	11		1
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	U	

			Lab ID Lab ID Sample ID Sample Date Sample Type	SI 45 BPOW6-5 6/2	14556 556-6RA 5-GW-062 5/2015 Indwater	515
Method	Analyte	CAS No	Units	Result	Qual	RC
	-				1	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	U	+
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	U	
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	0.5	U	-
8260C	1,1,2-TRICHLOROETHANE	79-00-5 75-34-3	UG_L	0.5	U	+
8260C	1,1-DICHLOROETHANE		UG_L	0.5	U	+
8260C	1,1-DICHLOROETHENE 1,2,4-TRICHLOROBENZENE	75-35-4	UG_L	0.5	U	-
8260C		120-82-1	UG_L	0.5 0.75	U	+
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L			+
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	<u> </u>
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	U	
8260C	ACETONE	67-64-1	UG_L	2.5	U	-
8260C	BENZENE	71-43-2	UG_L	0.5	U	<u> </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	UJ	С
8260C	CHLOROFORM	67-66-3	UG_L	0.5	U	
8260C	CHLOROMETHANE	74-87-3	UG_L	1	U	
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	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	U	
8260C	TOLUENE	108-88-3	UG_L	0.76	J	
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	U	T
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	U	
8260C	TRICHLOROETHENE	79-01-6	UG_L	0.5	U	1
8260C	TRICHLOROFLUOROMETHANE	75-69-4	UG L	1	Ü	
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	Ü	1
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	Ü	1
3270D_SIM	1,4-DIOXANE	123-91-1	UG_L	0.17	Ü	1

		•	Lab ID Sample ID Sample Date	SI45 BPOW6-6	4556 556-7RA -GW-062 5/2015	515
			Sample Type	Grou	ndwater	
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	U	
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	U	
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	0.5	U	
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.5	U	
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	0.5	U	
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	0.5	U	1
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	I
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	L
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	U	
8260C	BENZENE	71-43-2	UG_L	0.5	U	1
8260C	BROMODICHLOROMETHANE	75-27-4	UG_L	0.5	U	Ī
8260C	BROMOFORM	75-25-2	UG L	0.5	U	
8260C	BROMOMETHANE	74-83-9	UG_L	1	U	Ì
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	U	Ì
8260C	CARBON TETRACHLORIDE	56-23-5	UG_L	0.5	U	Ì
8260C	CHLOROBENZENE	108-90-7	UG_L	0.5	U	
8260C	CHLOROETHANE	75-00-3	UG_L	1	UJ	С
8260C	CHLOROFORM	67-66-3	UG_L	0.5	U	1
8260C	CHLOROMETHANE	74-87-3	UG_L	1	U	Ī
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	0.5	U	1
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG L	0.5	U	Ī
8260C	CYCLOHEXANE	110-82-7	UG_L	0.5	U	1
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG_L	0.5	U	1
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	1	U	1
8260C	ETHYLBENZENE	100-41-4	UG_L	0.5	U	
8260C	ISOPROPYLBENZENE	98-82-8	UG_L	0.5	U	1
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	Ī
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	U	T
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	U	T
8260C	METHYLENE CHLORIDE	75-09-2	UG_L	2.5	U	1
8260C	O-XYLENE	95-47-6	UG_L	0.5	U	
8260C	STYRENE	100-42-5	UG_L	0.5	U	T
8260C	TETRACHLOROETHENE	127-18-4	UG_L	0.5	U	1
8260C	TOLUENE	108-88-3	UG_L	1		
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	U	T
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	1
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	U	Ī
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	Ü	
3270D_SIM	1,4-DIOXANE	123-91-1	UG_L	0.17	U	

		·	Lab ID Sample ID Sample Date Sample Type	SI4 RE117D1 6/2	4556 4556-8 -GW-062! 5/2015 ndwater	515
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	U	INC
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	U	+
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	0.5	U	+
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.5	U	+
8260C	1,1-DICHLOROETHANE	75-34-3	UG L	0.5	U	+
8260C	1,1-DICHLOROETHENE	75-35-4	UG L	0.5	U	+
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG L	0.5	U	+
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG L	0.75	U	1
8260C	1,2-DIBROMOETHANE	106-93-4	UG L	0.5	U	1
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	1
8260C	1,2-DICHLOROETHANE	107-06-2	UG L	0.5	U	1
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	\vdash
8260C	1,4-DICHLOROBENZENE	106-46-7	UG L	0.5	U	
8260C	2-BUTANONE	78-93-3	UG L	2.1	J	
8260C	2-HEXANONE	591-78-6	UG L	2.5	Ü	
8260C	4-METHYL-2-PENTANONE	108-10-1	UG L	2.5	U	1
8260C	ACETONE	67-64-1	UG L	2.5	Ü	1
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	U	1
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	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	U	
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	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	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	7.8		
8260C	TRICHLOROFLUOROMETHANE	75-69-4	UG_L	1	U	
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	U	
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	U	
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	0.17	U	

			Lab ID Sample ID Sample Date	SI4 TRIPBLA 6/2	14556 1556-9 ANK-0625 4/2015	15
			Sample Type		o Blank	1
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	U	
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	U	
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	0.5	U	
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.5	U	
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	0.5	U	
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	0.5	U	
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	U	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	U	
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	U	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	U	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	1	U	
8260C	1,2-DICHLOROPROPANE	78-87-5	UG L	0.5	U	
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	U	
8260C	1,4-DICHLOROBENZENE	106-46-7	UG L	0.5	Ü	
8260C	2-BUTANONE	78-93-3	UG L	2.5	Ü	
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	U	
8260C	BENZENE	71-43-2	UG L	0.5	U	
8260C	BROMODICHLOROMETHANE	75-27-4	UG L	0.5	U	
8260C	BROMOFORM	75-27-4	UG_L	0.5	U	
8260C 8260C	BROMOMETHANE	75-25-2	UG_L	1	U	
8260C 8260C	CARBON DISULFIDE	75-15-0	UG_L UG L	0.5		
					U	
8260C	CARBON TETRACHLORIDE	56-23-5	UG_L	0.5		
8260C	CHLOROBENZENE	108-90-7	UG_L	0.5	U	
8260C	CHLOROETHANE	75-00-3	UG_L	1	UJ	С
8260C	CHLOROFORM	67-66-3	UG_L	0.5	U	ļ
8260C	CHLOROMETHANE	74-87-3	UG_L	1	U	
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	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	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	
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	NA	U	
02/UU_3IIVI	I 1,4-DIOAANL	123-71-1	UU_L	INA	1	1

Notes: UG_L = NA = Qual = RC = Micrograms per liter Not analyzed Final qualifier (Refer to Attachment B) Reason code (Refer to Attachment C)



Data Validation Report

Project:	Regional Groundwater Investigation - I	NWIRP Bethpage
Laboratory:	Katahdin Analytical	
Service Request:	TH0114	
Analyses/Method:	EPA SW-846 Method 9060A for TOC a Organic Carbon by High-Temperature	
Validation Level:	3	
AECOM Project Number:	60266526.SA.DV	
Prepared by:	Dawn Brule/RESCON	Completed on: 01/06/2015
Reviewed by:	Lori Herberich/RESCON	File Name: TH0114_5310B and 9060A

SUMMARY

The samples listed below were collected by Resolution Consultants from the Regional Groundwater Investigation - NWIRP Bethpage site on November 18, 2014.

Sample ID	Matrix/Sample Type
BPOW 6-3-EQUIPMENT BLANK-111814	Equipment blank
BPOW 6-3-SOIL-D-111814-753-755	Field Duplicate of BPOW 6-3-SOIL-111814-753-755
BPOW 6-3-SOIL-111814-753-755	Soil

The samples were analyzed in accordance with:

- Standard Methods for the Examination of Water and Wastewater, Method SM5310B, Total Organic Carbon by High-Temperature Combustion
- Test Methods for Evaluating Solid Waste, Physical/Chemical Methods SW-846, Method 9060A, Total Organic Carbon (USEPA, 1996).

Data validation activities were conducted with reference to these methods, *USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Superfund Data Review* (January 2010), and Quality Systems Manual (QSM) for Environmental Laboratories, Version 4.2 (DoD, October 2010) where applicable. 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 review elements (where applicable to the method):

- ✓ Data completeness (chain-of-custody [COC])/sample integrity
- ✓ Holding times and sample preservation
- ✓ Initial calibration/continuing calibration verification

- ✓ Laboratory blanks/equipment blanks
- Matrix spike (MS)/matrix duplicate (MD) and/or matrix spike duplicate (MSD) results
- ✓ Laboratory control sample (LCS) results
- ✓ Field duplicates
- ✓ 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. The symbol () indicates that a QC nonconformance resulted in the qualification of data. Any QC nonconformance that resulted in the qualification of data is discussed below. In addition, nonconformances or other issues that were noted during validation, but did not result in qualification of data, may be discussed for informational purposes only.

The data appear valid as reported and may be used for decision making purposes. There were no data points qualified or rejected on the basis of this data review.

RESULTS

Data Completeness (COC)/Sample Integrity

The data package was reviewed and found to meet acceptance criteria for completeness:

- The COCs were reviewed for completeness of information relevant to the samples and requested analyses, and for signatures indicating transfer of sample custody.
- The laboratory sample login sheet(s) were reviewed for issues potentially affecting sample integrity, including the condition of sample containers upon receipt at the laboratory.
- Completeness of analyses was verified by comparing the reported results to the COC requests.

Due to limitations in the reporting system, the laboratory truncated the IDs for all the samples in the report. The submitted EDD file reflects the full sample ID.

Holding Times/Sample Preservation

Sample preservation and preparation/analysis holding times were reviewed for conformance with the QC acceptance criteria. The QC acceptance criteria were met.

Initial Calibration/Continuing Calibration Verification

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

- all criteria were met for the calibration curves
- the initial calibration verification (ICV) percent recovery (%R) criteria were met; and
- the continuing calibration verification standard (CCV) method %Rs were met

The QC acceptance criteria were met.

Laboratory Blanks/Equipment Blanks

Laboratory method blanks and equipment rinsate blanks were evaluated as to whether there were contaminants detected above the detection limit (DL).

Data validation qualifications for individual samples are based on the maximum contaminant concentration detected in all associated blanks.

Method and equipment rinsate results were reviewed for conformance with the QC acceptance criteria. Detected results in blanks are not discussed in this data validation report if the associated results were nondetect or if qualification of sample results was not required.

The QC acceptance criteria were met and/or qualification of the sample results was not required.

MS/MSD Results

All MS/MD/MSD criteria were met and/or qualification of the data was not required.

LCS Results

The LCS %Rs were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

Field Duplicate Results

Field duplicate RPDs were reviewed for conformance with the QC criterion of ≤60% for soil matrices. This criteria applies if both results were greater than five times the Limit of Quantitation (LOQ). All QC acceptance criteria were met.

Sample Results/Reporting Issues

Compounds that were not detected in the sample are reported as not detected (U) at the Limit of Detection (LOD).

Compounds detected at concentrations less than the LOQ but greater than the detection limit (DL) were qualified by the laboratory as estimated (J). This "J" qualifier was retained during data validation.

Any sample that was analyzed at a dilution due to high concentrations of target or non-target compounds or matrix interferences was checked to ensure that the results and/or sample specific LODs and LOQs were adjusted accordingly by the laboratory.

QUALIFICATION ACTIONS

No sample results were qualified as a result of this data review.



Resolution

Suite B-1

Valente

☐ FED EX

Sample Description

BPOW 6-3 - Equipmen RINSote Blank-1/1814

1000 6-3-501 - Find

BPOW6-3-5011-

111814-753-755

TEMP BLANK

WORK ORDER #:

100 Red Schoolhouse Ril

Client

Address

Purchase Order #

Bill (if different than above)

Sampler (Print / Sign)

LAB USE ONLY

REMARKS:

SHIPPING INFO:

AIRBILL NO: __ TEMP'C -

600 Technology Way Scarborough, ME 04074 Tel: (207) 874-2400 Fax: (207) 775-4029

Consulta

KATAHDIN PROJECT NUMBER THO 11 4

D UPS

□ INTACT

Date / Time

coll'd

NOVIS

2014/

Contact

Address

City

Proj. Name / No.

CLIENT

Matrix

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O NOT INTACT

No. of Cntrs.

