

**2014 OU2 GROUNDWATER INVESTIGATION  
VPB 152  
BETHPAGE, NY**

Prepared for:



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

**Comprehensive Long-Term Environmental Action Navy  
Contract Number N62470-11-D-8013**

**CTO WE15**

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## List of Acronyms and Abbreviations

AOC	Area of Concern
bgs	below ground surface
DoD	Department of Defense
ELAP	Environmental Laboratory Accreditation Program
EPA	Environmental Protection Agency, United States
ft	feet
GOCO	Government-Owned Contractor-Operated
IDW	Investigation Derived Waste
IR	Installation Restoration
Katahdin	Katahdin Analytical Services, Inc
NAD	North American Datum
NAVD	North American Vertical Datum
NAVFAC	Naval Facilities Engineering Command
NG	Northrop Grumman
NWIRP	Naval Weapons Industrial Reserve Plant
NYSDEC	New York State Department of Environmental Conservation
OU	Operable Unit
PCBs	Polychlorinated Biphenyls
PCE	Tetrachloroethene
PID	Photoionization Detector
POTW	Publicly Owned Treatment Works
PPE	Personal Protective Equipment
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
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, Mid-Atlantic under contract task order WE15 Contract N62470-11-D-8013. This report describes vertical profile boring (VPB) installation activities (specifically at the VPB 152 location) in 2014 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 data summary report provides information on the installation of VPB 152. The purpose of the VPB 152 investigation was to ascertain subsurface conditions and contaminant levels upgradient of the South Farmington Water District Plant 6 wellfield and to design outpost wells that will be used to provide early warning of plume migration toward the South Farmington Water District Plant 6 wellfield. VPB locations within the general vicinity of VPB 152 are shown in Figure 2. VPB 152 was completed to 997 feet (ft) below ground surface (bgs).

Field tasks were conducted in 2014 in accordance with the *United Federal Programs Sampling and Analysis Plan (UFP SAP)*, Bethpage, New York and the UFP SAP Addendum Installation of Vertical Profile Borings and Monitoring Wells (Resolution Consultants, 2013). The field investigation included completing one vertical profile boring, groundwater grab samples, geophysical logging and surveying.

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

### 1.2 Site History

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

The Navy's property originally totaled 109.5 acres and was formerly a Government-Owned Contractor-Operated (GOCO) facility that was operated by Northrop Grumman (NG) until September 1998. Prior to 2002, the NWIRP property was bordered on the north, west, and south

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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 Nassau County property. 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 ft of Cretaceous 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 bgs observed onsite. The Magothy is characterized by fine to medium sands and silts interbedded with zones of clays, silty sands and sandy clays. Sand and gravel lenses are found in some areas between depths of 600 and 945 ft 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 in the offsite plume to date. 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.

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 regional groundwater flow in the area is to the south-southeast.

## **2.0 FIELD PROGRAM**

Field investigation activities at VPB 152 consisted of drilling, sampling, soil/groundwater analysis, geophysical logging, and surveying. Drilling during this investigation was performed by Delta Well and Pump Company of Ronkonkoma, New York. A description of these tasks is provided below.

### **2.1 Vertical Profile Borings**

One vertical profile boring (VPB 152) was completed during this field effort between May 23, 2014 and July 7, 2014. The total depth of VPB 152 was 997 ft. The location is shown in Figure 2 and details are summarized in Table 1.

#### **2.1.1 Drilling**

VPB 152 was installed by drilling an 8-inch diameter hole using mud rotary drilling techniques. Drilling mud consisted of potable water and polymer-free sodium bentonite or equivalent. Drilling mud was contained and re-circulated in baffled, high capacity mud tubs. A sand separator was used intermittently to remove fines from circulation.

#### **2.1.2 Sampling**

A total of eight split spoon samples were collected from ground surface to the bottom of the boring. A change in geology was observed by the field geologist at 980 ft bgs and three split spoon samples were subsequently collected to confirm the presence of the Raritan Clay. Samples were logged by the field geologist and screened for Volatile Organic Compounds (VOCs) utilizing a photoionization detector (PID). A detailed boring log for VPB 152 is included in Appendix A.

Groundwater grab samples were collected every 50 ft for the first 200 ft of borehole depth. After the first 200 ft, groundwater grab samples were collected approximately every 20 ft until the boring terminated in the Raritan. Groundwater grab samples were collected with a hydropunch sampler and analyzed for VOCs using Environmental Protection Agency (EPA) Method 8260C. The groundwater grab samples were analyzed by Katahdin Analytical Services (Katahdin), a Department of Defense (DoD), Environmental Laboratory Accreditation Program (ELAP), and New York State Department of Environmental Conservation (NYSDEC)-certified laboratory. During the collection of groundwater grab samples, field parameters were measured (pH, temperature, specific conductivity, oxidation reduction potential, dissolved oxygen, and turbidity). Data validation was performed by Resolution Consultants. Groundwater grab sample logs, data validation packages, and analytical data tables are included in Appendix A.

One soil sample was collected for laboratory analysis for total organic carbon (TOC) by EPA series SW-846 method 9060A. During drilling, air sampling was conducted under a Community Air Monitoring Plan. One air sample was collected per VPB using Summa canisters and submitted for laboratory analysis by EPA Method TO-15. All analyses were performed or sub-contracted by Katahdin. Data validation of both TOC and air data was performed by Resolution Consultants. Data validation packages and analytical data tables are included in Appendix A.

### **2.1.3 Geophysics**

Borehole geophysical logs (gamma) were recorded after the borehole was drilled but prior to the removal of drill rods. A Mount Sopris Instrument model 2PGA-100 poly gamma was used. Starting at the top of the hole, the probe was advanced at a maximum rate of 12 ft per minute. A copy of the log was printed in the field for review once the probe reached the bottom of the borehole. The instrument was then raised to the top of the boring and a second log was generated and printed in the field. The down hole gamma log sheets and plots comparing the gamma log with trichloroethene (TCE) and tetrachloroethene (PCE) concentrations from hydropunch samples are included in Appendix A.

## **2.2 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, split spoons, and the hydropunch 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.

As part of the IDW management practices and in accordance with the SAP, the investigation waste (consisting of soil cuttings, drilling muds, IDW fluids, and personal protective equipment (PPE)) generated during the boring installation was containerized and staged at NWIRP Bethpage. IDW solids were characterized and disposed of properly. 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 water 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 waters). To the extent feasible, soil and water were not mixed. All analytical criteria were met for disposal of soil and water.

### **2.3 Surveying**

A survey of the boring location 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 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

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

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

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

Resolution Consultants, 2013. UFP SAP Addendum, *Groundwater Sampling Using Low Stress (Low Flow) Purging and Sampling Protocol*. November.

Resolution Consultants, 2013. UFP SAP Addendum, *Installation of Vertical Profile Borings and Monitoring Wells*. December.

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

## Tables

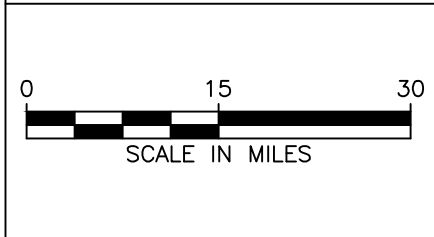
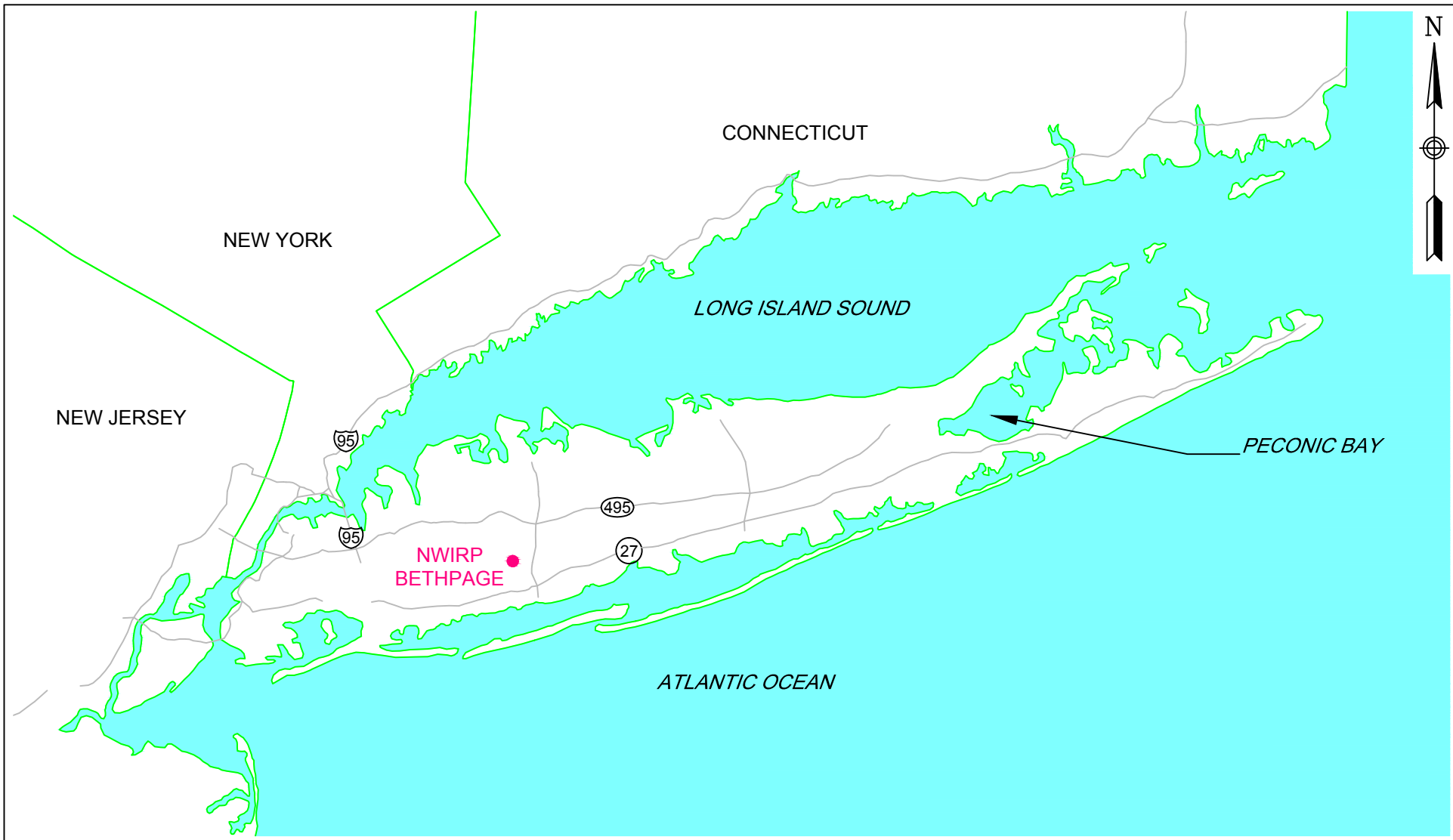


**TABLE 1  
 VERTICAL PROFILE BORING SUMMARY  
 2014 OU2 GROUNDWATER INVESTIGATION  
 NWIRP BETHPAGE, NY**

<b>BORING</b>	<b>BORING START DATE</b>	<b>BORING COMPLETION DATE</b>	<b>GROUND ELEVATION (MSL)</b>	<b>TOTAL DEPTH (ft bgs)</b>	<b>SURFACE CASING SET AT (ft bgs)</b>	<b>NO. OF SPOON SAMPLES</b>	<b>GAMMA LOG (ft bgs)</b>	<b>NO. GW SAMPLES COLLECTED/ ATTEMPTED</b>	<b>TOC SAMPLES</b>	<b>DATE OF AIR SAMPLE</b>	<b>MONITORING WELLS INSTALLED AT LOCATION</b>
VP 152	5/23/2014	7/7/2014	58.16	997	53.5	8	993	42/52*	325 - 327 ft bgs	6/24/2014	Pending