Eleanor Vivaudou

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http://www.katahdinlab.com



Report of Analytical Results

Client: Rick Purdy

AECOM

Wakefield,MA 01880 701 Edgewater Drive

Lab Sample ID: TH0114-1

Report Date: 15-DEC-14 Client PO: 16518

Project: Navy Clean WE15-03-0 SDG: TH0114

Date Received 26-NOV-14 18-NOV-14 09:45:00 Date Sampled Matrix AQ

BPOW6-3-EB-111814 Sample Description

Prep. Date Footnotes X/A Prep. Method N/A WG155518 10-DEC-14 18:42:12 Anal. Date QC.Batch Anal. Method SM5310B Adj LOQ Adj MDL Adj LOD 0.10 0.1 J0.35 mg/L Result Total Organic Carbon Parameter

Report of Analytical Results

/W Katahdin ANALYTICAL SERVICES

Client: Rick Purdy

AECOM

701 Edgewater Drive Wakefield,MA 01880

Lab Sample ID: TH0114-2 Report Date: 15-DEC-14

Client PO: 16518

Project: Navy Clean WE15-03-0 SDG: TH0114

Date Received 26-NOV-14 18-NOV-14 09:35:00 Date Sampled Matrix SL OW6-3-1118-753-755 Sample Description

8		
Footnotes		
Prep. Date	N/A	03-DEC-14
Prep. Method	N/A	SM2540G
Anal. Date	09-DEC-14 14:23:47	04-DEC-14 11:13:15
QC.Batch	WG155500 (WG155033
Adj MDL Adj LOD Anal. Method QC.Batch	SW846 9060A W Mod.	SM2540G
Adj LOD	550	N/A
Adj MDL	150	
Adj LOQ	730	-
Result	1200 ug/gdrywt	67. %

TOC In Soil

Parameter

Total Solids



Report of Analytical Results

Client: Rick Purdy

AECOM

Wakefield,MA 01880 701 Edgewater Drive

Lab Sample ID: TH0114-3

Report Date: 15-DEC-14 Client PO: 16518

Project: Navy Clean WE15-03-0 SDG: TH0114

Date Received 26-NOV-14 18-NOV-14 09:35:00 Date Sampled Matrix SI OW6-3-1118-753-755D Sample Description

Result	Adj 1.00	Adj MDL	Adj LOD	LOQ Adj MDL Adj LOD Anal, Method QC.Batch	QC.Batch	Anal. Date	Prep. Method Prep. Date Footnotes	Prep. Date	Footnotes	
1300 ug/gdrywt	1200	250	880	SW846 9060A Mod.	WG155500	SW846 9060A WG155500 09-DEC-14 15:06:16 Mod.	N/A	N/A		ł
78.%	-		N/A	SM2540G	WG155033	WG155033 04-DEC-14 11:13:30	SM2540G	03-DEC-14		

TOC la Soil Parameter

Total Solids

http://www.katahdinlab.com



DATA VALIDATION REPORT

Project:	Regional Groundwater Investigat	ion — NWIRP Bethpage
Laboratory:	Katahdin Analytical	
Sample Delivery Group:	TH0531	
Analyses/Method:	Total Organic Carbon by U.S. EPA	A SW-846 Method 9060A
Validation Level:	3	
Project Number:	0888812477.SA.DV	
Prepared by:	Dana Miller/Resolution Consultants Completed on: 02/06/2015 Revised on: 10/27/2015	
Reviewed by:	Tina Cantwell/Resolution Consultants	File Name: TH0531_9060A

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 December 2014 in accordance with the following Sampling and Analysis Plans:

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

Sample ID	Matrix/Sample Type	Analysis
BPOW6-4-SOIL-120514-563-565	Soil	9060A
BPOW6-4-SOIL-D-120814	Soil	9060A
BPOW6-4-EB-120814	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), *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 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 FLEMENTS

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

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

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

Field Duplicate

Field duplicate RPDs were reviewed for conformance with the Resolution Consultants QC criteria of $\leq 50\%$ for solid matrices and $\leq 30\%$ for aqueous matrices. These criteria apply if both results were greater than two times the limit of quantitation (LOQ). Non-conformance is summarized in Attachment A in Table A-1. Data qualification to the analytes associated with the specific field duplicate RPDs was as follows:



Field Duplicate Non-conformances Chart:

Criteria	RPD	Action		
Criteria	KPD	Detected	Non-detected	
Sample and duplicate are nondetect	Not calculable (NC)	No qualification	No qualification	
Sample and duplicate results ≥2x LOQ	>30 (aqueous)	-30 (aqueous)		
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	IJ	
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

Qualifications Actions

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

No results were rejected; therefore, analytical completeness was calculated to be 100 percent. Data not qualified during data review are considered usable by the project. The remaining results qualified as estimated may be high or low, but the data are usable for their intended purpose, according to U.S. 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. Katahdin Analytical discovered a QC error in Total Organic Carbon calculations, informed us, and has implemented corrective action provided in Attachment D. Attachment E 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: Katahdin Analytical Corrective Action Report

Attachment E: Final Results after Data Review

Attachment A Non-Conformance Summary Tables

	Field Dup	Table A-1 licate Non-Co	nformance				
Sample ID	Duplicate ID	Compound	Sample Result (UG_G)	Duplicate Result (UG_G)	RPD	RPD Limit	Qualifiers
BPOW6-4-SOIL-120514- 563-565	BPOW6-4-SOIL-D-120814	TOC	1400	720	64.2	50	J-both

Notes:

ID = Identification

TOC = Total organic carbon

UG_G = Micrograms per gram

RPD = Relative percent difference

J = 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
bt	Trip blank contamination
С	Calibration issue
d	Reporting limit raised due to chromatographic interference
fd	Field duplicate relative percent difference
h	Holding times
i	Internal standard areas
k	Estimated Maximum Possible Concentration
I	Laboratory control sample
lc	Labeled compound recovery
ld	Laboratory duplicate relative percent difference
lp	Laboratory control sample/laboratory control sample duplicate relative percent difference
m	Matrix spike recovery
mc	Method compliance non-conformance
md	Matrix spike/matrix spike duplicate relative percent difference
nb	Negative laboratory blank contamination
р	Chemical preservation issue
r	Dual column relative percent difference
q	Quantitation issue
S	Surrogate recovery
su	Ion suppression
t	Temperature preservation issue
Х	Percent solids
У	Serial dilution results
Z	Interference check sample results (metals)

Attachment D Katahdin Analytical Corrective Action Report

KATAHDIN ANALYTICAL SERVICES, INC. - CORRECTIVE ACTION REPORT

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

KATAHDIN ANALYTICAL SERVICES, INC. - CORRECTIVE ACTION REPORT

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

Corrective Action Plan: Name: Leslie Dimond

Date: 08/28/2015

Details of Corrective Action Plan -

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

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

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

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

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

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

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

KATAHDIN ANALYTICAL SERVICES, INC. – CORRECTIVE ACTION REPORT

senior management or the Quality Assurance Officer must also check the new calculations.								
Review & Approval of Corrective Action Plan								
Supervisor Approval:		Date: 39・11・15						
Operations Manager Approval:	lan	Date: 9:11-15						
Quality Assurance Officer: Luseic T		Date: 09-11-15						
Monitoring of Corrective Action (To be completed by QA Officer and/or Operations Manager): List details of follow-up								
Corrective Action Effective	Return to Control –	Yes No	Further Monitoring Needed/Additional Corrective Action					
QA Approval:			Date:					
Additional Information								
Additional Information:								

Attachment E Final Results after Data Review

	S	ery Group	TH0531			TH0531			TH0531			
Lab ID			TH0531-1			TH0531-2			TH0531-3			
Sample ID			BP0W6-4-SOII	BPOW6-4-SOIL-120814-563-565 BPOW6-4-SOIL-D-120814			314	BPOW6-4-EB-120814				
	Sample Date			12/8/2014			12/	8/2014		12/8/2014		
		Sa	mple Type	Soil			Field	Duplicate		Equipment Blank		
Method	Analyte	CAS No	Units	Result	Qual	RC	Result	Qual	RC	Result	Qual	RC
5310B	5310B TOTAL ORGANIC CARBON -28		MG_L	NA			NA			0.34	J	
9060A	TOTAL ORGANIC CARBON	-28	UG_G	1400	J	fd	720	J	fd	NA		

Notes:

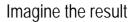
MG_L = UG_G =

NA

Milligrams per liter Micrograms per gram Not analyzed Final qualifier (Refer to Attachment B) Reason code (Refer to Attachment C) Qual RC =

Section 6

Analytical Data Validation - Arcadis





Northrop Grumman Corporation- Operable Unit 2

Data Review

BETHPAGE, NEW YORK

Volatile Analysis

SDG #JB97683

Analyses Performed By: Accutest Laboratories Dayton, New Jersey

Report #23906R July 17, 2015 Review Level: Tier II Project #NY001496.1514.NAVI4

SUMMARY

This data quality assessment summarizes the review of Sample Delivery Group (SDG) #JB97683 for samples collected in association with the Northrop Grumman-Bethpage Site. The review was conducted as a Tier II evaluation and included review of data package completeness. Only analytical data associated with constituents of concern were reviewed for this validation. Field documentation was not included in this review. Included with this assessment are the validation annotated sample result sheets, and chain of custody. Analyses were performed on the following samples:

			Sample			,	Analysis		
Sample ID	Lab ID	Matrix	Collection Date	Parent Sample	voc	svoc	РСВ	MET	MISC
BPOW 6-1	JB97683-1	Water	06/22/2015		Х				
TB062215KV1	JB97683-2	Water	06/22/2015		Х				
FB062215KV1	JB97683-3	Water	06/22/2015		Х				

ANALYTICAL DATA PACKAGE DOCUMENTATION

GENERAL INFORMATION

Items Reviewed	Rep	orted		mance ptable	Not	
	No	Yes	No	Yes	Required	
1. Sample receipt condition		Х		Х		
2. Requested analyses and sample results		Х		Х		
3. Collection Technique (grab, composite, etc.)		Х		Х		
4. Methods of analysis		Х		Х		
5. Reporting limits		Х		Х		
6. Sample collection date		Х		Х		
7. Laboratory sample received date		Х	Х			
8. Sample preservation verification (as applicable)		Х		X		
9. Sample preparation/extraction/analysis dates		Х		Х		
10. Fully executed Chain-of-Custody (COC) form completed		Х		Х		
11. Narrative summary of QA or sample problems provided		Х		Х		
12. Data Package Completeness and Compliance		Х		Х		

QA - Quality Assurance

VOLATILE ANALYSIS INTRODUCTION

Analyses were performed according to United States Environmental Protection Agency (USEPA), Method 524.2. Data were reviewed in accordance with USEPA National Functional Guidelines of October 1999.

The data review process is an evaluation of data on a technical basis rather than a determination of contract compliance. As such, the standards against which the data are being weighed may differ from those specified in the analytical method. It is assumed that the data package represents the best efforts of the laboratory and had already been subjected to adequate and sufficient quality review prior to submission.

During the review process, laboratory qualified and unqualified data are verified against the supporting documentation. Based on this evaluation, qualifier codes may be added, deleted, or modified by the data reviewer. Results are qualified with the following codes in accordance with USEPA National Functional Guidelines:

- Concentration (C) Qualifiers
 - U The compound was analyzed for but not detected. The associated value is the compound quantitation limit.
 - B The compound has been found in the sample as well as its associated blank, its presence in the sample may be suspect.
- Quantitation (Q) Qualifiers
 - E The compound was quantitated above the calibration range.
 - D Concentration is based on a diluted sample analysis.
- Validation Qualifiers
 - J The compound was positively identified; however, the associated numerical value is an estimated concentration only.
 - UJ The compound was not detected above the reported sample quantitation limit. However, the reported limit is approximate and may or may not represent the actual limit of quantitation.
 - JN The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification. The associated numerical value is an estimated concentration only.
 - UB Compound considered non-detect at the listed value due to associated blank contamination.
 - N The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification.
 - R The sample results are rejected.

Two facts should be noted by all data users. First, the "R" flag means that the associated value is unusable. In other words, due to significant quality control (QC) problems, the analysis is invalid and provides no information as to whether the compound is present or not. "R" values should not appear on data tables because they cannot be relied upon, even as a last resort. The second fact to keep in mind is that no compound concentration, even if it has passed all QC tests, is guaranteed to be accurate. Strict QC serves to increase confidence in data but any value potentially contains error.

VOLATILE ORGANIC COMPOUND (VOC) ANALYSES

1. Holding Times

The specified holding times for the following methods are presented in the following table.

Method	Matrix	Holding Time	Preservation
EPA 524.2	Water	14 days from collection to analysis	Cool to < 6°C; preserved to a pH of less than 2 s.u

s.u. Standard units

All samples were analyzed within the specified holding time and temperature criteria.

2. Blank Contamination

Quality assurance (QA) blanks (i.e., method and rinse blanks) are prepared to identify any contamination which may have been introduced into the samples during sample preparation or field activity. Method blanks measure laboratory contamination. Rinse blanks measure contamination of samples during field operations.

A blank action level (BAL) of five times the concentration of a detected compound in an associated blank (common laboratory contaminant compounds are calculated at ten times) is calculated for QA blanks containing concentrations greater than the method detection limit (MDL). The BAL is compared to the associated sample results to determine the appropriate qualification of the sample results, if needed.

All compounds associated with the QA blanks exhibited a concentration less than the MDL, with the exception of the compounds listed in the following table. Sample results associated with QA blank contamination that were greater than the BAL resulted in the removal of the laboratory qualifier (B) of data. Sample results less than the BAL associated with the following sample locations were qualified as listed in the following table.

Sample Locations	Analytes	Sample Result	Qualification
PDOW 6.1	Acetone	Detected sample results >RL and <bal< td=""><td>"UB" at detected sample concentration</td></bal<>	"UB" at detected sample concentration
BPOW 6-1	TICs: Isopropyl alcohol (RT: 7.46) Propanal, 2-methyl (RT: 8.71)	Detected sample results less than 5 times blank result	R

RL Reporting limit

3. Surrogates/System Monitoring Compounds

All samples to be analyzed for organic compounds are spiked with surrogate compounds prior to sample preparation to evaluate overall laboratory performance and efficiency of the analytical technique. VOC analysis requires that all surrogates associated with the analysis exhibit recoveries within the laboratory-established acceptance limits.

All surrogate recoveries were within control limits.

4. Matrix Spike/Matrix Spike Duplicate (MS/MSD) Analysis

MS/MSD data are used to assess the precision and accuracy of the analytical method. The compounds used to perform the MS/MSD analysis must exhibit a percent recovery within the laboratory-established acceptance limits. The relative percent difference (RPD) between the MS/MSD recoveries must exhibit an RPD within the laboratory-established acceptance limits.

Note: The MS/MSD recovery control limits do not apply for MS/MSD performed on sample locations where the compound concentration detected in the parent sample exceeds the MS/MSD concentration by a factor of four or greater.

A MS analysis was not performed on a sample location associated with this SDG.

5. Laboratory Control Sample (LCS) Analysis

The LCS analysis is used to assess the precision and accuracy of the analytical method independent of matrix interferences. The compounds associated with the LCS analysis must exhibit a percent recovery within the laboratory-established acceptance limits.

All compounds associated with the LCS analysis exhibited recoveries within the control limits.

6. Field Duplicate Analysis

Field duplicate analysis is used to assess the overall precision of the field sampling procedures and analytical method. A control limit of 30% for water matrices is applied to the RPD between the parent sample and the field duplicate. In the instance when the parent and/or duplicate sample concentrations are less than or equal to 5 times the RL, a control limit of two times the RL is applied for water matrices.

A field duplicate was not collected on a sample location associated with this SDG.

7. Laboratory Duplicate Analysis

The laboratory duplicate relative percent difference (RPD) criterion is applied when parent and duplicate sample concentrations are greater than or equal to 5 times the RL. A control limit of 20% for water matrices is applied when the criteria above is true. In the instance when the parent and/or duplicate sample concentrations are less than or equal to 5 times the RL, a control limit of two times the RL is applied for water matrices.

A laboratory duplicate was not performed on a sample location associated with this SDG.

8. System Performance and Overall Assessment

Tentatively identified compounds (TICs) were identified in sample locations BPOW 6-1 and FB062215KV1. The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification. TICs are qualified as estimated (JN).

Overall system performance was noted by the laboratory in the case narrative as acceptable. Other than for those deviations specifically mentioned in this review, the overall data quality is within the guidelines specified in the method.

DATA VALIDATION CHECKLIST FOR VOCs

VOCs: 524.2	Rep	orted		mance ptable	Not Required			
	No	Yes	No	Yes	Required			
GAS CHROMATOGRAPHY/MASS SPECTROMETRY (GC/MS)								
Tier II Validation								
Holding times & Temperature		Х		Х				
Reporting limits (units)		Х		Х				
Blanks								
A. Method blanks		Х		Х				
B. Equipment blanks		Х	Х					
C. Trip blanks		Х	Х					
Surrogate (%R)		Х		Х				
Laboratory Control Sample (%R)		Х		Х				
Laboratory Control Sample Duplicate(LCSD)					Х			
LCS/LCSD Precision (RPD)					Х			
Matrix Spike (MS)					Х			
Matrix Spike Duplicate(MSD)					Х			
MS/MSD Precision (RPD)					Х			
Field/Lab Duplicate (RPD)					Х			
Dilution Factor		Х		Х				
Moisture Content					Х			

[%]R Percent Recovery RPD Relative Percent Difference

VALIDATION PERFORMED BY: Lisa Horton

SIGNATURE: XISA TTORION

DATE: _ July 17, 2015

PEER REVIEW BY: Todd Church

DATE: July 20, 2015

CHAIN OF CUSTODY/ CORRECTED SAMPLE ANALYSIS DATA SHEETS

GWBB

CHAIN OF CUSTODY Accutest New Jersey/SPL Environmental 2235 Route 130 Dayton NI 08810

PAGE ___OF ___

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JB97683: Chain of Custody Page 1 of 3



Page 1 of 2

Date Sampled: 06/22/15

Report of Analysis

Client Sample ID: BPOW 6-1 Lab Sample ID: JB97683-1

 Matrix:
 AQ - Water
 Date Received:
 06/23/15

 Method:
 EPA 524.2 REV 4.1
 Percent Solids:
 n/a

Project: Northrop Grumman, OU2 Hydro, Bethpage, NY

File ID DF Analyzed By Prep Date Prep Batch Analytical Batch
Run #1 1B97958.D 1 06/29/15 MD n/a n/a V1B4647

Run #2

Purge Volume

Run #1 5.0 ml

Run #2

Special VOA List

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	9.3	5.0	0.91	ug/l	UB
78-93-3	2-Butanone	ND	5.0	0.57	ug/l	
71-43-2	Benzene	ND	0.50	0.057	ug/l	
75-27-4	Bromodichloromethane	ND	0.50	0.082	ug/l	
75-25-2	Bromoform	ND	0.50	0.046	ug/l	
74-83-9	Bromomethane	ND	0.50	0.077	ug/l	
75-15-0	Carbon disulfide	ND	0.50	0.028	ug/l	
108-90-7	Chlorobenzene	ND	0.50	0.027	ug/l	
75-00-3	Chloroethane	ND	0.50	0.037	ug/l	
67-66-3	Chloroform	ND	0.50	0.031	ug/l	
74-87-3	Chloromethane	0.51	0.50	0.044	ug/l	
56-23-5	Carbon tetrachloride	ND	0.50	0.074	ug/l	
75-34-3	1,1-Dichloroethane	ND	0.50	0.039	ug/l	
75-35-4	1,1-Dichloroethylene	ND	0.50	0.054	ug/l	
107-06-2	1,2-Dichloroethane	ND	0.50	0.034	ug/l	
78-87-5	1,2-Dichloropropane	ND	0.50	0.082	ug/l	
124-48-1	Dibromochloromethane	ND	0.50	0.042	ug/l	
156-60-5	trans-1,2-Dichloroethylene	ND	0.50	0.039	ug/l	
156-59-2	cis-1,2-Dichloroethylene	ND	0.50	0.081	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.033	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.063	ug/l	
100-41-4	Ethylbenzene	ND	0.50	0.033	ug/l	
76-13-1	Freon 113	ND	1.0	0.10	ug/l	
591-78-6	2-Hexanone	ND	2.0	0.084	ug/l	
75-09-2	Methylene chloride	ND	0.50	0.047	ug/l	
108-10-1	4-Methyl-2-pentanone	ND	2.0	0.27	ug/l	
100-42-5	Styrene	ND	0.50	0.028	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	0.50	0.050	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.035	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	0.50	0.052	ug/l	
127-18-4	Tetrachloroethylene	ND	0.50	0.091	ug/l	
108-88-3	Toluene	ND	0.50	0.044	ug/l	

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

N = Indicates presumptive evidence of a compound



Date Sampled: 06/22/15

Date Received: 06/23/15

n/a

Percent Solids:

Report of Analysis

Client Sample ID: BPOW 6-1 Lab Sample ID: JB97683-1

Matrix: AQ - Water Method: EPA 524.2 REV 4.1

Project: Northrop Grumman, OU2 Hydro, Bethpage, NY

Special VOA List

CAS No.	Compound	Result	RL	MDL	Units	Q		
79-01-6 75-01-4	Trichloroethylene Vinyl chloride m,p-Xylene	ND ND ND	0.50 0.50 0.50	0.024 0.032 0.13	ug/l ug/l ug/l			
95-47-6 CAS No.	o-Xylene Surrogate Recoveries	ND Run# 1	0.50 Run# 2	0.029 Limi	ug/l ts			
2199-69-1 460-00-4	1,2-Dichlorobenzene-d4 4-Bromofluorobenzene	93% 86%		78-1 77-1				
CAS No.	Tentatively Identified Compo	ounds	R.T.	Est.	Conc.	Units	Q	
67-63-0	Isopropyl Alcohol Propanal, methyl- Propanol, methyl- Furan, tetrahydro-tetramethyl- Hexanol, ethyl- Total TIC, Volatile		7.46 8.71 10.84 12.87 17.36 5.12	.98 2.1 2.2 .82 66.1		ug/l ug/l ug/l ug/l ug/l ug/l	JN J N J N J N J N	R R

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank N = Indicates presumptive evidence of a compound



Report of Analysis

Client Sample ID: FB062215KV1

Lab Sample ID:JB97683-2Date Sampled:06/22/15Matrix:AQ - Field Blank WaterDate Received:06/23/15Method:EPA 524.2 REV 4.1Percent Solids:n/a

Project: Northrop Grumman, OU2 Hydro, Bethpage, NY

File ID DF **Prep Date Analytical Batch** Analyzed By **Prep Batch** Run #1 1B97959.D 1 06/29/15 MD V1B4647 n/an/aRun #2

Purge Volume Run #1 5.0 ml

Run #2

Special VOA List

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	1.1	5.0	0.91	ug/l	J
78-93-3	2-Butanone	ND	5.0	0.57	ug/l	
71-43-2	Benzene	ND	0.50	0.057	ug/l	
75-27-4	Bromodichloromethane	ND	0.50	0.082	ug/l	
75-25-2	Bromoform	ND	0.50	0.046	ug/l	
74-83-9	Bromomethane	ND	0.50	0.077	ug/l	
75-15-0	Carbon disulfide	ND	0.50	0.028	ug/l	
108-90-7	Chlorobenzene	ND	0.50	0.027	ug/l	
75-00-3	Chloroethane	ND	0.50	0.037	ug/l	
67-66-3	Chloroform	0.15	0.50	0.031	ug/l	J
74-87-3	Chloromethane	ND	0.50	0.044	ug/l	
56-23-5	Carbon tetrachloride	ND	0.50	0.074	ug/l	
75-34-3	1,1-Dichloroethane	ND	0.50	0.039	ug/l	
75-35-4	1,1-Dichloroethylene	ND	0.50	0.054	ug/l	
107-06-2	1,2-Dichloroethane	ND	0.50	0.034	ug/l	
78-87-5	1,2-Dichloropropane	ND	0.50	0.082	ug/l	
124-48-1	Dibromochloromethane	ND	0.50	0.042	ug/l	
156-60-5	trans-1,2-Dichloroethylene	ND	0.50	0.039	ug/l	
156-59-2	cis-1,2-Dichloroethylene	ND	0.50	0.081	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.033	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.063	ug/l	
100-41-4	Ethylbenzene	ND	0.50	0.033	ug/l	
76-13-1	Freon 113	ND	1.0	0.10	ug/l	
591-78-6	2-Hexanone	ND	2.0	0.084	ug/l	
75-09-2	Methylene chloride	ND	0.50	0.047	ug/l	
108-10-1	4-Methyl-2-pentanone	ND	2.0	0.27	ug/l	
100-42-5	Styrene	ND	0.50	0.028	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	0.50	0.050	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.035	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	0.50	0.052	ug/l	
127-18-4	Tetrachloroethylene	ND	0.50	0.091	ug/l	
108-88-3	Toluene	ND	0.50	0.044	ug/l	

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

N = Indicates presumptive evidence of a compound



Page 2 of 2

4

Report of Analysis

Client Sample ID: FB062215KV1

Lab Sample ID:JB97683-2Date Sampled:06/22/15Matrix:AQ - Field Blank WaterDate Received:06/23/15Method:EPA 524.2 REV 4.1Percent Solids:n/a

Project: Northrop Grumman, OU2 Hydro, Bethpage, NY

Special VOA List

CAS No.	Compound	Result	RL	MDL	Units	Q	
79-01-6 75-01-4 95-47-6	Trichloroethylene Vinyl chloride m,p-Xylene o-Xylene	ND ND ND ND	0.50 0.50 0.50 0.50	0.024 0.032 0.13 0.029	ug/l ug/l ug/l ug/l		
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limi	ts		
2199-69-1 460-00-4	1,2-Dichlorobenzene-d4 4-Bromofluorobenzene	91% 86%		78-11 77-11			
CAS No.	Tentatively Identified Compo	ounds	R.T.	Est.	Conc.	Units	Q
67-63-0	Isopropyl Alcohol Propanal, methyl- Total TIC, Volatile		7.47 8.72	4.6 1.3 5.9		ug/l ug/l ug/l	JN J _N J _N

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank N = Indicates presumptive evidence of a compound



Report of Analysis

Client Sample ID: TB062215KV1

Lab Sample ID:JB97683-3Date Sampled:06/22/15Matrix:AQ - Trip Blank WaterDate Received:06/23/15Method:EPA 524.2 REV 4.1Percent Solids:n/a

Project: Northrop Grumman, OU2 Hydro, Bethpage, NY

	File ID	DF	Analyzed	$\mathbf{B}\mathbf{y}$	Prep Date	Prep Batch	Analytical Batch
Run #1	1B97907.D	1	06/27/15	MD	n/a	n/a	V1B4644
Dun #2							

	Purge Volume	
Run #1	5.0 ml	
Run #2		

Special VOA List

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	5.0	0.91	ug/l	
78-93-3	2-Butanone	ND	5.0	0.57	ug/l	
71-43-2	Benzene	ND	0.50	0.057	ug/l	
75-27-4	Bromodichloromethane	ND	0.50	0.082	ug/l	
75-25-2	Bromoform	ND	0.50	0.046	ug/l	
74-83-9	Bromomethane	ND	0.50	0.077	ug/l	
75-15-0	Carbon disulfide	ND	0.50	0.028	ug/l	
108-90-7	Chlorobenzene	ND	0.50	0.027	ug/l	
75-00-3	Chloroethane	ND	0.50	0.037	ug/l	
67-66-3	Chloroform	0.17	0.50	0.031	ug/l	J
74-87-3	Chloromethane	ND	0.50	0.044	ug/l	
56-23-5	Carbon tetrachloride	ND	0.50	0.074	ug/l	
75-34-3	1,1-Dichloroethane	ND	0.50	0.039	ug/l	
75-35-4	1,1-Dichloroethylene	ND	0.50	0.054	ug/l	
107-06-2	1,2-Dichloroethane	ND	0.50	0.034	ug/l	
78-87-5	1,2-Dichloropropane	ND	0.50	0.082	ug/l	
124-48-1	Dibromochloromethane	ND	0.50	0.042	ug/l	
156-60-5	trans-1,2-Dichloroethylene	ND	0.50	0.039	ug/l	
156-59-2	cis-1,2-Dichloroethylene	ND	0.50	0.081	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.033	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.063	ug/l	
100-41-4	Ethylbenzene	ND	0.50	0.033	ug/l	
76-13-1	Freon 113	ND	1.0	0.10	ug/l	
591-78-6	2-Hexanone	ND	2.0	0.084	ug/l	
75-09-2	Methylene chloride	ND	0.50	0.047	ug/l	
108-10-1	4-Methyl-2-pentanone	ND	2.0	0.27	ug/l	
100-42-5	Styrene	ND	0.50	0.028	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	0.50	0.050	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.035	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	0.50	0.052	ug/l	
127-18-4	Tetrachloroethylene	ND	0.50	0.091	ug/l	
108-88-3	Toluene	ND	0.50	0.044	ug/l	

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank N = Indicates presumptive evidence of a compound



Page 2 of 2

Report of Analysis

Client Sample ID: TB062215KV1

Lab Sample ID: JB97683-3 **Date Sampled:** 06/22/15 Matrix: AQ - Trip Blank Water **Date Received:** 06/23/15 Method: EPA 524.2 REV 4.1 **Percent Solids:** n/a

Project: Northrop Grumman, OU2 Hydro, Bethpage, NY

Special VOA List

CAS No.	Compound	Result	RL	MDL	Units	Q	
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l		
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l		
	m,p-Xylene	ND	0.50	0.13	ug/l		
95-47-6	o-Xylene	ND	0.50	0.029	ug/l		
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limi	its		
		2221		=0.4			
2199-69-1	1,2-Dichlorobenzene-d4	88%		78-1	,.		
460-00-4	4-Bromofluorobenzene	88%		77-1	15%		
		_		_			_
CAS No.	Tentatively Identified Compo	ounds	R.T.	Est.	Conc.	Units	Q
	m . 1 m/G . 1/1 . 11			0			
	Total TIC, Volatile			0		ug/l	