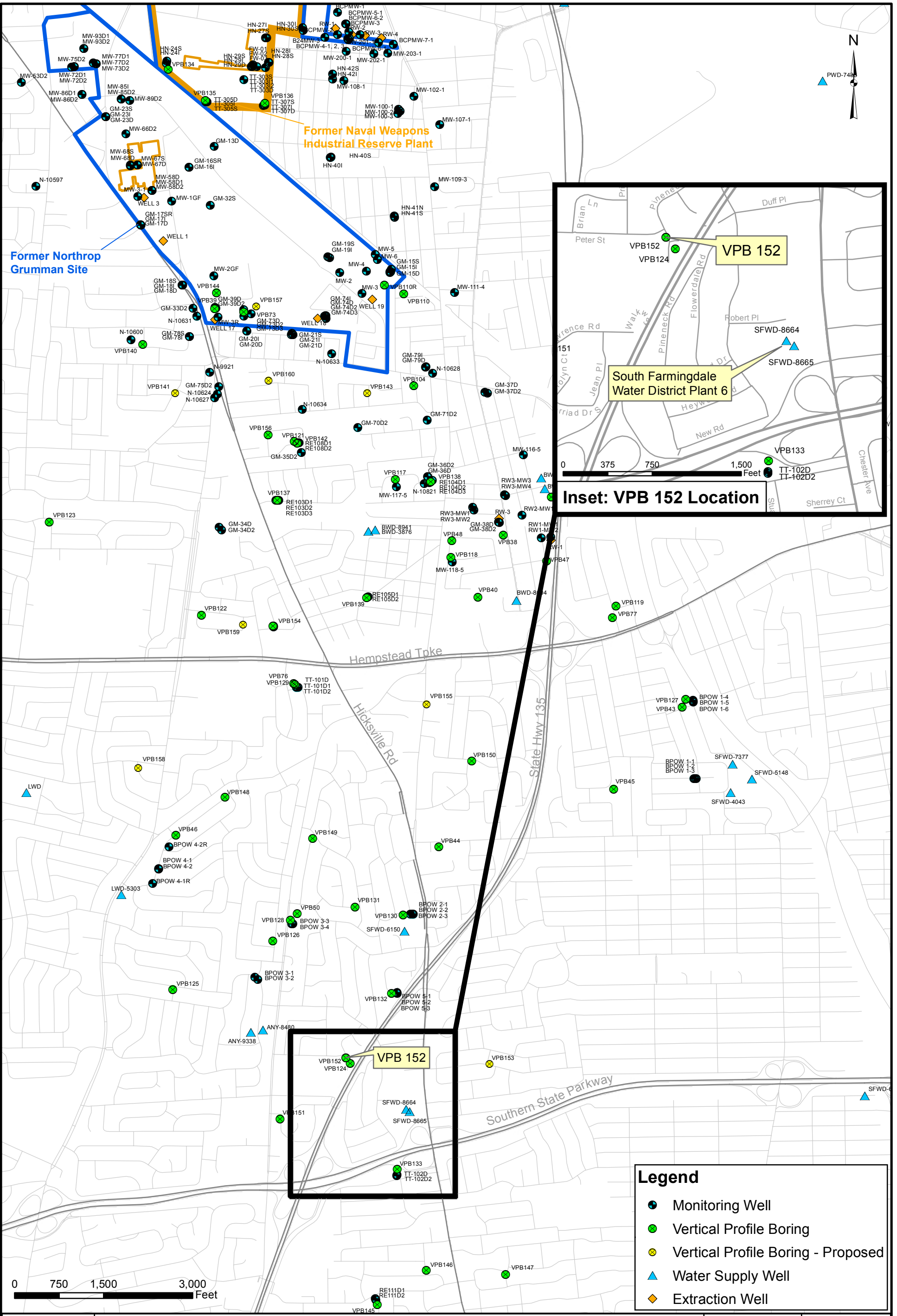
\* includes 2 field duplicates

## Figures



GENERAL LOCATION MAP  
NWIRP BETHPAGE  
BETHPAGE, NEW YORK

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



**VPB 152 LOCATION MAP**  
**NAVAL WEAPONS INDUSTRIAL RESERVE PLANT**  
**BETHPAGE, NEW YORK**

CONTRACT NUMBER N62470-11-D8013	CTO NUMBER WE15
APPROVED BY PS	DATE 12/10/2014
APPROVED BY	DATE
FIGURE NO. 2	REV 0

## **Appendix A**

**VPB 152**

**Section 1**

**VPB 152 Boring and Gamma Logs**

<b>Client:</b> Department of the Navy, Naval Facilities Engineering Command, Mid-Atlantic			<b>Logged By:</b> V. Varricchio		
<b>Location:</b> Pineneck Road, Bethpage		<b>Northing:</b> 197311.98		<b>Easting:</b> 1126289.12	
<b>Project #:</b> 60266526		<b>Ground Elevation (ft amsl):</b> 58.16		<b>Drilling Company:</b> Delta Well & Pump	
<b>Start Date:</b> 5/23/2014		<b>Drilling Method:</b> Auger (0-50' bgs) Mud Rotary (>50' bgs)		<b>Well Screen Interval (ft):</b> NA	
<b>Finish Date:</b> 7/7/2014				<b>Water Level (ft):</b> NA	
				<b>Total Depth (ft):</b> 997.0	

Mud Rotary Drilling Note: Unless denoted by a splitspoon sample (indicated by the presence of a PID reading), boundaries between strata are approximate only and may be transitional because they are based on screened wash samples collected during mud rotary drilling at 5 ft. intervals.

DEPTH (ft)	Gamma Ray	PID (ppm)	TCE (ug/L)	PCE (ug/L)	Formation	USCS	GRAPHIC LOG	MATERIAL DESCRIPTION
0								
2					Upper Glacial	SW		Top Soil
4						SW		Strong brown (7.5 YR 4/6) well graded SAND with some fine to coarse subrounded Gravel, trace silt
6						SW		Strong brown (7.5 YR 4/6) well graded SAND with fine to coarse subrounded Gravel, trace silt
8						SW		Strong brown (7.5 YR 4/6) well graded SAND with fine to coarse subrounded Gravel, trace silt
10						SW		Strong brown (7.5 YR 4/6) well graded SAND with fine to coarse subrounded Gravel, trace silt
12						SW		Strong brown (7.5 YR 4/6) well graded SAND with fine to coarse subrounded Gravel, trace silt
14						SW		Strong brown (7.5 YR 4/6) well graded SAND with fine to coarse subrounded Gravel, trace silt
16						SW		Strong brown (7.5 YR 4/6) well graded SAND with fine to coarse subrounded Gravel, trace silt
18						SW		Strong brown (7.5 YR 4/6) well graded SAND with fine to coarse subrounded Gravel, trace silt
20						SW		Strong brown (7.5 YR 4/6) well graded SAND with fine to coarse subrounded Gravel, trace silt
22						SW		Strong brown (7.5 YR 4/6) well graded SAND with fine to coarse subrounded Gravel, trace silt
24						SW		Strong brown (7.5 YR 4/6) well graded SAND with fine to coarse subrounded Gravel, trace silt
26						SW		Strong brown (7.5 YR 4/6) well graded SAND with fine to coarse subrounded Gravel, trace silt
28						SW		Strong brown (7.5 YR 4/6) well graded SAND with fine to coarse subrounded Gravel, trace silt
30						SW		Strong brown (7.5 YR 4/6) well graded SAND with fine to coarse subrounded Gravel, trace silt
32						SW		Strong brown (7.5 YR 4/6) well graded SAND with fine to coarse subrounded Gravel, trace silt
34						SW		Strong brown (7.5 YR 4/6) well graded SAND with fine to coarse subrounded Gravel, trace silt
36						SW		Strong brown (7.5 YR 4/6) well graded SAND with fine to coarse subrounded Gravel
38						SW		Strong brown (7.5 YR 4/6) well graded SAND with fine to coarse subrounded Gravel, trace silt
40						SW		Strong brown (7.5 YR 4/6) well graded SAND with fine to coarse subrounded Gravel, trace silt
42						SW		Yellowish brown (10 YR 5/6) well graded SAND with fine to coarse subrounded Gravel, trace silt
44						SW		Yellowish brown (10 YR 5/6) well graded SAND with fine to coarse subrounded Gravel, trace silt
46						SW		Yellowish brown (10 YR 5/6) well graded SAND with fine to coarse subrounded Gravel, trace silt
48						SW		Yellowish brown (10 YR 5/6) well graded SAND with fine to coarse subrounded Gravel, trace silt
50						SW		Yellowish brown (10 YR 5/6) well graded SAND with fine to coarse subrounded Gravel, trace silt
52						GW		Yellowish brown (10 YR 5/6) well graded fine to coarse subrounded GRAVEL with some medium to coarse Sand
54								

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DEPTH (ft)	Gamma Ray	PID (ppm)	TCE (ug/L)	PCE (ug/L)	Formation	USCS	GRAPHIC LOG	MATERIAL DESCRIPTION
54	30 60 90							
56					Upper Glacial	GW		Yellowish brown (10 YR 5/6) fine subrounded GRAVEL with few coarse Sand
58				GP				
60			< 0.50	< 0.50				
62						GP		Yellowish brown (10 YR 5/4) poorly graded fine subrounded GRAVEL with medium to coarse Sand
64								
66						GW		Yellowish brown (10 YR 5/4) well graded subrounded GRAVEL with few coarse Sand, trace lignite
68								
70								
72						SP		Very pale brown (10 YR 7/3) poorly graded coarse SAND with some fine Gravel, trace silt
74								
76					SP		Very pale brown (10 YR 7/3) poorly graded coarse SAND with some fine Gravel, trace silt	
78								
80								
82					SP		Very pale brown (10 YR 7/3) poorly graded coarse SAND with some fine Gravel, trace silt	
84								
86					GW		Grayish brown (10 YR 5/2) well graded subrounded GRAVEL and angular iron nodules with some medium to coarse Sand, trace silt	
88								
90								
92					SW		Grayish brown (10 YR 5/2) well graded SAND with fine subrounded Gravel, trace silt, iron nodules and lignite	
94								
96								
98					SW		Grayish brown (10 YR 5/2) well graded SAND with fine subrounded Gravel, trace silt, iron nodules and lignite	
100			< 0.50	< 0.50				
102					Magothy	SW-SM		Light brownish gray (10 YR 6/2) well graded SAND with Silt and subrounded gravel, trace lignite and iron nodules
104								
106						SW		Light brownish gray (10 YR 6/2) well graded SAND with subrounded Gravel, trace silt, iron nodules and lignite
108								
110								
112					SW		Dark gray (10 YR 4/1) well graded SAND and Lignite, trace iron nodules and silt	
114								

(Continued Next Page)