ND = Not detected MDL = Method Detection Limit

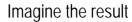
RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank N = Indicates presumptive evidence of a compound







Northrop Grumman Corporation- Operable Unit 2

Data Review

BETHPAGE, NEW YORK

Volatile Analysis

SDG #JB97745

Analyses Performed By: Accutest Laboratories Dayton, New Jersey

Report #23913R July 10, 2015 Review Level: Tier II Project #NY001496.1514.NAVI4

SUMMARY

This data quality assessment summarizes the review of Sample Delivery Group (SDG) #JB97745 for samples collected in association with the Northrop Grumman-Bethpage Site. The review was conducted as a Tier II evaluation and included review of data package completeness. Only analytical data associated with constituents of concern were reviewed for this validation. Field documentation was not included in this review. Included with this assessment are the validation annotated sample result sheets, and chain of custody. Analyses were performed on the following samples:

			Sample Collection Parent Sample			,	Analysis	S	
Sample ID	Lab ID	Matrix	Date	Parent Sample	voc	svoc	РСВ	MET	MISC
BPOW 6-2	JB97745-1	Water	06/23/2015		Х				
FB062315KV1	JB97745-2	Water	06/23/2015		Χ				
TB062315KV1	JB97745-3	Water	06/23/2015		Х				
BPOW 6-R	JB97745-4	Water	06/23/2015	BPOW 6-2	Х				

ANALYTICAL DATA PACKAGE DOCUMENTATION

GENERAL INFORMATION

Items Reviewed	Reported		Performance Acceptable		Not
	No	Yes	No	Yes	Required
1. Sample receipt condition		Х		Х	
2. Requested analyses and sample results		Х		Х	
3. Collection Technique (grab, composite, etc.)		Х		Х	
4. Methods of analysis		Х		Х	
5. Reporting limits		Х		Х	
6. Sample collection date		Х		Х	
7. Laboratory sample received date		Х	Х		
8. Sample preservation verification (as applicable)		Х		X	
9. Sample preparation/extraction/analysis dates		Х		Х	
10. Fully executed Chain-of-Custody (COC) form completed		Х		Х	
11. Narrative summary of QA or sample problems provided		Х		Х	
12. Data Package Completeness and Compliance		Х		Х	

QA - Quality Assurance

VOLATILE ANALYSIS INTRODUCTION

Analyses were performed according to United States Environmental Protection Agency (USEPA), Method 524.2. Data were reviewed in accordance with USEPA National Functional Guidelines of October 1999.

The data review process is an evaluation of data on a technical basis rather than a determination of contract compliance. As such, the standards against which the data are being weighed may differ from those specified in the analytical method. It is assumed that the data package represents the best efforts of the laboratory and had already been subjected to adequate and sufficient quality review prior to submission.

During the review process, laboratory qualified and unqualified data are verified against the supporting documentation. Based on this evaluation, qualifier codes may be added, deleted, or modified by the data reviewer. Results are qualified with the following codes in accordance with USEPA National Functional Guidelines:

- Concentration (C) Qualifiers
 - U The compound was analyzed for but not detected. The associated value is the compound quantitation limit.
 - B The compound has been found in the sample as well as its associated blank, its presence in the sample may be suspect.
- Quantitation (Q) Qualifiers
 - E The compound was quantitated above the calibration range.
 - D Concentration is based on a diluted sample analysis.
- Validation Qualifiers
 - J The compound was positively identified; however, the associated numerical value is an estimated concentration only.
 - UJ The compound was not detected above the reported sample quantitation limit. However, the reported limit is approximate and may or may not represent the actual limit of quantitation.
 - JN The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification. The associated numerical value is an estimated concentration only.
 - UB Compound considered non-detect at the listed value due to associated blank contamination.
 - N The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification.
 - R The sample results are rejected.

Two facts should be noted by all data users. First, the "R" flag means that the associated value is unusable. In other words, due to significant quality control (QC) problems, the analysis is invalid and provides no information as to whether the compound is present or not. "R" values should not appear on data tables because they cannot be relied upon, even as a last resort. The second fact to keep in mind is that no compound concentration, even if it has passed all QC tests, is guaranteed to be accurate. Strict QC serves to increase confidence in data but any value potentially contains error.

VOLATILE ORGANIC COMPOUND (VOC) ANALYSES

1. Holding Times

The specified holding times for the following methods are presented in the following table.

Method	Matrix	Holding Time	Preservation
EPA 524.2	Water	14 days from collection to analysis	Cool to < 6°C; preserved to a pH of less than 2 s.u

s.u. Standard units

All samples were analyzed within the specified holding time and temperature criteria.

2. Blank Contamination

Quality assurance (QA) blanks (i.e., method and rinse blanks) are prepared to identify any contamination which may have been introduced into the samples during sample preparation or field activity. Method blanks measure laboratory contamination. Rinse blanks measure contamination of samples during field operations.

A blank action level (BAL) of five times the concentration of a detected compound in an associated blank (common laboratory contaminant compounds are calculated at ten times) is calculated for QA blanks containing concentrations greater than the method detection limit (MDL). The BAL is compared to the associated sample results to determine the appropriate qualification of the sample results, if needed.

All compounds associated with the QA blanks exhibited a concentration less than the MDL, with the exception of the compounds listed in the following table. Sample results associated with QA blank contamination that were greater than the BAL resulted in the removal of the laboratory qualifier (B) of data. Sample results less than the BAL associated with the following sample locations were qualified as listed in the following table.

Sample Locations	Analytes	Sample Result	Qualification
BPOW 6-2 BPOW 6-R	Acetone	Detected sample results >RL and <bal< td=""><td>"UB" at detected sample concentration</td></bal<>	"UB" at detected sample concentration
BPOW 6-R	TICs: Isopropyl alcohol (RT: 7.46/7.47)	Detected sample results less than 5 times blank result	R

RL Reporting limit

3. Surrogates/System Monitoring Compounds

All samples to be analyzed for organic compounds are spiked with surrogate compounds prior to sample preparation to evaluate overall laboratory performance and efficiency of the analytical technique. VOC analysis requires that all surrogates associated with the analysis exhibit recoveries within the laboratory-established acceptance limits.

All surrogate recoveries were within control limits.

4. Matrix Spike/Matrix Spike Duplicate (MS/MSD) Analysis

MS/MSD data are used to assess the precision and accuracy of the analytical method. The compounds used to perform the MS/MSD analysis must exhibit a percent recovery within the laboratory-established acceptance limits. The relative percent difference (RPD) between the MS/MSD recoveries must exhibit an RPD within the laboratory-established acceptance limits.

Note: The MS/MSD recovery control limits do not apply for MS/MSD performed on sample locations where the compound concentration detected in the parent sample exceeds the MS/MSD concentration by a factor of four or greater.

A MS analysis was not performed on a sample location associated with this SDG.

5. Laboratory Control Sample (LCS) Analysis

The LCS analysis is used to assess the precision and accuracy of the analytical method independent of matrix interferences. The compounds associated with the LCS analysis must exhibit a percent recovery within the laboratory-established acceptance limits.

All compounds associated with the LCS analysis exhibited recoveries within the control limits.

6. Field Duplicate Analysis

Field duplicate analysis is used to assess the overall precision of the field sampling procedures and analytical method. A control limit of 30% for water matrices is applied to the RPD between the parent sample and the field duplicate. In the instance when the parent and/or duplicate sample concentrations are less than or equal to 5 times the RL, a control limit of two times the RL is applied for water matrices.

Results for duplicate samples are summarized in the following table.

Sample ID/Duplicate ID	Compound	Sample Result	Duplicate Result	RPD
BPOW 6-2/ BPOW 6-R	Acetone	2.0 J	2.9 J	AC

AC Acceptable

The calculated RPDs between the parent sample and field duplicate were acceptable.

7. Laboratory Duplicate Analysis

The laboratory duplicate relative percent difference (RPD) criterion is applied when parent and duplicate sample concentrations are greater than or equal to 5 times the RL. A control limit of 20% for water matrices is applied when the criteria above is true. In the instance when the parent and/or duplicate sample concentrations are less than or equal to 5 times the RL, a control limit of two times the RL is applied for water matrices.

A laboratory duplicate was not performed on a sample location associated with this SDG.

8. System Performance and Overall Assessment

Tentatively identified compounds (TICs) were identified in sample locations BPOW 6-2, FB062315KV1 and BPOW 6-R. The analysis indicates the presence of a compound for which there is presumptive evidence to

make a tentative identification. TICs are qualified as estimated (JN).

Overall system performance was noted by the laboratory in the case narrative as acceptable. Other than for those deviations specifically mentioned in this review, the overall data quality is within the guidelines specified in the method.

DATA VALIDATION CHECKLIST FOR VOCs

VOCs: 524.2	Rep	orted	Performance Acceptable		Not Required	
	No	Yes	No	Yes	Nequired	
GAS CHROMATOGRAPHY/MASS SPECTROME	TRY (GC/	MS)				
Tier II Validation						
Holding times & Temperature		Х		Х		
Reporting limits (units)		Х		Х		
Blanks						
A. Method blanks		Х		Х		
B. Equipment blanks		Х	Х			
C. Trip blanks		Х	Х			
Surrogate (%R)		Х		Х		
Laboratory Control Sample (%R)		Х		Х		
Laboratory Control Sample Duplicate(LCSD)					Х	
LCS/LCSD Precision (RPD)					Х	
Matrix Spike (MS)					Х	
Matrix Spike Duplicate(MSD)					Х	
MS/MSD Precision (RPD)					Х	
Field/Lab Duplicate (RPD)		Х		Х		
Dilution Factor		Х		Х		
Moisture Content					Х	

[%]R Percent Recovery RPD Relative Percent Difference

VALIDATION PERFORMED BY: Lisa Horton

SIGNATURE: Xisa Horland

DATE: _ July 10, 2015

PEER REVIEW BY: Todd Church

DATE: _ July 16, 2015

CHAIN OF CUSTODY/ CORRECTED SAMPLE ANALYSIS DATA SHEETS

ACC	UTEST.
	LABORATORIES

av
WTB
WFB

CHAIN OF CUSTODY

PAGE OF

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JB97745: Chain of Custody Page 1 of 2



Report of Analysis

Client Sample ID: BPOW 6-2 Lab Sample ID: JB97745-1

 Lab Sample ID:
 JB97745-1
 Date Sampled:
 06/23/15

 Matrix:
 AQ - Ground Water
 Date Received:
 06/24/15

 Method:
 EPA 524.2 REV 4.1
 Percent Solids:
 n/a

Project: Northrop Grumman, OU2 Hydro, Bethpage, NY

File IDDFAnalyzedByPrep DatePrep BatchAnalytical BatchRun #11B97914.D106/27/15MDn/an/aV1B4645

Run #2

Purge Volume

Run #1 5.0 ml

Run #2

Special VOA List

CAS No.	Compound	Result	RL	MDL	Units Q
67-64-1	Acetone 5.0	-2.0	5.0	0.91	ug/l → UB
78-93-3	2-Butanone	ND	5.0	0.57	ug/l
71-43-2	Benzene	ND	0.50	0.057	ug/l
75-27-4	Bromodichloromethane	ND	0.50	0.082	ug/l
75-25-2	Bromoform	ND	0.50	0.046	ug/l
74-83-9	Bromomethane	ND	0.50	0.077	ug/l
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74-87-3	Chloromethane	ND	0.50	0.044	ug/l
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75-34-3	1,1-Dichloroethane	ND	0.50	0.039	ug/l
75-35-4	1,1-Dichloroethylene	ND	0.50	0.054	ug/l
107-06-2	1,2-Dichloroethane	ND	0.50	0.034	ug/l
78-87-5	1,2-Dichloropropane	ND	0.50	0.082	ug/l
124-48-1	Dibromochloromethane	ND	0.50	0.042	ug/l
156-60-5	trans-1,2-Dichloroethylene	ND	0.50	0.039	ug/l
156-59-2	cis-1,2-Dichloroethylene	ND	0.50	0.081	ug/l
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.033	ug/l
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.063	ug/l
100-41-4	Ethylbenzene	ND	0.50	0.033	ug/l
76-13-1	Freon 113	ND	1.0	0.10	ug/l
591-78-6	2-Hexanone	ND	2.0	0.084	ug/l
75-09-2	Methylene chloride	ND	0.50	0.047	ug/l
108-10-1	4-Methyl-2-pentanone	ND	2.0	0.27	ug/l
100-42-5	Styrene	ND	0.50	0.028	ug/l
71-55-6	1,1,1-Trichloroethane	ND	0.50	0.050	ug/l
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.035	ug/l
79-00-5	1,1,2-Trichloroethane	ND	0.50	0.052	ug/l
127-18-4	Tetrachloroethylene	ND	0.50	0.091	ug/l
108-88-3	Toluene	ND	0.50	0.044	ug/l

ND = Not detected

MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

N = Indicates presumptive evidence of a compound



Date Sampled: 06/23/15

06/24/15

n/a

Date Received:

Percent Solids:

Report of Analysis

Client Sample ID: BPOW 6-2 Lab Sample ID: JB97745-1

Matrix: AQ - Ground Water
Method: EPA 524.2 REV 4.1

Project: Northrop Grumman, OU2 Hydro, Bethpage, NY

Special VOA List

CAS No.	Compound	Result	RL	MDL	Units	Q		
79-01-6	Trichloroethylene	ND ND	0.50	0.024	ug/l			
75-01-4	Vinyl chloride m,p-Xylene	ND ND	0.50 0.50	0.032 0.13	ug/l ug/l			
95-47-6	o-Xylene	ND	0.50	0.029	ug/l			
CAS No.	Surrogate Recoveries Run		Run# 2	Limi	ts			
2199-69-1	1,2-Dichlorobenzene-d4	90%		78-1	14%			
460-00-4	4-Bromofluorobenzene	87%		77-1	15%			
CAS No.	Tentatively Identified Compounds			Est.	Conc.	Units	Q	
67-63-0	Isopropyl Alcohol		7.46	1.9		ug/l	JN_	R
78-84-2	Propanal, 2-methyl-		8.71	.79		ug/l	JN	
	Total TIC, Volatile		0.79	2.69	-	ug/l	J N	

 $ND = Not detected \qquad MDL = 1$

MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

 $N = \ \ Indicates \ presumptive \ evidence \ of \ a \ compound$



Report of Analysis

Client Sample ID: FB062315KV1

Lab Sample ID:JB97745-2Date Sampled:06/23/15Matrix:AQ - Field Blank WaterDate Received:06/24/15Method:EPA 524.2 REV 4.1Percent Solids:n/a

Project: Northrop Grumman, OU2 Hydro, Bethpage, NY

File ID DF **Prep Date Analytical Batch** Analyzed By **Prep Batch** Run #1 1B97955.D 1 06/29/15 MD n/aV1B4647 n/aRun #2

Purge Volume

Run #1 5.0 ml

Run #2

Special VOA List

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	1.1	5.0	0.91	ug/l	J
78-93-3	2-Butanone	ND	5.0	0.57	ug/l	
71-43-2	Benzene	ND	0.50	0.057	ug/l	
75-27-4	Bromodichloromethane	ND	0.50	0.082	ug/l	
75-25-2	Bromoform	ND	0.50	0.046	ug/l	
74-83-9	Bromomethane	ND	0.50	0.077	ug/l	
75-15-0	Carbon disulfide	ND	0.50	0.028	ug/l	
108-90-7	Chlorobenzene	ND	0.50	0.027	ug/l	
75-00-3	Chloroethane	ND	0.50	0.037	ug/l	
67-66-3	Chloroform	0.13	0.50	0.031	ug/l	J
74-87-3	Chloromethane	ND	0.50	0.044	ug/l	
56-23-5	Carbon tetrachloride	ND	0.50	0.074	ug/l	
75-34-3	1,1-Dichloroethane	ND	0.50	0.039	ug/l	
75-35-4	1,1-Dichloroethylene	ND	0.50	0.054	ug/l	
107-06-2	1,2-Dichloroethane	ND	0.50	0.034	ug/l	
78-87-5	1,2-Dichloropropane	ND	0.50	0.082	ug/l	
124-48-1	Dibromochloromethane	ND	0.50	0.042	ug/l	
156-60-5	trans-1,2-Dichloroethylene	ND	0.50	0.039	ug/l	
156-59-2	cis-1,2-Dichloroethylene	ND	0.50	0.081	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.033	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.063	ug/l	
100-41-4	Ethylbenzene	ND	0.50	0.033	ug/l	
76-13-1	Freon 113	ND	1.0	0.10	ug/l	
591-78-6	2-Hexanone	ND	2.0	0.084	ug/l	
75-09-2	Methylene chloride	ND	0.50	0.047	ug/l	
108-10-1	4-Methyl-2-pentanone	ND	2.0	0.27	ug/l	
100-42-5	Styrene	ND	0.50	0.028	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	0.50	0.050	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.035	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	0.50	0.052	ug/l	
127-18-4	Tetrachloroethylene	ND	0.50	0.091	ug/l	
108-88-3	Toluene	ND	0.50	0.044	ug/l	

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

N = Indicates presumptive evidence of a compound



Page 2 of 2

Report of Analysis

Client Sample ID: FB062315KV1

Lab Sample ID: JB97745-2 **Date Sampled:** 06/23/15 Matrix: AQ - Field Blank Water **Date Received:** 06/24/15 Method: **Percent Solids:** EPA 524.2 REV 4.1 n/a

Project: Northrop Grumman, OU2 Hydro, Bethpage, NY

Special VOA List

CAS No.	Compound	Result	RL	MDL	Units	Q	
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l		
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l		
	m,p-Xylene	ND	0.50	0.13	ug/l		
95-47-6	o-Xylene	ND	0.50	0.029	ug/l		
CACNo	Cuma acta Decembra	Run# 1	Run# 2	Limi	4		
CAS No.	Surrogate Recoveries	Kun# 1	Kun# Z	LIIII	ıs		
2199-69-1	1,2-Dichlorobenzene-d4	98%		78-1	14%		
460-00-4	4-Bromofluorobenzene	87%		77-1	15%		
CAS No.	Tentatively Identified Compo	ounds	R.T.	Est.	Conc.	Units	Q
55.52. 0				4.5			T. T
67-63-0	Isopropyl Alcohol		7.47	1.7		ug/l	JN
	Total TIC, Volatile			1.7		ug/l	JΝ

ND = Not detected MDL = Method Detection Limit J = Indicates an estimated value

RL = Reporting Limit

E = Indicates value exceeds calibration range

B = Indicates analyte found in associated method blank N = Indicates presumptive evidence of a compound



Report of Analysis

Client Sample ID: TB062315KV1

Lab Sample ID:JB97745-3Date Sampled:06/23/15Matrix:AQ - Trip Blank WaterDate Received:06/24/15Method:EPA 524.2 REV 4.1Percent Solids:n/a

Project: Northrop Grumman, OU2 Hydro, Bethpage, NY

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B97956.D	1	06/29/15	MD	n/a	n/a	V1B4647
Run #2							

Purge Volume
Run #1 5.0 ml
Run #2

Special VOA List

Compound	Result	RL	MDL	Units	Q
Acetone	ND	5.0	0.91	ug/l	
2-Butanone	ND	5.0	0.57	ug/l	
Benzene	ND	0.50	0.057	ug/l	
Bromodichloromethane	ND	0.50	0.082	ug/l	
Bromoform	ND	0.50	0.046	ug/l	
Bromomethane	ND	0.50	0.077	ug/l	
Carbon disulfide	ND	0.50	0.028	ug/l	
Chlorobenzene	ND	0.50	0.027	ug/l	
Chloroethane	ND	0.50	0.037	ug/l	
Chloroform	0.17	0.50	0.031	ug/l	J
Chloromethane	ND	0.50	0.044	ug/l	
Carbon tetrachloride	ND	0.50	0.074	ug/l	
1,1-Dichloroethane	ND	0.50	0.039	ug/l	
1,1-Dichloroethylene	ND	0.50	0.054	ug/l	
1,2-Dichloroethane	ND	0.50	0.034	ug/l	
1,2-Dichloropropane	ND	0.50	0.082	ug/l	
Dibromochloromethane	ND	0.50	0.042	ug/l	
trans-1,2-Dichloroethylene	ND	0.50	0.039	ug/l	
cis-1,2-Dichloroethylene	ND	0.50	0.081	ug/l	
	ND	0.50	0.033	ug/l	
cis-1,3-Dichloropropene	ND	0.50	0.063	ug/l	
Ethylbenzene		0.50	0.033		
Freon 113		1.0	0.10		
2-Hexanone	ND	2.0	0.084	-	
	ND	0.50	0.047	-	
4-Methyl-2-pentanone	ND	2.0	0.27	ug/l	
Styrene	ND	0.50	0.028	ug/l	
	ND	0.50	0.050	ug/l	
	ND		0.035	-	
	ND		0.052	ug/l	
· ·	ND	0.50	0.091	ug/l	
Toluene	ND	0.50	0.044	ug/l	
	Acetone 2-Butanone Benzene Bromodichloromethane Bromoform Bromomethane Carbon disulfide Chlorobenzene Chloroethane Chloroform Chloromethane Carbon tetrachloride 1,1-Dichloroethane 1,1-Dichloroethylene 1,2-Dichloropropane Dibromochloromethane trans-1,2-Dichloroethylene trans-1,2-Dichloroethylene trans-1,3-Dichloropropene cis-1,3-Dichloropropene Ethylbenzene Freon 113 2-Hexanone Methylene chloride 4-Methyl-2-pentanone Styrene 1,1,1-Trichloroethane 1,1,2,2-Tetrachloroethane 1,1,2-Trichloroethane Tetrachloroethylene	Acetone 2-Butanone Benzene ND Bromodichloromethane ND Bromoform ND Bromomethane ND Carbon disulfide ND Chlorobenzene ND Chloroethane ND Carbon tetrachloride ND Carbon tetrachloride ND 1,1-Dichloroethane ND 1,2-Dichloroethane ND Dibromochloromethane ND Dibromochloromethane ND Cis-1,2-Dichloroethylene ND trans-1,2-Dichloroethylene ND trans-1,3-Dichloropropene ND Ethylbenzene ND Freon 113 ND 2-Hexanone ND Methylene chloride ND 1,1,1-Trichloroethane ND ND ND Styrene ND ND ND ND ND ND ND ND ND ND ND ND ND	Acetone 2-Butanone Benzene ND 5.0 Benzene ND 0.50 Bromodichloromethane ND 0.50 Bromoform ND 0.50 Bromomethane ND 0.50 Carbon disulfide ND 0.50 Chlorobenzene ND 0.50 Chlorothane ND 0.50 Chlorothane ND 0.50 Chlorothane ND 0.50 Chlorothane ND 0.50 Chlorothane ND 0.50 Chlorothane ND 0.50 Chlorothane ND 0.50 Carbon tetrachloride ND 0.50 1,1-Dichloroethane ND 0.50 1,2-Dichloroethylene ND 0.50 1,2-Dichloropropane ND 0.50 Dibromochloromethane ND 0.50 trans-1,2-Dichloroethylene ND 0.50 trans-1,3-Dichloropropene ND 0.50 Ethylbenzene ND 0.50 Freon 113 ND 1.0 2-Hexanone ND 0.50 Methylene chloride ND 0.50 ND 0.50 Styrene ND 0.50	Acetone 2-Butanone ND 5.0 0.57 Benzene ND 0.50 0.057 Bromodichloromethane ND 0.50 0.082 Bromoform ND 0.50 0.046 Bromomethane ND 0.50 0.046 Bromomethane ND 0.50 0.046 Bromomethane ND 0.50 0.077 Carbon disulfide ND 0.50 0.028 Chlorobenzene ND 0.50 0.027 Chloroethane ND 0.50 0.037 Chloroform 0.17 0.50 0.031 Chloromethane ND 0.50 0.031 Chloromethane ND 0.50 0.044 Carbon tetrachloride ND 0.50 0.074 1,1-Dichloroethane ND 0.50 0.039 1,1-Dichloroethylene ND 0.50 0.034 1,2-Dichloroethane ND 0.50 0.034 1,2-Dichloromethane ND 0.50 0.039 cis-1,2-Dichloroethylene ND 0.50 0.039 cis-1,2-Dichloroethylene ND 0.50 0.039 cis-1,3-Dichloropropene ND 0.50 0.033 Freon 113 ND 1.0 0.10 2-Hexanone ND 0.50 0.028 ND 0.50 0.033 Freon 113 ND 1.0 0.10 2-Hexanone ND 0.50 0.028 1,1,1-Trichloroethane ND 0.50 0.028 1,1,1-Trichloroethane ND 0.50 0.035 1,1,2-Tetrachloroethane ND 0.50 0.035 1,1,2-Tetrachloroethane ND 0.50 0.035 1,1,2-Trichloroethane ND 0.50 0.035 Tetrachloroethylene ND 0.50 0.035 Tetrachloroethylene ND 0.50 0.035 1,1,2-Trichloroethane ND 0.50 0.035 Tetrachloroethylene ND 0.50 0.035 Tetrachloroethylene ND 0.50 0.035 Tetrachloroethylene ND 0.50 0.035 Tetrachloroethylene ND 0.50 0.055 0.055 Tetrachloroethylene ND 0.50 0.055 0.055	Acetone ND 5.0 0.91 ug/l 2-Butanone ND 5.0 0.57 ug/l Benzene ND 0.50 0.057 ug/l Bromodichloromethane ND 0.50 0.082 ug/l Bromoform ND 0.50 0.046 ug/l Bromoform ND 0.50 0.046 ug/l Bromomethane ND 0.50 0.077 ug/l Carbon disulfide ND 0.50 0.028 ug/l Chlorobenzene ND 0.50 0.028 ug/l Chlorobenzene ND 0.50 0.027 ug/l Chlorothane ND 0.50 0.037 ug/l Chlorothane ND 0.50 0.037 ug/l Chlorothane ND 0.50 0.031 ug/l Chloromethane ND 0.50 0.031 ug/l Chloromethane ND 0.50 0.031 ug/l Chloromethane ND 0.50 0.044 ug/l 1,1-Dichlorothane ND 0.50 0.044 ug/l 1,1-Dichlorothane ND 0.50 0.039 ug/l 1,1-Dichlorothane ND 0.50 0.034 ug/l 1,2-Dichlorothane ND 0.50 0.034 ug/l 1,2-Dichlorothane ND 0.50 0.082 ug/l 1,2-Dichloropropane ND 0.50 0.042 ug/l trans-1,2-Dichlorothylene ND 0.50 0.042 ug/l trans-1,2-Dichlorothylene ND 0.50 0.039 ug/l cis-1,2-Dichlorothylene ND 0.50 0.031 ug/l trans-1,3-Dichloropropene ND 0.50 0.033 ug/l trans-1,3-Dichloropropene ND 0.50 0.033 ug/l Ethylbenzene ND 0.50 0.033 ug/l Ethylbenzene ND 0.50 0.033 ug/l Freon 113 ND 1.0 0.10 ug/l 2-Hexanone ND 2.0 0.084 ug/l ND 0.50 0.035 ug/l Styrene ND 0.50 0.028 ug/l 1,1,2-Trichloroethane ND 0.50 0.035 ug/l 1,1,2-Trichloroethane ND 0.50 0.035 ug/l 1,1,2-Trichloroethylene ND 0.50 0.052 ug/l Tetrachloroethylene ND 0.50 0.052 ug/l Tetrachloroethylene ND 0.50 0.052 ug/l 1,1,2-Trichloroethylene ND 0.50 0.052 ug/l Tetrachloroethylene ND 0.50 0.051 ug/l

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

N = Indicates presumptive evidence of a compound



Report of Analysis

Client Sample ID: TB062315KV1

Lab Sample ID: JB97745-3 **Date Sampled:** 06/23/15 Matrix: AQ - Trip Blank Water Date Received: 06/24/15 Method: EPA 524.2 REV 4.1 **Percent Solids:** n/a