DEPTH (ft)	Gamma Ray	PID (ppm)	TCE (ug/L)	PCE (ug/L)	Formation	USCS	GRAPHIC LOG	MATERIAL DESCRIPTION
116	30 60 90				Magothy	SW		Dark gray (10 YR 4/1) well graded SAND and Lignite, trace silt and fine subrounded gravel ( <i>continued</i> )
118				SW				
120								
122						SP		Dark gray (10 YR 4/1) poorly graded fine SAND with some lignite, trace Silt
124								
126						SP-SM		Dark gray (10 YR 4/1) poorly graded fine SAND mixed with Silt and fine subrounded gravel, trace lignite
128								
130						SM		Dark gray (10 YR 4/1) SILTY SAND with few fine subrounded Gravel, trace lignite
132								
134						ML		Gray (10 YR 5/1) SANDY SILT with few fine subrounded Gravel, trace lignite
136								
138						SW		Gray (10 YR 6/1) Well graded SAND, trace subrounded fine Gravel, clay and iron nodules
140								
142						SW		Gray (10 YR 6/1) Well graded SAND, trace subrounded fine Gravel, clay and iron nodules
144								
146						SW		Gray (10 YR 6/1) Well graded SAND, trace subrounded fine Gravel, clay and iron nodules
148								
150						SP		Gray (10 YR 6/1) poorly graded medium SAND with lignite, trace Clay
152			< 0.50	< 0.50				
154						SP		Gray (10 YR 6/1) poorly graded medium SAND with lignite, trace Clay
156								
158					CL		Gray (10 YR 5/1) lean CLAY with some fine Sand, trace silt and lignite	
160								
162					ML		Gray (10 YR 6/1) sandy SILT with trace rounded Gravel	
164								
166					SM		Gray (10 YR 5/1) silty SAND, trace lignite	
168								
170					MH-OL		Gray (10 YR 5/1) micaceous SILT mixed with poorly graded fine to medium subangular Sand and lignite	
172								
174								
176								

(Continued Next Page)

DEPTH (ft)	Gamma Ray 30 60 90	PID (ppm)	TCE (ug/L)	PCE (ug/L)	Formation	USCS	GRAPHIC LOG	MATERIAL DESCRIPTION
178					Magothy	MH-OL		Gray (10 YR 5/1) micaceous SILT mixed with poorly graded fine to medium subangular Sand and lignite (continued)
180						CH		Dark gray (10 YR 4/1) fine to medium sandy fat CLAY and lignite, trace iron nodules
182						CH		Dark gray (10 YR 4/1) well graded fine sandy fat CLAY and abundant wood chips, trace lignite
184						CH		Dark Gray (10 YR 4/1) poorly graded fine and some medium sandy fat CLAY, abundant lignite and some wood chips
186						CH		Dark Gray (10 YR 4/1) poorly graded fine and some medium sandy fat CLAY, abundant lignite and some wood chips
188						CH		Dark Gray (10 YR 4/1) poorly graded fine and some medium sandy fat CLAY, abundant lignite and some wood chips
190						CH		Dark Gray (10 YR 4/1) poorly graded fine and some medium sandy fat CLAY, abundant lignite and some wood chips
192						CH		Dark Gray (10 YR 4/1) poorly graded fine and some medium sandy fat CLAY, abundant lignite and some wood chips
194						CH		Dark Gray (10 YR 4/1) poorly graded fine and some medium sandy fat CLAY, abundant lignite and some wood chips
196						SM		Very dark gray (10 YR 3/1) silty poorly graded fine and some medium SAND with abundant lignite and wood chips
198						SM		Very dark gray (10 YR 3/1) silty well graded fine to medium SAND with abundant lignite
200						SM		Very dark gray (10 YR 3/1) silty well graded fine to medium SAND with abundant lignite
202			< 0.50	< 0.50		SM		Very dark gray (10 YR 3/1) silty well graded fine to medium SAND with abundant lignite
204						SM		Very dark gray (10 YR 3/1) silty well graded fine to medium SAND with abundant lignite
206						SM		Very dark gray (10 YR 3/1) silty well graded fine to medium SAND with abundant lignite
208						SM		Very dark gray (10 YR 3/1) silty well graded fine to medium SAND with abundant lignite
210						SM		Very dark gray (10 YR 3/1) silty well graded fine to medium SAND with abundant lignite
212						SM		Very dark gray (10 YR 3/1) silty well graded fine to medium SAND with abundant lignite
214						SM		Very dark gray (10 YR 3/1) silty well graded fine to medium SAND with abundant lignite
216						SM		Very dark gray (10 YR 3/1) silty poorly graded fine SAND with abundant lignite, trace gravel
218					SM		Very dark gray (10 YR 3/1) silty poorly graded fine SAND with abundant lignite, trace gravel	
220					SM		Very dark gray (10 YR 3/1) silty poorly graded fine SAND with abundant lignite, trace gravel	
222			< 0.50	< 0.50	SM		Very dark gray (10 YR 3/1) silty fine to medium SAND with abundant lignite	
224					SM		Very dark gray (10 YR 3/1) silty fine to medium SAND with abundant lignite	
226					SM		Very dark gray (10 YR 3/1) silty well graded fine to coarse SAND with abundant lignite, trace Clay and rounded gravel	
228					SM		Very dark gray (10 YR 3/1) silty well graded fine to coarse SAND with abundant lignite, trace Clay and rounded gravel	
230					SM-SC		Very dark gray (10 YR 3/1) silty clayey well graded fine to medium SAND with abundant lignite	
232					SM-SC		Very dark gray (10 YR 3/1) silty clayey well graded fine to medium SAND with abundant lignite	
234					SM-SC		Very dark gray (10 YR 3/1) silty clayey well graded fine to medium SAND with abundant lignite	
236					SM-SC		Very dark gray (10 YR 3/1) silty clayey well graded fine to medium SAND with abundant lignite	
238					SM-SC		Very dark gray (10 YR 3/1) silty clayey well graded fine to medium SAND with abundant lignite	

(Continued Next Page)

DEPTH (ft)	Gamma Ray 30 60 90	PID (ppm)	TCE (ug/L)	PCE (ug/L)	Formation	USCS	GRAPHIC LOG	MATERIAL DESCRIPTION
240					Magothy			
242			< 0.50	< 0.50		SM		Very dark gray (10 YR 3/1) silty well graded fine to medium SAND with abundant lignite
244						SM		
246						SM		Very dark gray (10 YR 3/1) silty well graded fine to medium SAND with abundant lignite
248						SM		
250						SM		
252						SM		Very dark gray (10 YR 3/1) silty poorly graded fine SAND with abundant lignite
254						SM		
256						SM		Very dark gray (10 YR 3/1) silty poorly graded fine SAND with abundant lignite
258						SM		
260						SM		
262			< 0.50	< 0.50		CL		Very dark gray (10 YR 3/1) sandy lean CLAY with abundant Lignite
264						CL		
266						CL		Very dark gray (10 YR 3/1) lean CLAY with fine to medium well graded Sand and abundant lignite
268						CL		
270						ML		Very dark gray (10 YR 3/1) SILT with poorly graded fine Sand and abundant lignite
272						ML		
274						ML		Very dark gray (10 YR 3/1) SILT with poorly graded fine SAND and abundant lignite, trace clay
276						ML		
278						ML		
280						ML		Very dark gray (10 YR 3/1) sandy SILT with abundant lignite
282			< 0.50	< 0.50		ML		
284						ML		
286						SM-SC		Very dark gray (10 YR 3/1) silty clayey poorly graded fine SAND with some lignite
288						SM-SC		
290						SM-SC		
292						SW-SM		Very dark gray (10 YR 3/1) well graded fine to medium SAND mixed with Silt and some lignite, trace clay
294						SW-SM		
296						SW-SM		
298						SP		Very dark gray (10 YR 3/1) poorly graded fine SAND, trace lignite
300			< 0.50	< 0.50		SP		

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DEPTH (ft)	Gamma Ray 30 60 90	PID (ppm)	TCE (ug/L)	PCE (ug/L)	Formation	USCS	GRAPHIC LOG	MATERIAL DESCRIPTION	
302			< 0.50	< 0.50	Magothy	SP		Very dark gray (10 YR 3/1) poorly graded fine SAND, trace lignite (continued)	
304						SP			
306						SM			Very dark gray (10 YR 3/1) silty poorly graded fine SAND with some lignite
308									
310						SM			Very dark gray (10 YR 3/1) silty poorly graded fine SAND with lignite
312									
314						ML			Very dark gray (10 YR 3/1) SILT with poorly graded fine Sand and lignite
316									
318						CL			Very dark gray (10 YR 3/1) sandy lean CLAY with abundant lignite
320			< 0.50	< 0.50					
322						CL			Very dark gray (10 YR 3/1) interbedded sandy lean CLAY with abundant lignite
324									
326		0.0				SM			Very dark gray (10 YR 3/1) silty poorly graded fine SAND with some lignite
328									
330						ML			Very dark gray (10 YR 3/1) SILT with poorly graded fine Sand and some lignite
332									
334					ML		Very dark gray (10 YR 3/1) SILT with well graded fine to medium Sand and abundant lignite		
336									
338					ML		Very dark gray (10 YR 3/1) SILT with poorly graded fine Sand and some lignite		
340									
342					ML		Very dark gray (10 YR 3/1) SILT with trace fine Sand, clay and lignite		
344									
346			< 0.50	< 0.50	ML		Very dark gray (10 YR 3/1) SILT with poorly graded fine Sand, trace lignite		
348									
350					ML		Very dark gray (10 YR 3/1) silty poorly graded fine SAND with lignite		
352									
354					SM		Very dark gray (10 YR 3/1) silty poorly graded fine SAND with lignite		
356									
358					ML		Very dark gray (10 YR 3/1) SILT with few poorly graded fine Sand, trace lignite		
360									
362			< 0.50	< 0.50					

(Continued Next Page)

DEPTH (ft)	Gamma Ray 30    60    90	PID (ppm)	TCE (ug/L)	PCE (ug/L)	Formation	USCS	GRAPHIC LOG	MATERIAL DESCRIPTION		
364					Magothy	ML		Very dark gray (10 YR 3/1) SILT with few poorly graded fine Sand, trace lignite (continued)		
366						ML		Very dark gray (10 YR 3/1) SILT, trace poorly graded fine Sand		
368						ML		Very dark gray (10 YR 3/1) SILT, trace poorly graded fine Sand and lignite		
370						ML		Very dark gray (10 YR 3/1) SILT, trace poorly graded fine Sand and lignite		
372						ML		Very dark gray (10 YR 3/1) SILT, trace poorly graded fine Sand and lignite		
374						ML		Very dark gray (10 YR 3/1) SILT, trace poorly graded fine Sand and lignite		
376						ML		Very dark gray (10 YR 3/1) SILT, trace poorly graded fine Sand and lignite		
378						ML		Very dark gray (10 YR 3/1) SILT, trace poorly graded fine Sand and lignite		
380						ML		Very dark gray (10 YR 3/1) SILT, trace poorly graded fine Sand and lignite		
382						< 0.50		< 0.50	ML	Gray (10 YR 5/1) poorly graded fine sandy SILT, trace lignite
384						ML		Gray (10 YR 5/1) lean CLAY with trace poorly graded fine Sand and lignite		
386						CL		Gray (10 YR 5/1) CLAYEY SILT with trace poorly graded fine Sand and lignite		
388						ML-CL		Very dark gray (10 YR 3/1) poorly graded fine sandy CLAY with lignite		
390						CL		Dark gray (10 YR 4/1) poorly graded fine sandy SILT, trace lignite		
392						ML		Dark gray (10 YR 4/1) silty well graded fine to medium SAND, trace Clay and lignite		
394						SM		Dark gray (10 YR 4/1) SILT with some poorly graded fine Sand		
396						ML		Gray (10 YR 5/1) silty poorly graded fine SAND, trace lignite		
398						CL		Dark gray (10 YR 4/1) SILT with poorly graded fine Sand, trace lignite		
400	< 0.50	< 0.50	ML	Dark gray (10 YR 4/1) poorly graded fine sandy SILT, trace lignite						
402	ML	Dark gray (10 YR 4/1) silty well graded fine to medium SAND, trace Clay and lignite								
404	SM	Dark gray (10 YR 4/1) SILT with some poorly graded fine Sand								
406	ML	Gray (10 YR 5/1) silty poorly graded fine SAND, trace lignite								
408	SM	Dark gray (10 YR 4/1) SILT with poorly graded fine Sand, trace lignite								
410	ML	Dark gray (10 YR 4/1) SILT with poorly graded fine Sand, trace lignite								
412	ML	Dark gray (10 YR 4/1) SILT with poorly graded fine Sand, trace lignite								
414	ML	Dark gray (10 YR 4/1) SILT with poorly graded fine Sand, trace lignite								
416	SM	Dark gray (10 YR 4/1) SILT with poorly graded fine Sand, trace lignite								
418	SM	Dark gray (10 YR 4/1) SILT with poorly graded fine Sand, trace lignite								
420	ML	Dark gray (10 YR 4/1) SILT with poorly graded fine Sand, trace lignite								
422	ML	Dark gray (10 YR 4/1) SILT with poorly graded fine Sand, trace lignite								
424	ML	Dark gray (10 YR 4/1) SILT with poorly graded fine Sand, trace lignite								