Project: Northrop Grumman, OU2 Hydro, Bethpage, NY

Special VOA List

CAS No.	Compound	Result	RL	MDL	Units	Q	
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l		
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l		
	m,p-Xylene	ND	0.50	0.13	ug/l		
95-47-6	o-Xylene	ND	0.50	0.029	ug/l		
CAS No.	Surrogate Recoveries Run# 1		Run# 2	Limi	its		
2199-69-1	1,2-Dichlorobenzene-d4	95%		78-1	14%		
460-00-4	4-Bromofluorobenzene	91%		77-1	15%		
CAS No.	Tentatively Identified Compe	R.T.	Est.	Conc.	Units	Q	
	Total TIC, Volatile			0		ug/l	

ND = Not detected MDL = Method Detection Limit J = Indicates an estimated value

RL = Reporting Limit

E = Indicates value exceeds calibration range

B = Indicates analyte found in associated method blank N = Indicates presumptive evidence of a compound



Report of Analysis

Client Sample ID: BPOW 6-R Lab Sample ID: JB97745-4

 Lab Sample ID:
 JB97745-4
 Date Sampled:
 06/23/15

 Matrix:
 AQ - Ground Water
 Date Received:
 06/24/15

 Method:
 EPA 524.2 REV 4.1
 Percent Solids:
 n/a

Project: Northrop Grumman, OU2 Hydro, Bethpage, NY

File IDDFAnalyzedByPrep DatePrep BatchAnalytical BatchRun #11B97957.D106/29/15MDn/an/aV1B4647

Run #2

Purge Volume

Run #1 5.0 ml

Run #2

Special VOA List

CAS No.	Compound	Result	RL	MDL	Units	Q
	5.0					
67-64-1	Acetone 5.0	2.9	5.0	0.91	ug/l	→ UB
78-93-3	2-Butanone	ND	5.0	0.57	ug/l	
71-43-2	Benzene	ND	0.50	0.057	ug/l	
75-27-4	Bromodichloromethane	ND	0.50	0.082	ug/l	
75-25-2	Bromoform	ND	0.50	0.046	ug/l	
74-83-9	Bromomethane	ND	0.50	0.077	ug/l	
75-15-0	Carbon disulfide	ND	0.50	0.028	ug/l	
108-90-7	Chlorobenzene	ND	0.50	0.027	ug/l	
75-00-3	Chloroethane	ND	0.50	0.037	ug/l	
67-66-3	Chloroform	ND	0.50	0.031	ug/l	
74-87-3	Chloromethane	ND	0.50	0.044	ug/l	
56-23-5	Carbon tetrachloride	ND	0.50	0.074	ug/l	
75-34-3	1,1-Dichloroethane	ND	0.50	0.039	ug/l	
75-35-4	1,1-Dichloroethylene	ND	0.50	0.054	ug/l	
107-06-2	1,2-Dichloroethane	ND	0.50	0.034	ug/l	
78-87-5	1,2-Dichloropropane	ND	0.50	0.082	ug/l	
124-48-1	Dibromochloromethane	ND	0.50	0.042	ug/l	
156-60-5	trans-1,2-Dichloroethylene	ND	0.50	0.039	ug/l	
156-59-2	cis-1,2-Dichloroethylene	ND	0.50	0.081	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.033	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.063	ug/l	
100-41-4	Ethylbenzene	ND	0.50	0.033	ug/l	
76-13-1	Freon 113	ND	1.0	0.10	ug/l	
591-78-6	2-Hexanone	ND	2.0	0.084	ug/l	
75-09-2	Methylene chloride	ND	0.50	0.047	ug/l	
108-10-1	4-Methyl-2-pentanone	ND	2.0	0.27	ug/l	
100-42-5	Styrene	ND	0.50	0.028	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	0.50	0.050	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.035	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	0.50	0.052	ug/l	
127-18-4	Tetrachloroethylene	ND	0.50	0.091	ug/l	
108-88-3	Toluene	ND	0.50	0.044	ug/l	

ND = Not detected

MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



06/24/15

n/a

Report of Analysis

Client Sample ID: BPOW 6-R Lab Sample ID: JB97745-4

Date Sampled: 06/23/15 Matrix: AQ - Ground Water **Date Received:** Method: **Percent Solids:** EPA 524.2 REV 4.1

Project: Northrop Grumman, OU2 Hydro, Bethpage, NY

Special VOA List

CAS No.	Compound	Result	RL	MDL	Units	Q		
79-01-6 75-01-4 95-47-6	Trichloroethylene Vinyl chloride m,p-Xylene o-Xylene	ND ND ND ND	0.50 0.50 0.50 0.50	0.024 0.032 0.13 0.029	ug/l ug/l ug/l ug/l			
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limi	ts			
2199-69-1 460-00-4	1,2-Dichlorobenzene-d4 4-Bromofluorobenzene	99% 91%		78-11 77-11				
CAS No.	Tentatively Identified Compo	ounds	R.T.	Est.	Conc.	Units	Q	
67-63-0 78-84-2 15045-43-9	Isopropyl Alcohol Propanal, 2-methyl- Furan, tetrahydro-2,2,5,5-tetra Total TIC, Volatile		7.47 8.72 12.88 3.39	3.5 2.7 .69 6.89	-	ug/l ug/l ug/l ug/l	JN JN JN J N	R

ND = Not detected MDL = Method Detection Limit

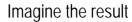
RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank N = Indicates presumptive evidence of a compound







Northrop Grumman Corporation-Operable Unit 2

Data Review

BETHPAGE, NEW YORK

Volatile Analysis

SDG #JB97890

Analyses Performed By: Accutest Laboratories Dayton, New Jersey

Report #23914R July 10, 2015 Review Level: Tier II Project #NY001496.1514.NAVI4

SUMMARY

This data quality assessment summarizes the review of Sample Delivery Group (SDG) #JB97890 for samples collected in association with the Northrop Grumman-Bethpage Site. The review was conducted as a Tier II evaluation and included review of data package completeness. Only analytical data associated with constituents of concern were reviewed for this validation. Field documentation was not included in this review. Included with this assessment are the validation annotated sample result sheets, and chain of custody. Analyses were performed on the following samples:

	Sample							,	Analysis	6	
Sample ID	Lab ID	Matrix	Collection Date	Parent Sample	voc	svoc	РСВ	MET	MISC		
BPOW 6-3	JB97890-1	Water	06/24/2015		Х						
TB062415KV1	JB97890-2	Water	06/24/2015		Х						
FB062415KV1	JB97890-3	Water	06/24/2015		Х						
BPOW 6-4	JB97890-4	Water	06/24/2015		Х				·		

Note:

1. The matrix spike/matrix spike duplicate (MS/MSD) analysis was performed on sample location BPOW 6-3.

ANALYTICAL DATA PACKAGE DOCUMENTATION

GENERAL INFORMATION

Items Reviewed	Rep	orted		mance ptable	Not
	No	Yes	No	Yes	Required
1. Sample receipt condition		Х		Х	
2. Requested analyses and sample results		Х		Х	
3. Collection Technique (grab, composite, etc.)		Х		Х	
4. Methods of analysis		Х		Х	
5. Reporting limits		Х		Х	
6. Sample collection date		Х		Х	
7. Laboratory sample received date		Х	Х		
8. Sample preservation verification (as applicable)		Х		X	
9. Sample preparation/extraction/analysis dates		Х		Х	
10. Fully executed Chain-of-Custody (COC) form completed		Х		Х	
11. Narrative summary of QA or sample problems provided		Х		Х	
12. Data Package Completeness and Compliance		Х		Х	

QA - Quality Assurance

VOLATILE ANALYSIS INTRODUCTION

Analyses were performed according to United States Environmental Protection Agency (USEPA), Method 524.2. Data were reviewed in accordance with USEPA National Functional Guidelines of October 1999.

The data review process is an evaluation of data on a technical basis rather than a determination of contract compliance. As such, the standards against which the data are being weighed may differ from those specified in the analytical method. It is assumed that the data package represents the best efforts of the laboratory and had already been subjected to adequate and sufficient quality review prior to submission.

During the review process, laboratory qualified and unqualified data are verified against the supporting documentation. Based on this evaluation, qualifier codes may be added, deleted, or modified by the data reviewer. Results are qualified with the following codes in accordance with USEPA National Functional Guidelines:

- Concentration (C) Qualifiers
 - U The compound was analyzed for but not detected. The associated value is the compound quantitation limit.
 - B The compound has been found in the sample as well as its associated blank, its presence in the sample may be suspect.
- Quantitation (Q) Qualifiers
 - E The compound was quantitated above the calibration range.
 - D Concentration is based on a diluted sample analysis.
- Validation Qualifiers
 - J The compound was positively identified; however, the associated numerical value is an estimated concentration only.
 - UJ The compound was not detected above the reported sample quantitation limit. However, the reported limit is approximate and may or may not represent the actual limit of quantitation.
 - JN The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification. The associated numerical value is an estimated concentration only.
 - UB Compound considered non-detect at the listed value due to associated blank contamination.
 - N The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification.
 - R The sample results are rejected.

Two facts should be noted by all data users. First, the "R" flag means that the associated value is unusable. In other words, due to significant quality control (QC) problems, the analysis is invalid and provides no information as to whether the compound is present or not. "R" values should not appear on data tables because they cannot be relied upon, even as a last resort. The second fact to keep in mind is that no compound concentration, even if it has passed all QC tests, is guaranteed to be accurate. Strict QC serves to increase confidence in data but any value potentially contains error.

VOLATILE ORGANIC COMPOUND (VOC) ANALYSES

1. Holding Times

The specified holding times for the following methods are presented in the following table.

Method	Matrix	Holding Time	Preservation
EPA 524.2	Water	14 days from collection to analysis	Cool to < 6°C; preserved to a pH of less than 2 s.u

s.u. Standard units

All samples were analyzed within the specified holding time and temperature criteria.

2. Blank Contamination

Quality assurance (QA) blanks (i.e., method and rinse blanks) are prepared to identify any contamination which may have been introduced into the samples during sample preparation or field activity. Method blanks measure laboratory contamination. Rinse blanks measure contamination of samples during field operations.

A blank action level (BAL) of five times the concentration of a detected compound in an associated blank (common laboratory contaminant compounds are calculated at ten times) is calculated for QA blanks containing concentrations greater than the method detection limit (MDL). The BAL is compared to the associated sample results to determine the appropriate qualification of the sample results, if needed.

All compounds associated with the QA blanks exhibited a concentration less than the MDL, with the exception of the compounds listed in the following table. Sample results associated with QA blank contamination that were greater than the BAL resulted in the removal of the laboratory qualifier (B) of data. Sample results less than the BAL associated with the following sample locations were qualified as listed in the following table.

Sample Locations	Analytes	Sample Result	Qualification
BPOW 6-3	Acetone	Detected sample results >RL and <bal< td=""><td>"UB" at detected sample concentration</td></bal<>	"UB" at detected sample concentration
	TICs: Isopropyl alcohol (RT: 7.46)	Detected sample results less than 5 times blank result	R

RL Reporting limit

3. Surrogates/System Monitoring Compounds

All samples to be analyzed for organic compounds are spiked with surrogate compounds prior to sample preparation to evaluate overall laboratory performance and efficiency of the analytical technique. VOC analysis requires that all surrogates associated with the analysis exhibit recoveries within the laboratory-established acceptance limits.

All surrogate recoveries were within control limits.

4. Matrix Spike/Matrix Spike Duplicate (MS/MSD) Analysis

MS/MSD data are used to assess the precision and accuracy of the analytical method. The compounds used to perform the MS/MSD analysis must exhibit a percent recovery within the laboratory-established acceptance limits. The relative percent difference (RPD) between the MS/MSD recoveries must exhibit an RPD within the laboratory-established acceptance limits.

Note: The MS/MSD recovery control limits do not apply for MS/MSD performed on sample locations where the compound concentration detected in the parent sample exceeds the MS/MSD concentration by a factor of four or greater.

The MS/MSD exhibited acceptable recoveries and RPD between the MS/MSD recoveries.

5. Laboratory Control Sample (LCS) Analysis

The LCS analysis is used to assess the precision and accuracy of the analytical method independent of matrix interferences. The compounds associated with the LCS analysis must exhibit a percent recovery within the laboratory-established acceptance limits.

All compounds associated with the LCS analysis exhibited recoveries within the control limits.

6. Field Duplicate Analysis

Field duplicate analysis is used to assess the overall precision of the field sampling procedures and analytical method. A control limit of 30% for water matrices is applied to the RPD between the parent sample and the field duplicate. In the instance when the parent and/or duplicate sample concentrations are less than or equal to 5 times the RL, a control limit of two times the RL is applied for water matrices.

A field duplicate was not collected on a sample location associated with this SDG.

7. Laboratory Duplicate Analysis

The laboratory duplicate relative percent difference (RPD) criterion is applied when parent and duplicate sample concentrations are greater than or equal to 5 times the RL. A control limit of 20% for water matrices is applied when the criteria above is true. In the instance when the parent and/or duplicate sample concentrations are less than or equal to 5 times the RL, a control limit of two times the RL is applied for water matrices.

A laboratory duplicate was not performed on a sample location associated with this SDG.

8. System Performance and Overall Assessment

Tentatively identified compounds (TICs) were identified in sample locations BPOW 6-3, TB062415KV1 and FB062415KV1. The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification. TICs are qualified as estimated (JN).

Overall system performance was noted by the laboratory in the case narrative as acceptable. Other than for those deviations specifically mentioned in this review, the overall data quality is within the guidelines specified in the method.

DATA VALIDATION CHECKLIST FOR VOCs

VOCs: 524.2	Rep	orted		mance ptable	Not Required
	No	Yes	No	Yes	Required
GAS CHROMATOGRAPHY/MASS SPECTROME	TRY (GC/	MS)			
Tier II Validation					
Holding times & Temperature		Х		Х	
Reporting limits (units)		Х		Х	
Blanks					
A. Method blanks		Х		Х	
B. Equipment blanks		Х	Х		
C. Trip blanks		Х	Х		
Surrogate (%R)		Х		Х	
Laboratory Control Sample (%R)		Х		Х	
Laboratory Control Sample Duplicate(LCSD)					Х
LCS/LCSD Precision (RPD)					Х
Matrix Spike (MS)		Х		Х	
Matrix Spike Duplicate(MSD)		Х		Х	
MS/MSD Precision (RPD)		Х		Х	
Field/Lab Duplicate (RPD)					Х
Dilution Factor		Х		Х	
Moisture Content					Х

[%]R Percent Recovery RPD Relative Percent Difference

VALIDATION PERFORMED BY: Lisa Horton

SIGNATURE: XISA 77

DATE: _ July 10, 2015

PEER REVIEW BY: Todd Church

DATE: July 16, 2015

CHAIN OF CUSTODY/ CORRECTED SAMPLE ANALYSIS DATA SHEETS

ACCUTEST.

CHAIN OF CUSTODY
Accutest New Jersey/SPL Environmental
2235 Route 130, Dayton, NJ 08810

PAGE	\perp	OF	1
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JB97890: Chain of Custody

Page 1 of 2



Date Sampled: 06/24/15

Report of Analysis

Client Sample ID: BPOW 6-3 Lab Sample ID: JB97890-1

 Matrix:
 AQ - Water
 Date Received:
 06/25/15

 Method:
 EPA 524.2 REV 4.1
 Percent Solids:
 n/a

Project: Northrop Grumman, OU2 Hydro, Bethpage, NY

File ID DF Analyzed By Prep Date Prep Batch Analytical Batch
Run #1 1B97954.D 1 06/29/15 MD n/a n/a V1B4647

Run #2

Purge Volume

Run #1 5.0 ml

Run #2

Special VOA List

CAS No.	Compound	Result	RL	MDL	Units	Q	
67-64-1	Acetone 5.0	1.0	5.0	0.91	ug/l	J	UB
78-93-3	2-Butanone	ND	5.0	0.57	ug/l	·	OD
71-43-2	Benzene	ND	0.50	0.057	ug/l		
75-27-4	Bromodichloromethane	ND	0.50	0.082	ug/l		
75-25-2	Bromoform	ND	0.50	0.046	ug/l		
74-83-9	Bromomethane	ND	0.50	0.077	ug/l		
75-15-0	Carbon disulfide	ND	0.50	0.028	ug/l		
108-90-7	Chlorobenzene	ND	0.50	0.027	ug/l		
75-00-3	Chloroethane	ND	0.50	0.037	ug/l		
67-66-3	Chloroform	ND	0.50	0.031	ug/l		
74-87-3	Chloromethane	ND	0.50	0.044	ug/l		
56-23-5	Carbon tetrachloride	ND	0.50	0.074	ug/l		
75-34-3	1,1-Dichloroethane	ND	0.50	0.039	ug/l		
75-35-4	1,1-Dichloroethylene	ND	0.50	0.054	ug/l		
107-06-2	1,2-Dichloroethane	ND	0.50	0.034	ug/l		
78-87-5	1,2-Dichloropropane	ND	0.50	0.082	ug/l		
124-48-1	Dibromochloromethane	ND	0.50	0.042	ug/l		
156-60-5	trans-1,2-Dichloroethylene	ND	0.50	0.039	ug/l		
156-59-2	cis-1,2-Dichloroethylene	ND	0.50	0.081	ug/l		
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.033	ug/l		
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.063	ug/l		
100-41-4	Ethylbenzene	ND	0.50	0.033	ug/l		
76-13-1	Freon 113	ND	1.0	0.10	ug/l		
591-78-6	2-Hexanone	ND	2.0	0.084	ug/l		
75-09-2	Methylene chloride	ND	0.50	0.047	ug/l		
108-10-1	4-Methyl-2-pentanone	ND	2.0	0.27	ug/l		
100-42-5	Styrene	ND	0.50	0.028	ug/l		
71-55-6	1,1,1-Trichloroethane	ND	0.50	0.050	ug/l		
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.035	ug/l		
79-00-5	1,1,2-Trichloroethane	ND	0.50	0.052	ug/l		
127-18-4	Tetrachloroethylene	ND	0.50	0.091	ug/l		
108-88-3	Toluene	ND	0.50	0.044	ug/l		

ND = Not detected

MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

N = Indicates presumptive evidence of a compound



Date Sampled: 06/24/15

Report of Analysis

Client Sample ID: BPOW 6-3 Lab Sample ID: JB97890-1 Matrix:

Method: EPA 524.2 REV 4.1

AQ - Water Date Received: 06/25/15 **Percent Solids:** n/a

Project: Northrop Grumman, OU2 Hydro, Bethpage, NY

Special VOA List

CAS No.	Compound	Result	RL	MDL	Units	Q		
79-01-6 75-01-4	Trichloroethylene Vinyl chloride m,p-Xylene	ND ND ND	0.50 0.50 0.50	0.024 0.032 0.13	ug/l ug/l ug/l			
95-47-6	o-Xylene	ND	0.50	0.029	ug/l			
CAS No.	Surrogate Recoveries	ogate Recoveries Run# 1 R		Lim	its			
2199-69-1 460-00-4	1,2-Dichlorobenzene-d4 4-Bromofluorobenzene	96% 91%		78-1 77-1				
CAS No.	Tentatively Identified Compo	ounds	R.T.	Est.	Conc.	Units	Q	
67-63-0	Isopropyl Alcohol Total TIC, Volatile		7.46	3.5 3.5		ug/l ug/l	JN J	R

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank N = Indicates presumptive evidence of a compound



Report of Analysis

Client Sample ID: TB062415AM1

Lab Sample ID:JB97890-2Date Sampled:06/24/15Matrix:AQ - Trip Blank WaterDate Received:06/25/15Method:EPA 524.2 REV 4.1Percent Solids:n/a

Project: Northrop Grumman, OU2 Hydro, Bethpage, NY

File ID DF Analyzed By Prep Date Prep Batch Analytical Batch
Run #1 1B97963.D 1 06/29/15 MD n/a n/a V1B4647

Run #2

Purge Volume

Run #1 5.0 ml

Run #2

Special VOA List

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	1.5	5.0	0.91	ug/l	J
78-93-3	2-Butanone	ND	5.0	0.57	ug/l	
71-43-2	Benzene	ND	0.50	0.057	ug/l	
75-27-4	Bromodichloromethane	ND	0.50	0.082	ug/l	
75-25-2	Bromoform	ND	0.50	0.046	ug/l	
74-83-9	Bromomethane	ND	0.50	0.077	ug/l	
75-15-0	Carbon disulfide	ND	0.50	0.028	ug/l	
108-90-7	Chlorobenzene	ND	0.50	0.027	ug/l	
75-00-3	Chloroethane	ND	0.50	0.037	ug/l	
67-66-3	Chloroform	0.17	0.50	0.031	ug/l	J
74-87-3	Chloromethane	ND	0.50	0.044	ug/l	
56-23-5	Carbon tetrachloride	ND	0.50	0.074	ug/l	
75-34-3	1,1-Dichloroethane	ND	0.50	0.039	ug/l	
75-35-4	1,1-Dichloroethylene	ND	0.50	0.054	ug/l	
107-06-2	1,2-Dichloroethane	ND	0.50	0.034	ug/l	
78-87-5	1,2-Dichloropropane	ND	0.50	0.082	ug/l	
124-48-1	Dibromochloromethane	ND	0.50	0.042	ug/l	
156-60-5	trans-1,2-Dichloroethylene	ND	0.50	0.039	ug/l	
156-59-2	cis-1,2-Dichloroethylene	ND	0.50	0.081	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.033	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.063	ug/l	
100-41-4	Ethylbenzene	ND	0.50	0.033	ug/l	
76-13-1	Freon 113	ND	1.0	0.10	ug/l	
591-78-6	2-Hexanone	ND	2.0	0.084	ug/l	
75-09-2	Methylene chloride	ND	0.50	0.047	ug/l	
108-10-1	4-Methyl-2-pentanone	ND	2.0	0.27	ug/l	
100-42-5	Styrene	ND	0.50	0.028	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	0.50	0.050	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.035	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	0.50	0.052	ug/l	
127-18-4	Tetrachloroethylene	ND	0.50	0.091	ug/l	
108-88-3	Toluene	ND	0.50	0.044	ug/l	

ND = Not detected

MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



Page 2 of 2

4

Report of Analysis

Client Sample ID: TB062415AM1

 Lab Sample ID:
 JB97890-2
 Date Sampled:
 06/24/15

 Matrix:
 AQ - Trip Blank Water
 Date Received:
 06/25/15

 Method:
 EPA 524.2 REV 4.1
 Percent Solids:
 n/a

Project: Northrop Grumman, OU2 Hydro, Bethpage, NY

Special VOA List

CAS No.	Compound	Result	RL	MDL	Units	Q	
79-01-6 75-01-4 95-47-6	Trichloroethylene Vinyl chloride m,p-Xylene o-Xylene	ND ND ND ND	0.50 0.50 0.50 0.50	0.024 0.032 0.13 0.029	ug/l ug/l ug/l ug/l		
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limi	ts		
2199-69-1 460-00-4	1,2-Dichlorobenzene-d4 4-Bromofluorobenzene	94% 86%		78-11 77-11			
CAS No.	Tentatively Identified Compo	ounds	R.T.	Est.	Conc.	Units	Q
67-63-0	Isopropyl Alcohol unknown Total TIC, Volatile		7.47 9.67	3.4 .7 4.1		ug/l ug/l ug/l	JN J _N J _N

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank N = Indicates presumptive evidence of a compound



Report of Analysis

Client Sample ID: FB062415AM1

Lab Sample ID:JB97890-3Date Sampled:06/24/15Matrix:AQ - Field Blank WaterDate Received:06/25/15Method:EPA 524.2 REV 4.1Percent Solids:n/a

Project: Northrop Grumman, OU2 Hydro, Bethpage, NY

File ID DF **Prep Date Analytical Batch** Analyzed By **Prep Batch** Run #1 1B97961.D 1 06/29/15 MD V1B4647 n/an/aRun #2

Purge Volume Run #1 5.0 ml

Run #2

Special VOA List

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	1.3	5.0	0.91	ug/l	J
78-93-3	2-Butanone	ND	5.0	0.57	ug/l	
71-43-2	Benzene	ND	0.50	0.057	ug/l	
75-27-4	Bromodichloromethane	ND	0.50	0.082	ug/l	
75-25-2	Bromoform	ND	0.50	0.046	ug/l	
74-83-9	Bromomethane	ND	0.50	0.077	ug/l	
75-15-0	Carbon disulfide	ND	0.50	0.028	ug/l	
108-90-7	Chlorobenzene	ND	0.50	0.027	ug/l	
75-00-3	Chloroethane	ND	0.50	0.037	ug/l	
67-66-3	Chloroform	0.099	0.50	0.031	ug/l	J
74-87-3	Chloromethane	ND	0.50	0.044	ug/l	
56-23-5	Carbon tetrachloride	ND	0.50	0.074	ug/l	
75-34-3	1,1-Dichloroethane	ND	0.50	0.039	ug/l	
75-35-4	1,1-Dichloroethylene	ND	0.50	0.054	ug/l	
107-06-2	1,2-Dichloroethane	ND	0.50	0.034	ug/l	
78-87-5	1,2-Dichloropropane	ND	0.50	0.082	ug/l	
124-48-1	Dibromochloromethane	ND	0.50	0.042	ug/l	
156-60-5	trans-1,2-Dichloroethylene	ND	0.50	0.039	ug/l	
156-59-2	cis-1,2-Dichloroethylene	ND	0.50	0.081	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.033	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.063	ug/l	
100-41-4	Ethylbenzene	ND	0.50	0.033	ug/l	
76-13-1	Freon 113	ND	1.0	0.10	ug/l	
591-78-6	2-Hexanone	ND	2.0	0.084	ug/l	
75-09-2	Methylene chloride	ND	0.50	0.047	ug/l	
108-10-1	4-Methyl-2-pentanone	ND	2.0	0.27	ug/l	
100-42-5	Styrene	ND	0.50	0.028	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	0.50	0.050	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.035	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	0.50	0.052	ug/l	
127-18-4	Tetrachloroethylene	ND	0.50	0.091	ug/l	
108-88-3	Toluene	ND	0.50	0.044	ug/l	

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank N = Indicates presumptive evidence of a compound

N = Indicates presumptive evidence of a compound



Page 2 of 2

Report of Analysis

Client Sample ID: FB062415AM1

Lab Sample ID:JB97890-3Date Sampled:06/24/15Matrix:AQ - Field Blank WaterDate Received:06/25/15Method:EPA 524.2 REV 4.1Percent Solids:n/a

Project: Northrop Grumman, OU2 Hydro, Bethpage, NY

Special VOA List

CAS No.	Compound	Result	RL	MDL	Units	Q	
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l		
75-01-4	Vinyl chloride m,p-Xylene	ND ND	0.50 0.50	0.032 0.13	ug/l ug/l		
95-47-6	o-Xylene	ND	0.50	0.029	ug/l		
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limi	its		
2199-69-1	1,2-Dichlorobenzene-d4	99%		78-1			
460-00-4	4-Bromofluorobenzene	91%		77-1	15%		
CAS No.	Tentatively Identified Compo	R.T.	Est.	Conc.	Units	Q	
67-63-0	Isopropyl Alcohol Total TIC, Volatile		7.48	1.5 1.5		ug/l ug/l	JN J <mark>N</mark>

ND = Not detected MDL = Method Detection Limit J = Indicates and J = Indicates

RL = Reporting Limit

E = Indicates value exceeds calibration range

 $J = \ Indicates \ an \ estimated \ value$

B = Indicates analyte found in associated method blank N = Indicates presumptive evidence of a compound



Report of Analysis

 Client Sample ID:
 BPOW 6-4

 Lab Sample ID:
 JB97890-4
 Date Sampled:
 06/24/15

 Matrix:
 AQ - Water
 Date Received:
 06/25/15

 Method:
 EPA 524.2 REV 4.1
 Percent Solids:
 n/a

Project: Northrop Grumman, OU2 Hydro, Bethpage, NY

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B97962.D	1	06/29/15	MD	n/a	n/a	V1B4647
Run #2							