(Continued Next Page)

DEPTH (ft)	Gamma Ray 30 60 90	PID (ppm)	TCE (ug/L)	PCE (ug/L)	Formation	USCS	GRAPHIC LOG	MATERIAL DESCRIPTION	
426					Magothy			Dark gray (10 YR 4/1) SILT, trace poorly graded fine Sand and lignite	
428						ML			
430									
432						ML			Dark gray (10 YR 4/1) SILT with some poorly graded fine Sand, trace lignite
434									
436						CL-ML			Dark gray (10 YR 4/1) poorly graded fine sandy clayey SILT with trace lignite
438									
440									
442			< 0.50	< 0.50		ML			Dark gray (10 YR 4/1) poorly graded fine sandy SILT
444									
446						ML			Very dark gray (10 YR 3/1) SILT with few poorly graded fine Sand, trace lignite and clay
448									
450						ML			Very dark gray (10 YR 3/1) SILT with some poorly graded fine Sand and lignite
452									
454					ML			Very dark gray (10 YR 3/1) poorly graded fine sandy clayey SILT, abundant lignite	
456									
458					ML-CL				
460									
462			< 0.50	< 0.50	ML			Very dark gray (10 YR 3/1) poorly graded fine sandy SILT, trace lignite	
464									
466					SC			Very dark gray (10 YR 3/1) well graded fine to medium silty SAND with abundant lignite	
468									
470									
472					ML			Dark gray (10 YR 4/1) SILT with poorly graded fine SAND and Lignite, trace clay	
474									
476					SW-SM			Dark gray (10 YR 4/1) well graded fine to medium SAND mixed with Silt, trace lignite	
478									
480									
482			< 0.50	< 0.50	SW			Dark gray (10 YR 4/1) well graded fine to medium SAND with some Silt, trace lignite	
484									
486					SM-SC				

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DEPTH (ft)	Gamma Ray	PID (ppm)	TCE (ug/L)	PCE (ug/L)	Formation	USCS	GRAPHIC LOG	MATERIAL DESCRIPTION
486	30 60 90				Magothy			
488						SM-SC		Very dark gray (10 YR 3/1) silty clayey poorly graded fine SAND, trace lignite (continued)
490						ML		Dark gray (10 YR 4/1) well graded fine to medium sandy SILT, trace lignite and clay
492								Dark gray (10 YR 4/1) clayey SILT mixed with few poorly graded fine Sand, trace lignite
494						ML-CL		Dark gray (10 YR 4/1) lean CLAY, trace poorly graded fine Sand and lignite
496								Dark gray (10 YR 4/1) poorly graded fine sandy lean CLAY, trace lignite
498						CL		Dark gray (10 YR 4/1) poorly graded fine sandy SILT, trace lignite
500								Dark gray (10 YR 4/1) poorly graded fine sandy clayey SILT with some lignite
502			< 0.50	< 0.50		CL		Gray (10 YR 5/1) poorly graded fine sandy SILT with lignite
504								Gray (10 YR 5/1) well graded fine to coarse subangular SAND with some lignite, trace Silt
506						CL		Gray (10 YR 5/1) well graded fine to coarse subangular SAND with some lignite and silt
508								Gray (10 YR 5/1) poorly graded fine SAND with some lignite
510						ML		Dark gray (10 YR 4/1) poorly graded fine SAND with some lignite, trace Silt
512								Poorly graded fine sandy clayey SILT with some lignite
514						ML-CL		
516								
518						ML		
520								
522			< 0.50	< 0.50		ML		
524								
526					SW			
528								
530					SW			
532								
534					SP			
536								
538					SP			
540								
542			< 0.50	< 0.50	SP			
544								
546					ML-CL			

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DEPTH (ft)	Gamma Ray	PID (ppm)	TCE (ug/L)	PCE (ug/L)	Formation	USCS	GRAPHIC LOG	MATERIAL DESCRIPTION	
548	30 60 90				Magothy	ML-CL		Poorly graded fine sandy clayey SILT with some lignite (continued)	
550						ML		Dark gray (10 YR 4/1) poorly graded fine sandy SILT with some Lignite	
552									
554									
556									
558						CL-ML		Dark gray (10 YR 4/1) silty CLAY with abundant lignite/wood	
560									
562			< 0.50	< 0.50			SW		Gray (10 YR 5/1) well graded fine to medium SAND, trace Clay and lignite
564									
566							SW		Dark gray (10 YR 4/1) well graded fine to medium SAND with Clay, trace lignite
568									
570									
572							CL		Dark gray (10 YR 4/1) lean CLAY with trace poorly graded fine Sand
574									
576									
578							ML		Dark gray (10 YR 4/1) SILT with little poorly graded fine Sand, trace lignite
580									
582			< 0.50	< 0.50			CH		Dark gray (10 YR 4/1) fat CLAY with little poorly graded fine Sand
584									
586		0.0					CH		Dark gray (10 YR 4/1) fat CLAY with little poorly graded fine Sand
588						SP		Gray (10 YR 5/1) poorly graded fine SAND	
590									
592						SW		Dark gray (10 YR 4/1) well graded fine to medium SAND, trace lignite	
594									
596						SW		Dark gray (10 YR 4/1) well graded fine to coarse subrounded SAND, trace lignite	
598									
600									
602						SW		Gray (10 YR 5/1) well graded fine to coarse subrounded SAND, trace Clay and lignite	
604									
606			< 0.50	< 0.50		SW		Gray (10 YR 5/1) well graded fine to medium SAND, trace lignite	
608									

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DEPTH (ft)	Gamma Ray 30 60 90	PID (ppm)	TCE (ug/L)	PCE (ug/L)	Formation	USCS	GRAPHIC LOG	MATERIAL DESCRIPTION
610					Magothy			
612						SW-SM		Gray (10 YR 5/1) well graded fine to medium SAND mixed with Silt, trace lignite
614								
616								
618						SW		Gray (10 YR 5/1) well graded fine to medium SAND with trace Silt and lignite
620								
622			< 0.50	< 0.50				
624						CL		Dark gray (10 YR 4/1) lean CLAY with little poorly graded fine Sand and trace lignite
626								
628						CL		Dark gray (10 YR 4/1) lean CLAY with little poorly graded fine Sand and trace lignite
630								
632								
634						SW-SC		Gray (10 YR 5/1) well graded fine to coarse subangular SAND mixed with some Clay, trace lignite
636								
638						SW-SC		Gray (10 YR 5/1) well graded fine to coarse subangular SAND mixed with some Clay and lignite
640								
642			< 0.50	< 0.50				
644						SM-SC		Gray (10 YR 5/1) clayey silty poorly graded fine SAND, trace Lignite
646								
648					SM-SC		Gray (10 YR 5/1) clayey silty poorly graded fine SAND, trace Lignite	
650								
652					SM-SC		Gray (10 YR 5/1) clayey silty poorly graded fine SAND, trace Lignite	
654								
656					SM-SC		Gray (10 YR 5/1) clayey silty well graded fine to medium SAND, trace lignite	
658								
660								
662			< 0.50	< 0.50				
664					CH		Gray (10 YR 5/1) fat CLAY, trace poorly graded fine Sand	
666								
668					CH		Gray (10 YR 5/1) fat CLAY, trace poorly graded fine Sand	
670								
					CH-SW			

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DEPTH (ft)	Gamma Ray 30 60 90	PID (ppm)	TCE (ug/L)	PCE (ug/L)	Formation	USCS	GRAPHIC LOG	MATERIAL DESCRIPTION	
672					Magothy	CH-SW		Gray (10 YR 5/1) interbedded fat CLAY and well graded subrounded fine to coarse Sand with some lignite (continued)	
674						CL		Gray (10 YR 5/1) lean CLAY with some poorly graded fine Sand, trace lignite	
676									
678									
680									
682			< 0.50	< 0.50		ML		Gray (10 YR 5/1) poorly graded fine sandy SILT	
684									
686									
688						CL		Gray (10 YR 5/1) lean CLAY	
690									
692		0.0				SP-CL		Gray (10 YR 5/1) interbedded poorly graded fine SAND and lean Clay	
694									
696						SW-CH		Gray (10 YR 5/1) interbedded well graded fine to coarse SAND and light gray (10 YR 7/1) fat Clay	
698									
700									
702			< 50	< 50	CL		Gray (10 YR 5/1) lean CLAY with some poorly graded fine Sand, trace lignite and fine subrounded gravel		
704									
706									
708					SW-CL		Gray (10 YR 5/1) interbedded well graded fine to coarse SAND and lean Clay, trace subrounded fine gravel		
710									
712					SW-CL		Gray (10 YR 5/1) interbedded well graded fine to coarse SAND and lean Clay, trace subrounded fine gravel		
714									
716									
718					SW-CL		Gray (10 YR 5/1) interbedded well graded fine to coarse SAND and lean Clay, trace subrounded fine gravel		
720									
722					SW-CL		Gray (10 YR 5/1) interbedded well graded fine to coarse SAND and lean Clay, trace subrounded fine gravel		
724									
726			< 25	< 25	GC-CL		Gray (10 YR 6/1) interbedded clayey fine angular GRAVEL and lean Clay		
728									
730									
732					GC-CL		Gray (10 YR 6/1) interbedded clayey fine angular GRAVEL and lean Clay		

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DEPTH (ft)	Gamma Ray 30 60 90	PID (ppm)	TCE (ug/L)	PCE (ug/L)	Formation	USCS	GRAPHIC LOG	MATERIAL DESCRIPTION
734					Magothy	GC-CL		Gray (10 YR 6/1) interbedded clayey fine angular GRAVEL and lean Clay (continued)
736						SW-CL		Gray (10 YR 5/1) interbedded well graded fine to coarse SAND and lean Clay, trace subangular fine gravel
738								
740								
742			< 0.50	< 0.50		ML		Gray (10 YR 5/1) well graded fine to coarse sandy SILT, trace subangular Gravel
744								
746						GC-CL		Gray (10 YR 6/1) clayey fine angular GRAVEL and lean Clay with some well graded fine to coarse sand
748								
750								
752						GP		Very pale brown (10 YR 7/3) poorly graded fine subangular GRAVEL with some well graded fine to coarse Sand
754								
756						GP		Very pale brown (10 YR 7/3) poorly graded fine subangular GRAVEL with some well graded fine to coarse Sand, trace lean clay
758						GP		Very pale brown (10 YR 7/3) poorly graded fine subangular GRAVEL with some well graded fine to coarse Sand, trace lean clay
760								
762					GP		Very pale brown (10 YR 7/3) poorly graded fine subangular GRAVEL with some well graded fine to coarse Sand, trace lean clay	
764								
766					GP-GM		Very pale brown (10 YR 7/3) poorly graded fine subangular GRAVEL mixed with white (10 YR 8/1) Clay and well graded fine to coarse sand	
768								
770					GP-CH		Very pale brown (10 YR 7/3) interbedded poorly graded fine subangular GRAVEL with little well graded Sand and white (10 YR 8/1) fat clay	
772								
774								
776					CH		White (10 YR 8/1) well graded fine to coarse sandy fat CLAY with fine subangular Gravel	
778								
780								
782			< 0.50	< 0.50	SW-SC		Very pale brown (10 YR 7/3) well graded fine to coarse SAND mixed with some white (10 YR 8/1) Clay	
784								
786					SW		Very pale brown (10 YR 7/3) well graded fine to coarse SAND with few white (10 YR 8/1) Clay	
788								
790								
792					CL		Dark gray (10 YR 4/1) well graded fine to coarse sandy lean CLAY with trace fine subrounded Gravel	
794								

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DEPTH (ft)	Gamma Ray 30 60 90	PID (ppm)	TCE (ug/L)	PCE (ug/L)	Formation	USCS	GRAPHIC LOG	MATERIAL DESCRIPTION	
796					Magothy			Dark gray (10 YR 4/1) lean CLAY with trace Sand	
798						CL			
800									
802			< 0.50	< 0.50					Dark gray (10 YR 4/1) lean CLAY with trace Sand
804						CL			
806									Gray (10 YR 5/1) sandy silty lean CLAY
808						CL-ML			
810									Gray (10 YR 5/1) sandy silty lean CLAY
812						CL-ML			
814									Gray (10 YR 5/1) fine subangular gravelly lean CLAY with well graded fine to coarse Sand
816						CL			
818									Gray (10 YR 6/1) poorly graded fine SAND mixed with little Clay, trace fine subangular gravel
820						SP-SC			
822									Gray (10 YR 5/1) well graded fine to medium sandy silty CLAY
824									
826						CL-ML			
828									Light gray (10 YR 7/1) well graded fine to medium sandy clayey SILT
830			< 0.50	< 0.50					
832						ML-CL			
834								Light gray (10 YR 7/1) well graded fine to medium clayey silty SAND	
836									
838					SM-SC				
840								Light gray (10 YR 7/1) well graded fine to coarse clayey silty SAND	
842					SM-SC				
844								Gray (10 YR 6/1) fat CLAY, trace poorly graded fine Sand	
846		0.0			CH				
848								Gray (10 YR 6/1) poorly graded fine SAND, trace Silt	
850					SP				
852			< 0.50	< 0.50				Gray (10 YR 6/1) poorly graded fine sandy lean CLAY, trace Lignite	
854					CL				
856									
					CL-CH				

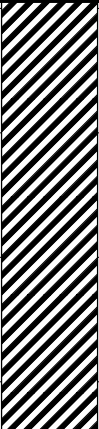
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DEPTH (ft)	Gamma Ray 30 60 90	PID (ppm)	TCE (ug/L)	PCE (ug/L)	Formation	USCS	GRAPHIC LOG	MATERIAL DESCRIPTION
858					Magothy	CL-CH		Gray (10 YR 5/1) mixed lean/fat CLAY with few poorly graded fine Sand ( <i>continued</i> )
860						CL-CH		Light gray (10 YR 7/1) mixed lean/fat CLAY, trace poorly graded fine Sand
862						CL-CH		Light gray (10 YR 7/1) mixed lean/fat CLAY, trace poorly graded fine Sand
864						CL-CH		Light gray (10 YR 7/1) mixed lean/fat CLAY, trace poorly graded fine Sand
866						CL-CH		Light gray (10 YR 7/1) mixed lean/fat CLAY, trace poorly graded fine Sand
868						CL-CH		Light gray (10 YR 7/1) mixed lean/fat CLAY, trace poorly graded fine Sand
870						CL-CH		Light gray (10 YR 7/1) mixed lean/fat CLAY, trace poorly graded fine Sand
872			< 0.50	< 0.50		CL-CH		Light gray (10 YR 7/1) mixed lean/fat CLAY, trace poorly graded fine Sand
874						CL-CH		Light gray (10 YR 7/1) mixed lean/fat CLAY, trace poorly graded fine Sand
876						CL-CH		Light gray (10 YR 7/1) mixed lean/fat CLAY, trace poorly graded fine Sand
878						CL-CH		Light gray (10 YR 7/1) mixed lean/fat CLAY, trace poorly graded fine Sand
880						CH		Gray (10 YR 5/1) fat CLAY, trace poorly graded fine Sand
882			< 0.50	< 0.50		CH		Gray (10 YR 5/1) fat CLAY, trace poorly graded fine Sand
884						CL-ML		Gray (10 YR 6/1) silty CLAY, trace poorly graded fine Sand and lignite
886						CL-ML		Gray (10 YR 6/1) silty CLAY with few poorly graded fine Sand
888					CH	Gray (10 YR 5/1) fat CLAY, trace poorly graded fine Sand		
890					CH	Gray (10 YR 6/1) fat CLAY		
892					CH	Gray (10 YR 6/1) fat CLAY, trace poorly graded fine Sand		
894					CH	Gray (10 YR 6/1) fat CLAY, trace poorly graded fine Sand		
896					CH	Gray (10 YR 6/1) fat CLAY, trace poorly graded fine Sand		
898					CH	Gray (10 YR 6/1) fat CLAY, trace poorly graded fine Sand		
900					CH	Gray (10 YR 6/1) fat CLAY, trace poorly graded fine Sand		
902					CH	Gray (10 YR 6/1) fat CLAY, trace poorly graded fine Sand		
904					CH	Gray (10 YR 6/1) fat CLAY, trace poorly graded fine Sand		
906					CH	Gray (10 YR 6/1) fat CLAY, trace poorly graded fine Sand		
908					CH	Gray (10 YR 6/1) fat CLAY, trace poorly graded fine Sand		
910					CH	Gray (10 YR 6/1) fat CLAY, trace poorly graded fine Sand		
912			< 0.50	< 0.50	CH	Gray (10 YR 6/1) fat CLAY, trace lignite		
914					CH	Gray (10 YR 6/1) fat CLAY, trace lignite		
916					CH	Gray (10 YR 6/1) fat CLAY, trace lignite		
918					CH	Gray (10 YR 6/1) fat CLAY, trace lignite		

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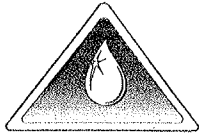
DEPTH (ft)	Gamma Ray	PID (ppm)	TCE (ug/L)	PCE (ug/L)	Formation	USCS	GRAPHIC LOG	MATERIAL DESCRIPTION	
918	30 60 90				Magothy			Gray (10 YR 6/1) fat CLAY, trace lignite <i>(continued)</i>	
920						CH		Gray (10 YR 5/1) fat CLAY, trace poorly graded fine Sand	
922						CH			
924									
926		0.0				CH			Gray (10 YR 5/1) fat CLAY, trace poorly graded fine Sand with microlamination
928									
930									
932			< 1.0	< 1.0		SM-SC			Gray (10 YR 6/1) silty clayey well graded fine to coarse SAND, trace lignite
934									
936						SM-SC			Gray (10 YR 6/1) silty clayey well graded fine to coarse SAND, trace lignite
938						SM-SC			
940									
942						SW			Gray (10 YR 6/1) well graded fine to medium SAND, trace Silt and clay
944									
946						SW			Gray (10 YR 6/1) well graded fine to coarse SAND with fine subrounded Gravel, trace silt and clay
948									
950									
952			< 1.0	< 1.0		SW-SC			Very dark gray (10 YR 3/1) well graded fine to coarse SAND mixed with fine subrounded Gravel and clay, trace lignite
954									
956					CH		Very dark gray (10 YR 3/1) well graded fine to coarse sandy fat CLAY with fine subrounded Gravel and lignite		
958									
960					CH		Very dark gray (10 YR 3/1) fat CLAY with little Sand and fine subrounded gravel, trace lignite		
962									
964					CH		Dark gray (10 YR 4/1) fat CLAY, trace Sand		
966									
968					CH		Gray (10 YR 6/1) fat CLAY, trace Sand		
970									
972					CH		Gray (10 YR 6/1) fat CLAY, trace Sand		
974									
976					CH		Gray (10 YR 6/1) fat CLAY, trace Sand		
978									

(Continued Next Page)

DEPTH (ft)	Gamma Ray	PID (ppm)	TCE (ug/L)	PCE (ug/L)	Formation	USCS	GRAPHIC LOG	MATERIAL DESCRIPTION	
980	30 60 90				Raritan			Dark gray (10 YR 4/1) hard fat CLAY, trace Sand	
982				CH				Very dark gray (10 YR 3/1) hard fat CLAY, laminated	
984									
986		0.0				CH			Very dark gray (10 YR 3/1) Hard fat CLAY, laminated
988									
990									
992		0.0						Very dark gray (10 YR 3/1) Hard fat CLAY, laminated	
994						CH			
996		0.0				CH		Gray (10 YR 5/1) Hard fat CLAY, laminated	

End of boring at 997.0 ft. bgs.

# DOWN HOLE RUN



COMPANY: DELTA WELL & PUMP CO., INC.

LOCATION: NWIRP PINENECK RD

Well: VPB-152

Depth Driller:

Depth Logger:

Date: 07/01/2014

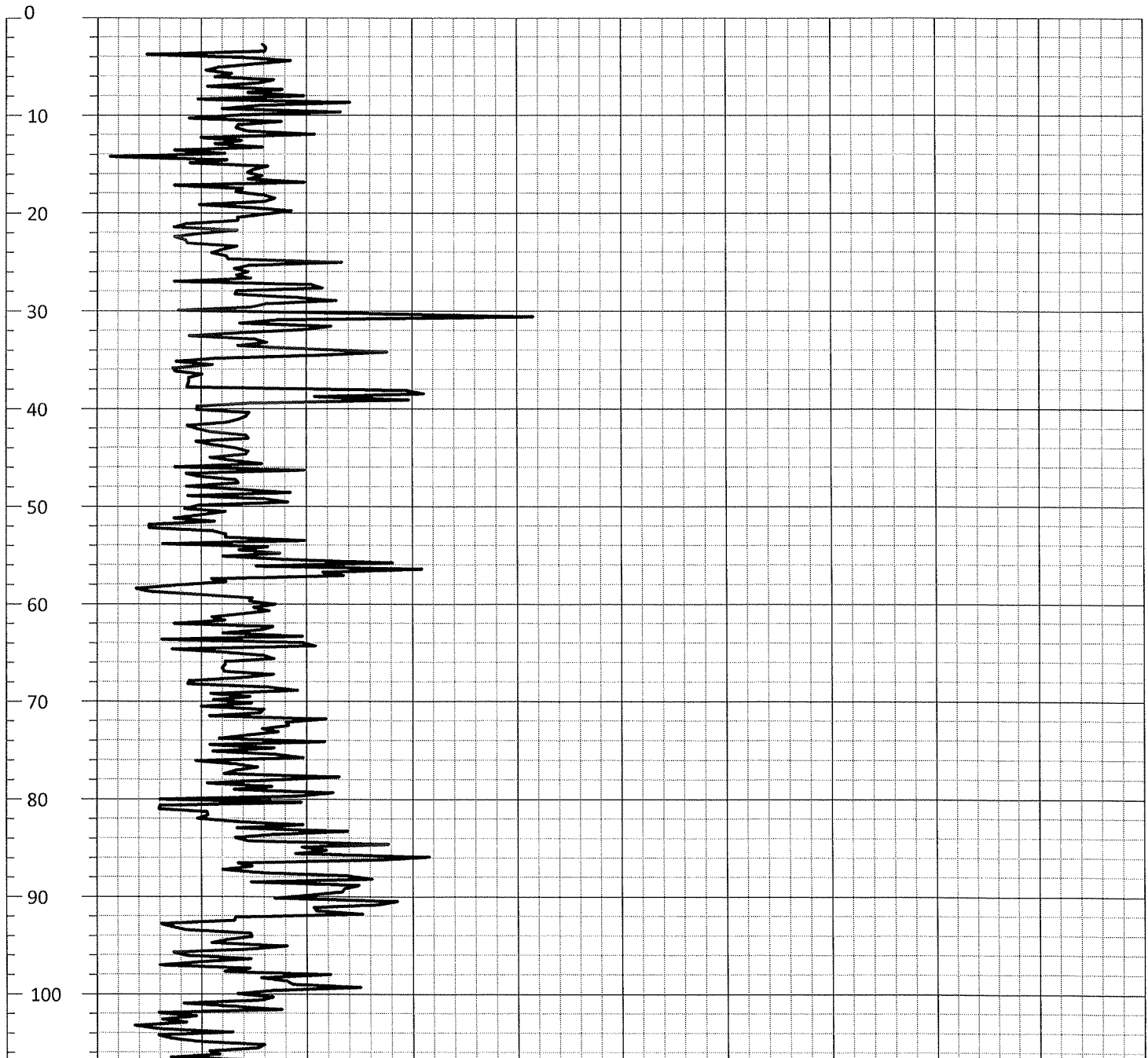
Time:

Logged by: CMO

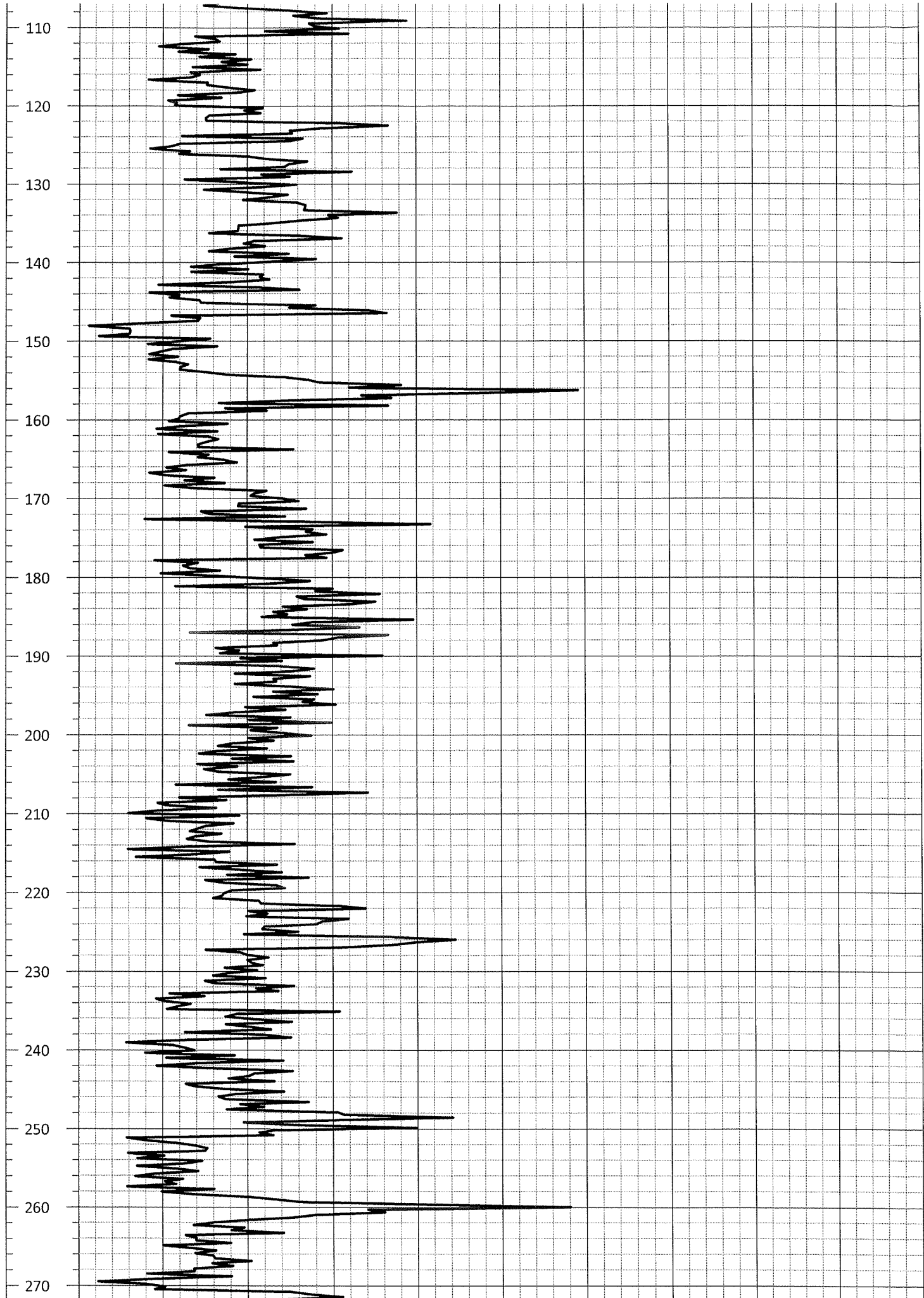
File Name: 739

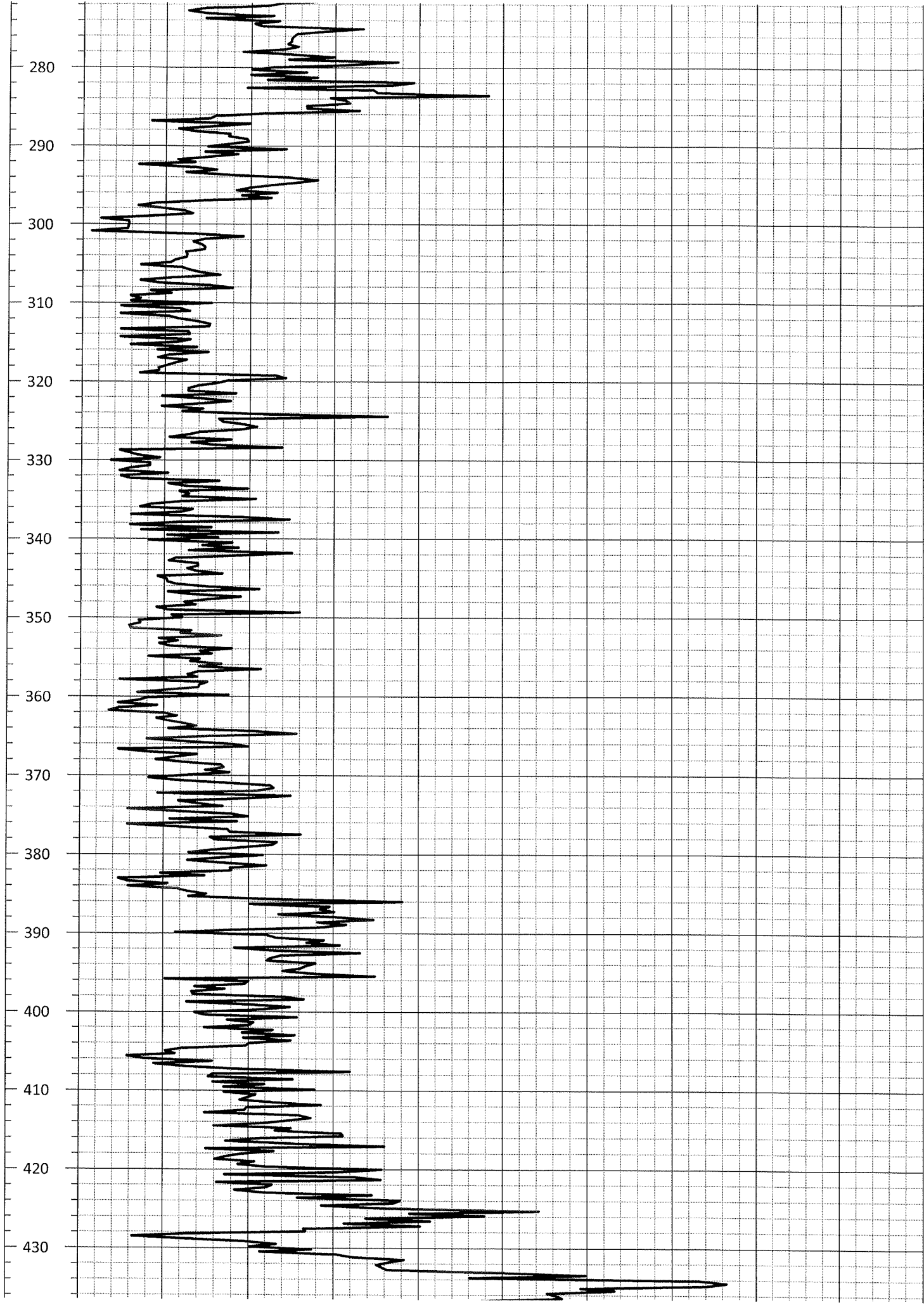
Witness: VINCE

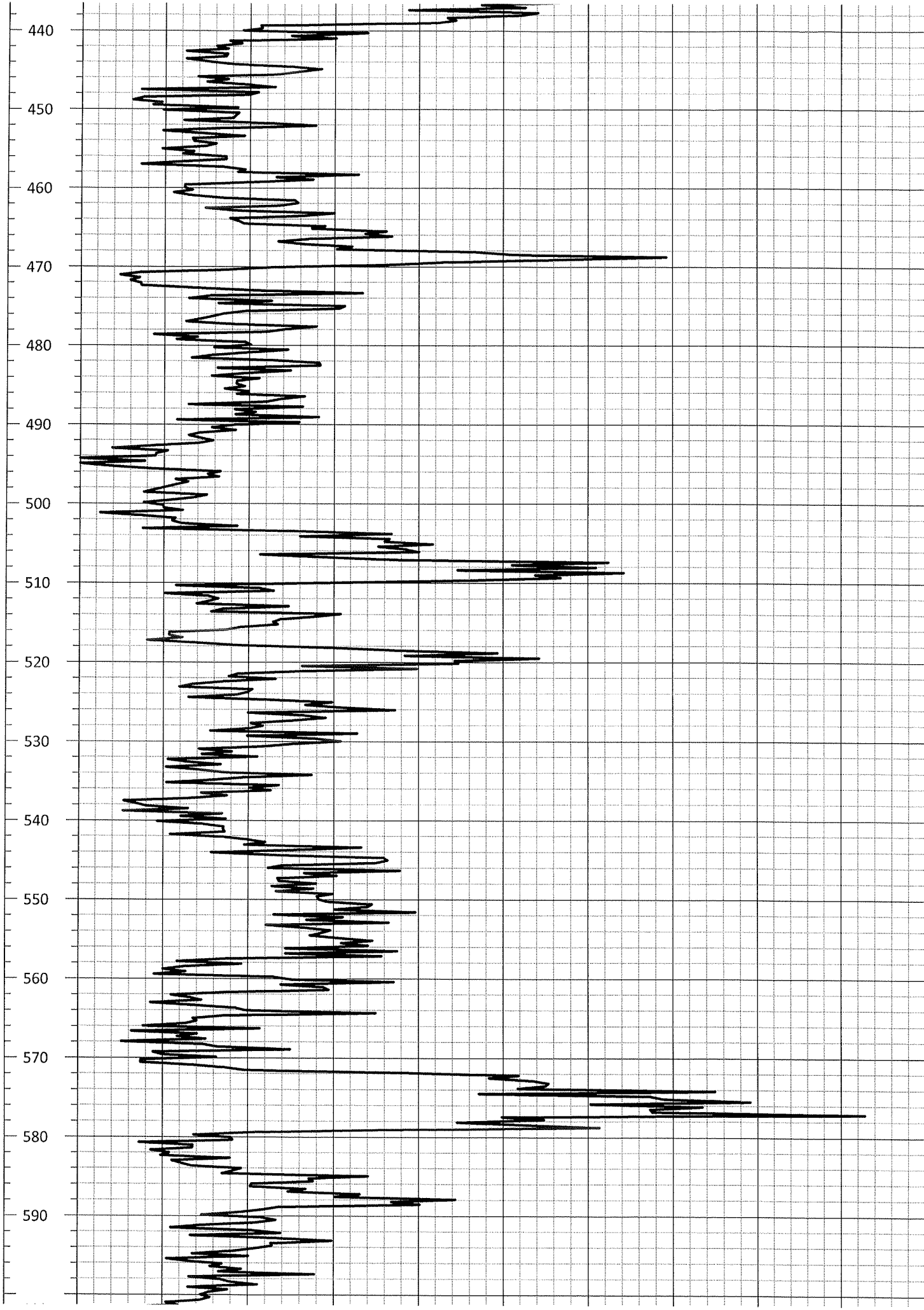
Depth (ft.)	0.0	GAMMA (cps)	100.0
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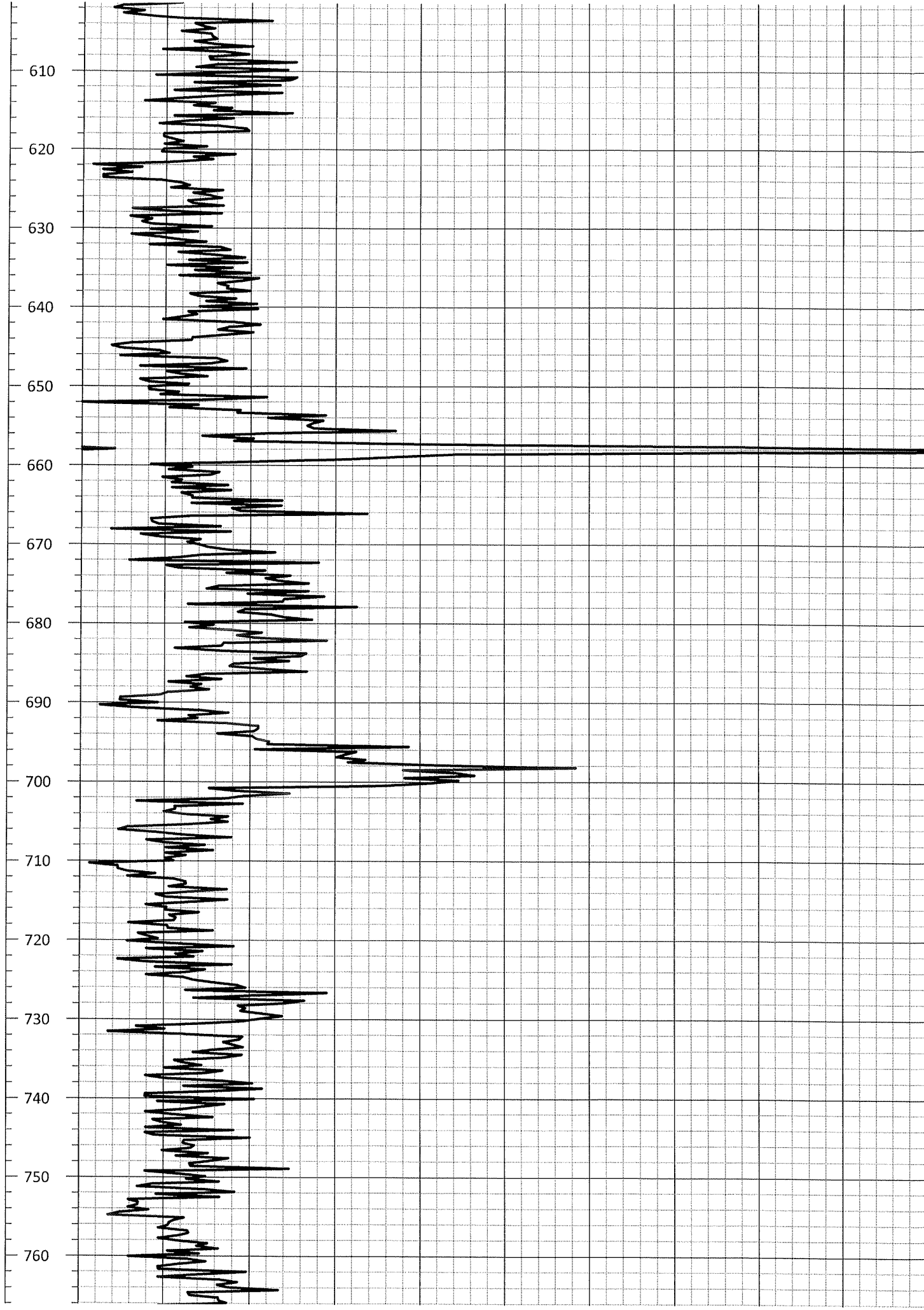