	Purge Volume		
Run #1	5.0 ml		
Run #2			

Special VOA List

CAS No. Compound		Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	5.0	0.91	ug/l	
78-93-3	2-Butanone	ND	5.0	0.57	ug/l	
71-43-2	Benzene	ND	0.50	0.057	ug/l	
75-27-4	Bromodichloromethane	ND	0.50	0.082	ug/l	
75-25-2	Bromoform	ND	0.50	0.046	ug/l	
74-83-9	Bromomethane	ND	0.50	0.077	ug/l	
75-15-0	Carbon disulfide	ND	0.50	0.028	ug/l	
108-90-7	Chlorobenzene	ND	0.50	0.027	ug/l	
75-00-3	Chloroethane	ND	0.50	0.037	ug/l	
67-66-3	Chloroform	ND	0.50	0.031	ug/l	
74-87-3	Chloromethane	ND	0.50	0.044	ug/l	
56-23-5	Carbon tetrachloride	ND	0.50	0.074	ug/l	
75-34-3	1,1-Dichloroethane	ND	0.50	0.039	ug/l	
75-35-4	1,1-Dichloroethylene	ND	0.50	0.054	ug/l	
107-06-2	1,2-Dichloroethane	ND	0.50	0.034	ug/l	
78-87-5	1,2-Dichloropropane	ND	0.50	0.082	ug/l	
124-48-1	Dibromochloromethane	ND	0.50	0.042	ug/l	
156-60-5	trans-1,2-Dichloroethylene	ND	0.50	0.039	ug/l	
156-59-2	cis-1,2-Dichloroethylene	ND	0.50	0.081	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.033	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.063	ug/l	
100-41-4	Ethylbenzene	ND	0.50	0.033	ug/l	
76-13-1	Freon 113	ND	1.0	0.10	ug/l	
591-78-6	2-Hexanone	ND	2.0	0.084	ug/l	
75-09-2	Methylene chloride	ND	0.50	0.047	ug/l	
108-10-1	4-Methyl-2-pentanone	ND	2.0	0.27	ug/l	
100-42-5	Styrene	ND	0.50	0.028	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	0.50	0.050	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.035	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	0.50	0.052	ug/l	
127-18-4	Tetrachloroethylene	ND	0.50	0.091	ug/l	
108-88-3	Toluene	ND	0.50	0.044	ug/l	

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank N = Indicates presumptive evidence of a compound



Page 2 of 2

Date Sampled: 06/24/15

06/25/15

n/a

Date Received:

Percent Solids:

Report of Analysis

Client Sample ID: BPOW 6-4 Lab Sample ID: JB97890-4 Matrix: AQ - Water

Method: EPA 524.2 REV 4.1

Northrop Grumman, OU2 Hydro, Bethpage, NY

Project:

Special VOA List

CAS No. Compound R		Result	RL	MDL	Units	Q	
79-01-6 75-01-4	Trichloroethylene Vinyl chloride m,p-Xylene	ND ND ND	0.50 0.50 0.50	0.024 0.032 0.13	ug/l ug/l ug/l		
95-47-6 CAS No.	o-Xylene Surrogate Recoveries	ND Run# 1	0.50 0.029 ug/l		C		
2199-69-1	1,2-Dichlorobenzene-d4			78-1			
460-00-4 CAS No.	4-Bromofluorobenzene 91% Tentatively Identified Compounds		R.T.	77-1 Est.	15% Conc.	Units	0
0120 1101	Total TIC, Volatile			0		ug/l	· ·

ND = Not detected MDL = Method Detection Limit J = Indicates an estimated value

RL = Reporting Limit

E = Indicates value exceeds calibration range

B = Indicates analyte found in associated method blank N = Indicates presumptive evidence of a compound





Table 1. Concentrations of Volatile Organic Compounds in Monitoring Wells BPOW 6-1, BPOW 6-2, BPOW 6-3 and BPOW 6-4, Second Quarter 2015, Operable Unit 2 (Groundwater), Bethpage, New York.

CONSTITUENT (Units in µg/L)	Well: Sample ID: Date:	BPOW 6-1 BPOW 6-1 6/22/2015	BPOW 6-2 BPOW 6-2 6/23/2015	BPOW 6-2 BPOW 6-R ⁽¹⁾ 6/23/2015	BPOW 6-3 BPOW 6-3 6/24/2015	BPOW 6-4 BPOW 6-4 6/24/2015
1,1,1-Trichloroethane		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
1,1,2,2-Tetrachloroethane		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
1,1,2-trichloro-1,2,2-trifluroethane		< 1.0	< 1.0	< 1.0	< 1.0	< 1.0
1,1,2-Trichloroethane		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
1,1-Dichloroethane		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
1,1-Dichloroethene		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
1,2-Dichloroethane		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
1,2-Dichloropropane		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
2-Butanone (MEK)		< 5.0	< 5.0	< 5.0	< 5.0	< 5.0
2-Hexanone		< 2.0	< 2.0	< 2.0	< 2.0	< 2.0
4-methyl-2-pentanone (MIK)		< 2.0	< 2.0	< 2.0	< 2.0	< 2.0
Acetone		< 9.3 B	< 5.0 B	< 5.0 B	< 5.0 B	< 5.0
Benzene		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Bromodichloromethane		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Bromoform		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Bromomethane		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Carbon Disulfide		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Carbon tetrachloride		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Chlorobenzene		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Chloroethane		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Chloroform		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Chloromethane		0.51	< 0.50	< 0.50	< 0.50	< 0.50
cis-1,2-dichloroethene		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
cis-1,3-dichloropropene		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Dibromochloromethane		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Ethylbenzene		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Methylene Chloride		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Styrene		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Tetrachloroethene		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Toluene		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
trans-1,2-dichloroethene		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
trans-1,3-dichloropropene		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Trichloroethylene		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Vinyl Chloride		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Xylene-o		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Xylenes - m,p		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Total VOCs		0.51	0	0	0	0

Notes and Abbreviations:

(1) BPOW 6-R is a blind duplicate sample.

Results validated following protocols specified in OU2 Groundwater Monitoring Plan (ARCADIS 2014)

Samples analyzed for the TCL VOCs using USEPA Method 524.2.

Total VOCs are rounded to two significant figures.

Bold value indicates a detection

TCL Target Compound List VOC Volatile Organic Compound

USEPA United States Environmental Protection Agency

μg/L Micrograms per liter

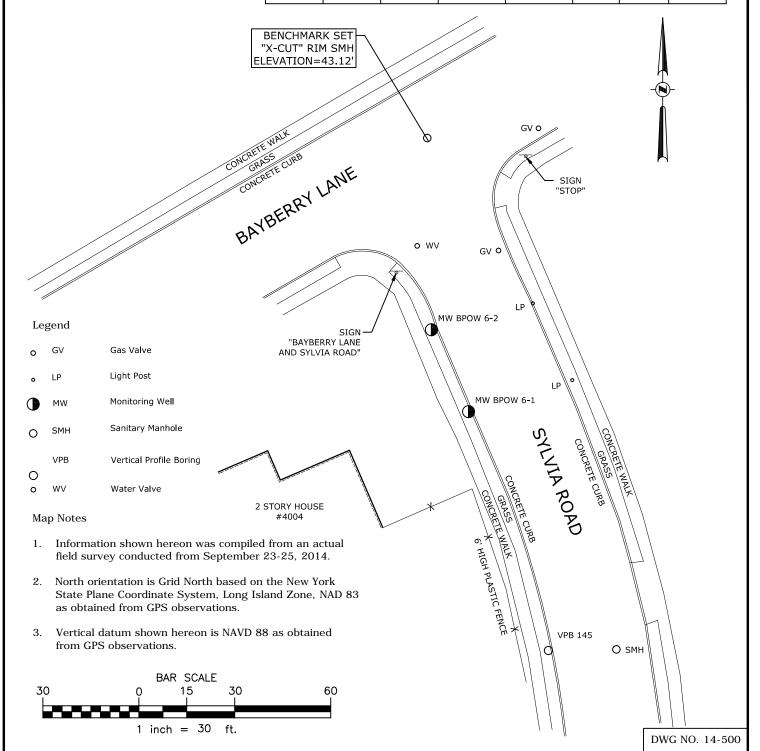
B Compound detected in associated blank sample

Section 7

Survey



Description	Northing	Easting	Latitude	Longitude	Ground	Rim	PVC
VPB 145	193153.53	1126821.88	N40-41-44.48	W73-29-09.07	44.20	NA	NA
MWBPOW6-1	193228.16	1126796.92	N40-41-45.22	W73-29-09.38	43.61	43.61	42.93
MWBPOW6-2	193253.75	1126785.34	N40-41-45.48	W73-29-09.53	43.58	43.58	43.08



~	Date	RECORD OF	WORK	Appr.		
1	7-15-15	WELL DESIGNATIONS REVISED		JFC		
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	Drafter:	Orafter: LMK Checker: JFC				
	Appr. by	7: JFC	Proj. No. 14.4121		Ş	

VERTICAL PROFILE BORING 145 SURVEY LOCATION 4004 BAYBERRY LANE

TOWN OF SEAFORD NASSAU COUNTY, NEW YORK

C.T. MALE ASSOCIATES

Engineering, Surveying, Architecture & Landscape Architecture, D.P.C.

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

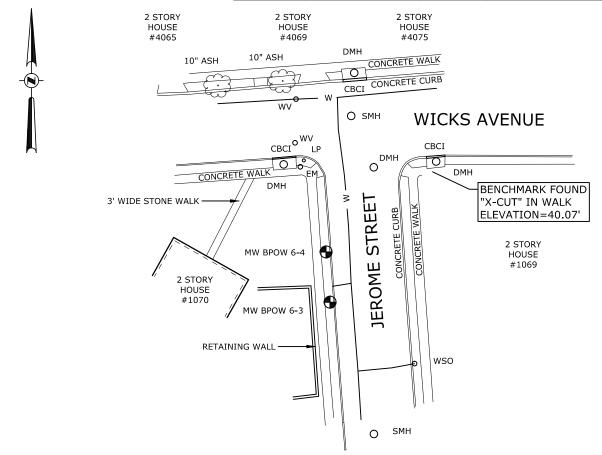


SCALE: 1"=30' DATE: SEPT. 24, 2014

Appr. by: JFC

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

Description	Northing	Easting	Latitude	Longitude	Ground	Rim	PVC
MW BPOW 6-4	194127.00	1127580.66	N40-41-54.06	W73-28-59.14	40.40	40.40	40.02
MW BPOW 6-3	194106.07	1127582.36	N40-41-53.85	W73-28-59.12	40.34	40.34	39.96



Legend

CBCI

Deciduous Tree

o EM
Electric Marker

o LP
Light Pole
MW BPOW 6-3
Monitoring Well

O SMH
Sanitary Manhole

o WSO
Water Shut Off

o WV
Water Valve

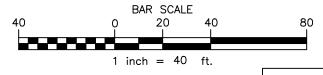
Catch Basin Curb Inlet

Water Line

14.4121

Map Notes

- Information shown hereon was compiled from an actual field survey conducted on March 26, 2015.
- North orientation is Grid North based on the New York State Plane Coordinate System, Long Island Zone, NAD 83 as obtained from GPS observations.
- 3. Vertical datum shown hereon is NAVD 88 as obtained from GPS observations.



DWG NO. 15-219

Date	RECORD OF WORK			Appr.		
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Drafter: LMK Checker: JFC]	

Proj. No.

VERTICAL PROFILE BORING 146 SURVEY LOCATION MW BPOW 6-3 AND MW BPOW 6-4 4099 JEROME STREET

TOWN OF BETHPAGE

SCALE: 1"=40'

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



DATE: MARCH 26, 2015

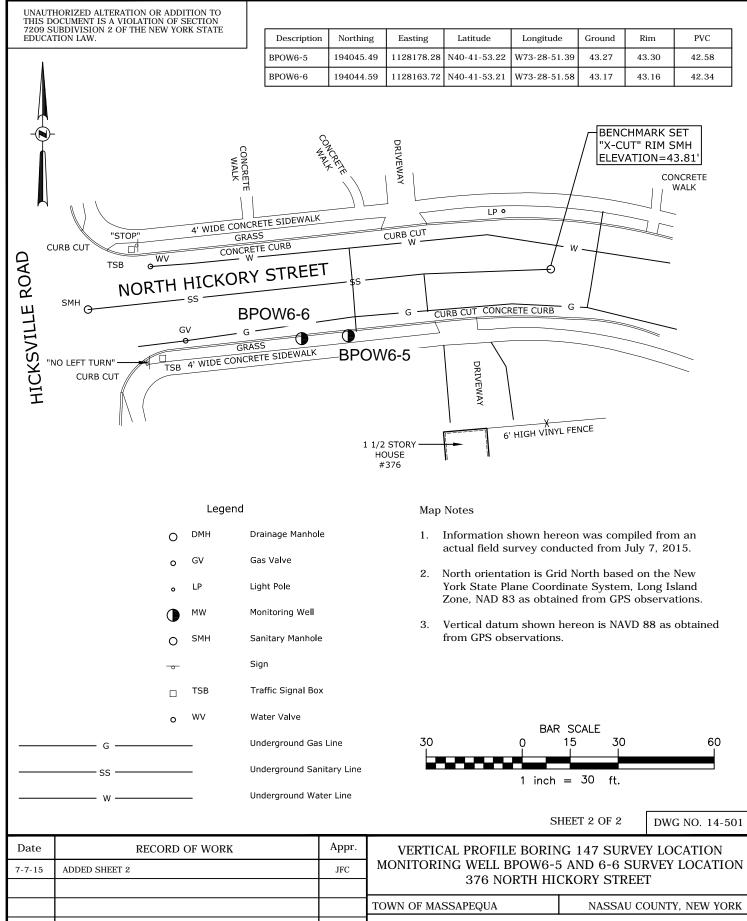
Drafter: LMK

Appr. by: JFC

Checker: JFC

Proj. No.

14.4121



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

50 CENTURY HILL DRIVE LATHAM NY 12110

SCALE: 1"=30'

518.786.7400 * FAX 518.786.7299

DATE: SEPT. 24, 2014