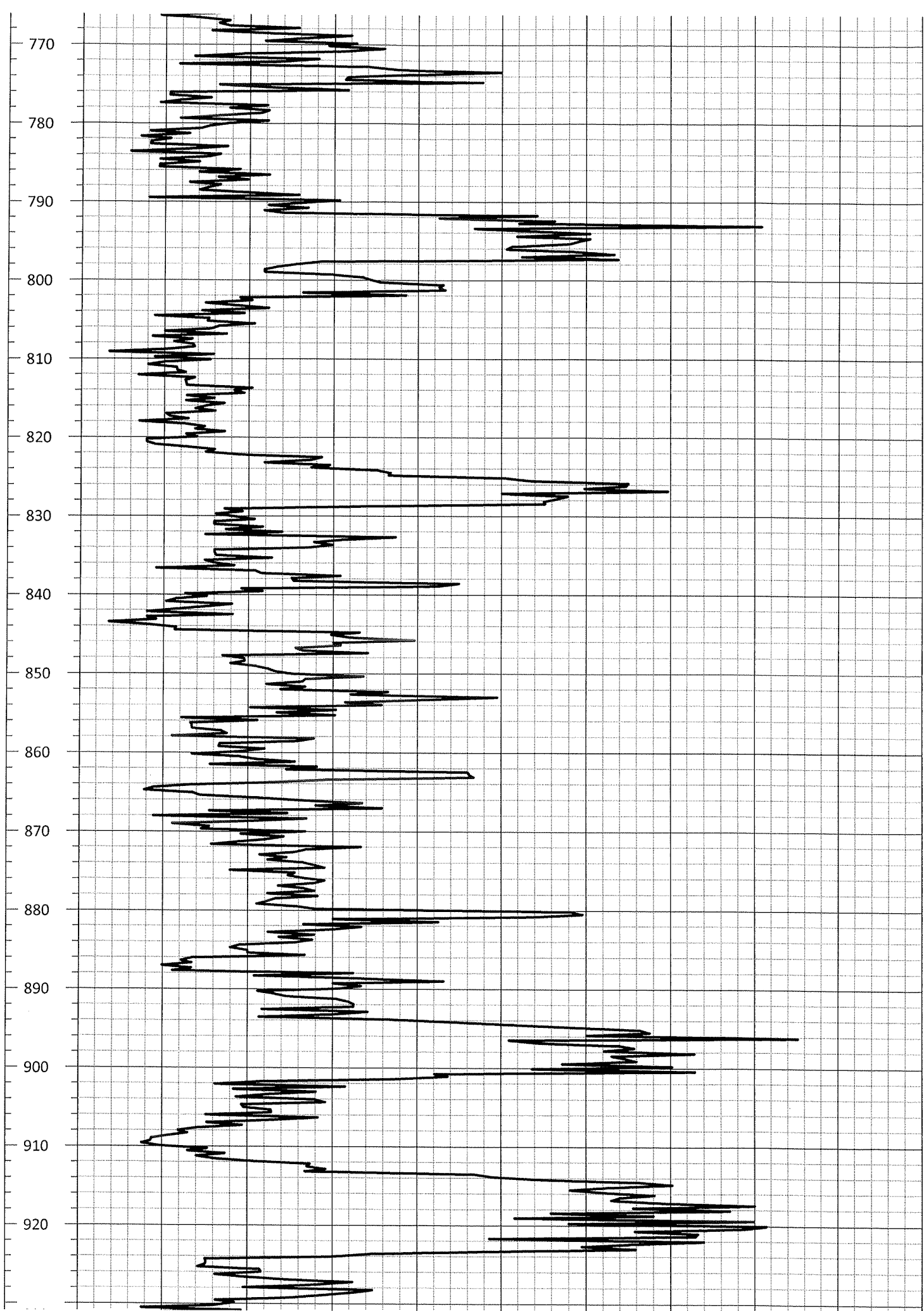


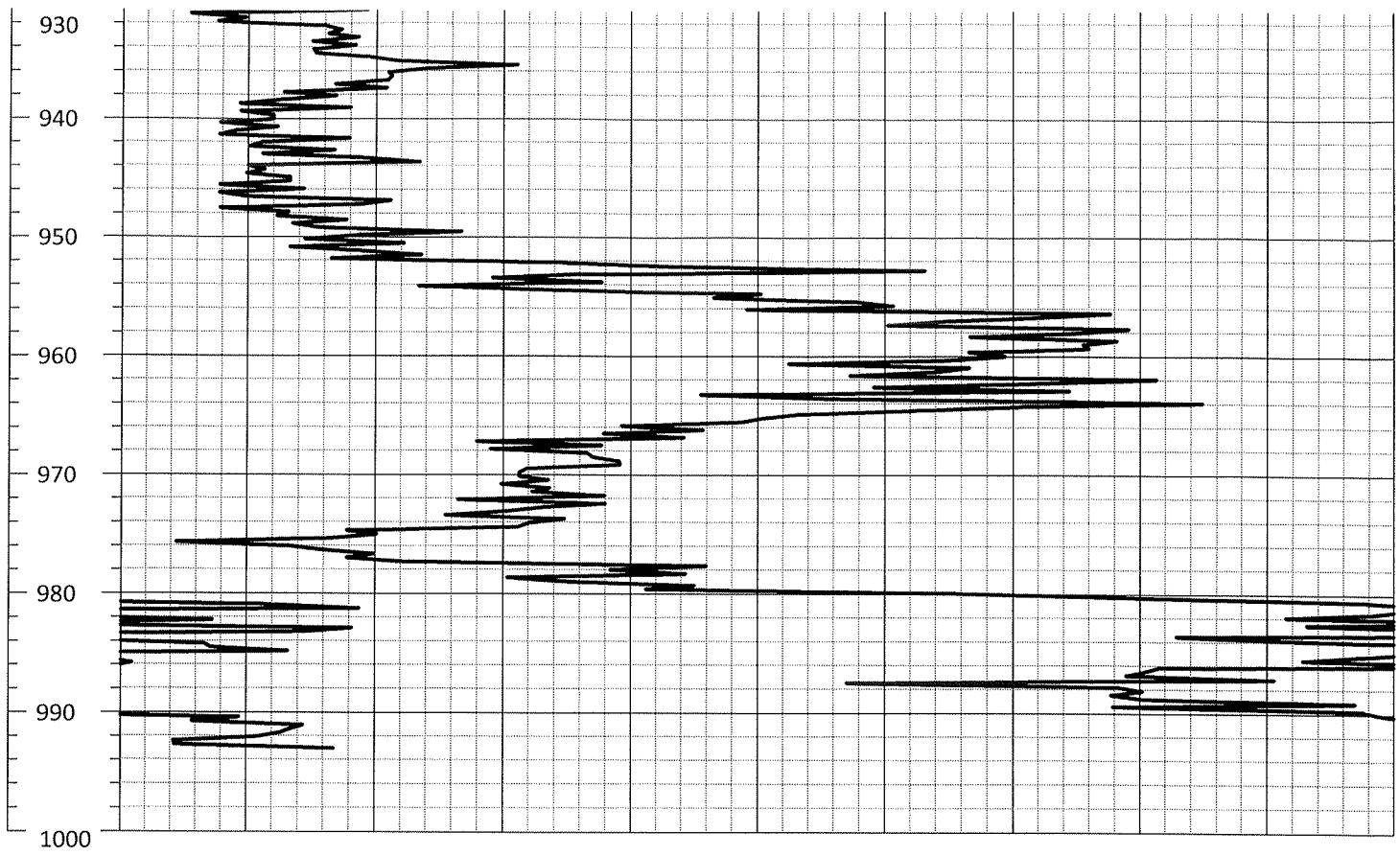












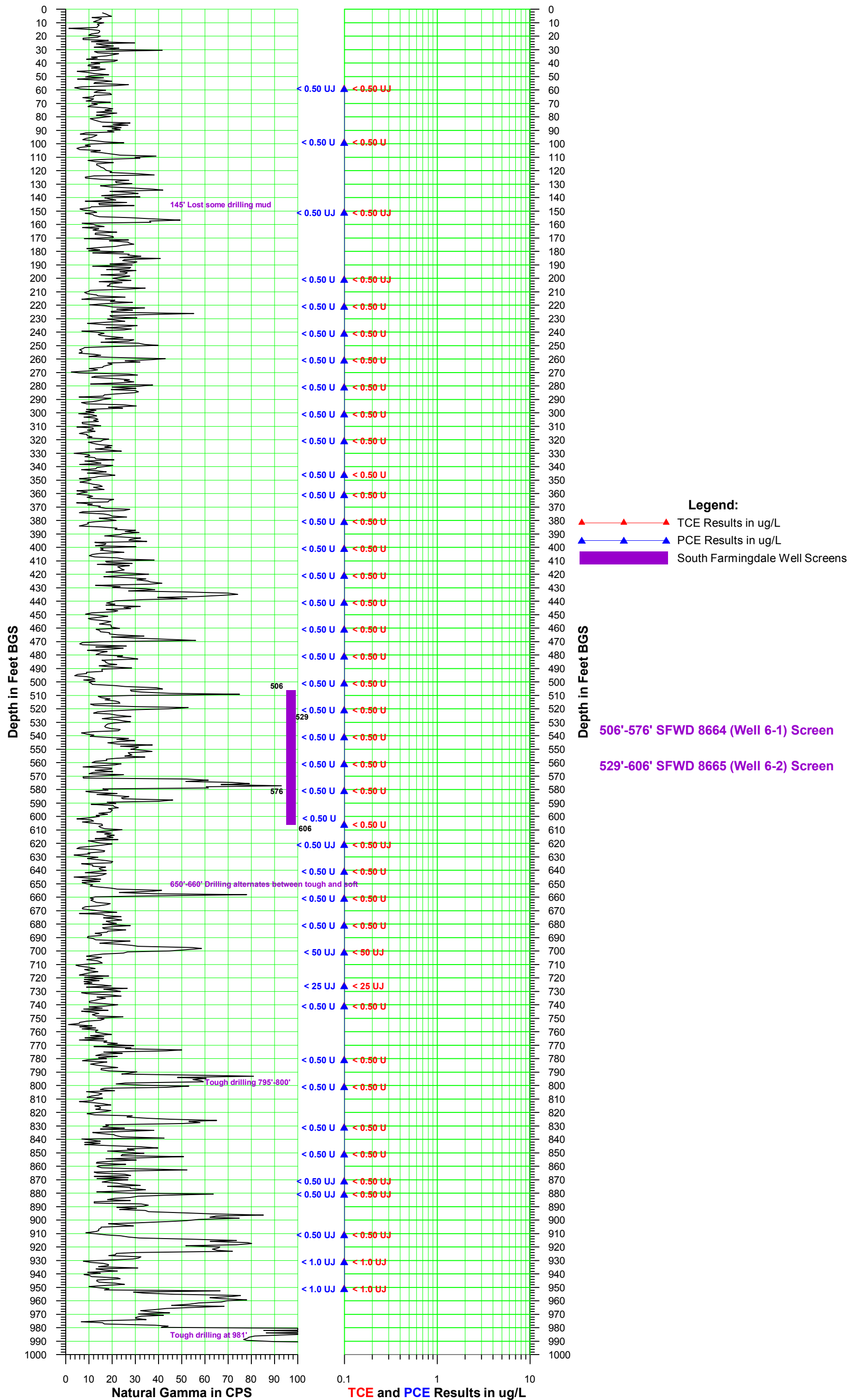
Depth (ft.)	0.0	GAMMA (cps)	100.0
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## **Section 2**

### **VPB 152 Gamma and PCE/TCE Plot**



# Vertical Profile Boring VPB-152 Downward Run - July 1, 2014 Validated Analytical Data



**Legend:**

- ▲ TCE Results in ug/L
- ▲ PCE Results in ug/L
- South Farmingdale Well Screens

506'-576' SFWD 8664 (Well 6-1) Screen

529'-606' SFWD 8665 (Well 6-2) Screen



**Section 3**

**VPB 152 Groundwater Sample Log Sheets**

## Hydropunch Sample

Client:  
Project No:  
Site Location:  
Weather Conds:

Navy - MWRP Bethpage  
06206526  
VPB - 152 Pinetree Rd

Date: 6/3-6/11/14  
VPB: 152  
Collector(s): W

Sample Date	Time	Temp (°C)	pH	Spec. Cond. (µS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Starting depth(ft)	Ending depth(ft)	Color
6/3/14	9:40	22.6	8.84	1041	6.03	65.8	71100	58	60	Brown
6/3/14	12:00	24.6	6.25	754	1.17	7.4	71400	98	100	Brown
6/3/14	1450	25.6	6.07	514	1.95	115.4	290	150	152	Slightly cloudy
6/4/14	1225	22.4	7.05	188	1.96	56.7	295	200	202	Slightly cloudy
6/4/14	1440	Not enough recovery for water parameters					Parameters	220	222	Slightly cloudy
6/5/14	940	18.1	5.84	355.6	3.56	119.4	266	240	242	Slightly cloudy
6/5/14	1145	Not enough recovery for water parameters					Parameters	240	<del>242</del> 242	Very dark brown
6/5/14	1355	18.5	5.60	290.7	3.09	125.2	541	280	285	cloudy
6/5/14	1600	20.8	5.72	369.5	3.99	74.5	246.5	300	302	Slightly cloudy
6/6/14	830	18.1	6.14	336	2.71	90.9	387.8	320	322	cloudy
6/6/14	1355	19.5	5.63	144.9	2.11	81.9	228.6	345	347	Slightly cloudy
6/9/14	940	19.1	5.81	240.3	2.44	84.7	352.1	360	362	cloudy
6/9/14	1135	16.4	6.21	356.2	1.78	93.7	236.8	380	382	Slightly cloudy
6/9/14	1340	16.2	5.47	333.5	1.70	78.5	104.1	400	402	Very slightly cloudy
6/9/14	1540	18.1	5.37	334.0	2.76	88.0	245.4	420	422	Slightly cloudy
6/10/14	940	14.7	5.53	315.9	2.68	102.9	420.4	440	442	cloudy
6/10/14	1140	19.3	3.83	413.4	4.21	263.2	112.4	460	462	Very slightly cloudy
6/10/14	1410	20.3	3.88	387.9	2.56	252.9	165	480	482	clear
6/10/14	1625	14.6	5.16	170	4.59	46.6	606.6	500	502	cloudy
6/11/14	1610	17	5.52	402.1	0.60	135.7	71100	520	522	brown
6/11/14	1235	18.4	4.63	106.3	2.51	17.5	71100	540	542	brown
			5.94							





## **Section 4**

### **VPB 152 Analytical Data Validation**

- Analytical Data Sheets
- Chain of Custody Records
- Validation Letter and Table



## Data Validation Report

Project:	Regional Groundwater Investigation - NWIRP Bethpage	
Laboratory:	Katahdin Analytical Services, Scarborough, Maine	
Service Request:	SH4831	
Analyses/Method:	EPA SW-846 Method 8260B for VOCs (GC/MS)	
Validation Level:	Limited	
RESCON Project Number:	60266526.SA.DV	
Prepared by:	Sheena Blair/RESCON	Completed on: 9/2/2014
Reviewed by:	Lori Herberich/RESCON	File Name: SH4831_8260B

### SUMMARY

The samples listed below were collected by Resolution Consultants (RESCON) from the Regional Groundwater Investigation - NWIRP Bethpage site on June 27 and 30, 2014.

Sample ID	Matrix/Sample Type
VPB152-GW-062714-950-952	Groundwater
VPB152-TRIP BLANK-063014	Trip Blank

Data validation activities were conducted with reference to *Test Methods for Evaluating Solid Waste, Physical/Chemical Methods SW846, specifically SW-846 Method 8260B, Volatile Organic Compounds by Gas Chromatography/Mass Spectrometry* (USEPA, 1996), *USEPA Contract Laboratory Program National Functional Guidelines for Superfund Organic Methods Data Review* (June 2008), and *Quality Systems Manual (QSM) for Environmental Laboratories, Version 4.2* (DoD, 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 review elements (where applicable to the method):

- X Data completeness (chain-of-custody [COC])/sample integrity
- ✓ Holding times and sample preservation
- ✓ GC/MS performance checks
- X Initial calibration/continuing calibration verification
- ✓ Laboratory blanks/equipment blanks/trip blanks
- X Surrogate spike recoveries
- NA Matrix spike (MS) and/or matrix spike duplicate (MSD) results
- ✓ Laboratory control sample (LCS) results
- NA Field duplicate results
- ✓ Internal standard results
- ✓ 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. Selected data points were estimated due to nonconformances of certain QC criteria (see discussion below). Qualified sample results are presented in Table 1.

## **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 "VPB" prefix and the "GW" from the sample ID, and truncated the ID for the Trip Blank in the report. The submitted EDD file reflects the full sample ID.

Sample VPB152-GW-062714-950-952 was mostly soil and had very little standing water. The laboratory decanted the water from three individual vials into one vial as a composite. Due to a limited sample volume the composite was analyzed at a 2-fold dilution. Positive and nondetect results for this sample were qualified as estimated (J and UJ, respectively), due to possible loss of sample integrity during the decanting procedure. Qualified sample results are presented in Table 1.

### **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.

### **GC/MS Performance Checks**

The data were reviewed to ensure that the 4-bromofluorobenzene (BFB) tuning was performed at the correct frequency and that the method acceptance criteria were met. 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:

- the initial calibration (ICAL) percent relative standard deviation (%RSD), correlation coefficient (r)/coefficient of determination ( $r^2$ ), and/or response factor method acceptance criteria were met;

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

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

**ICV Recovery Nonconformances:**

Nonconformance	Actions	
	Detected Compounds	Nondetected Compounds
%R > 120%	J	No qualification
20% < %R < 80%	J	UJ
%R < 20% (see note)	J	R*

Notes: Based on NFG 2008 VOC guidance, professional judgment is used to reject (R) nondetects in all associated samples for any analyte with < 20% recovery. Also, professional judgment is used to estimate (UJ) rather the reject (R) sample results previously negated (U) on the basis of blank contamination.

**CCV Linearity Nonconformances:**

Nonconformance	Actions	
	Detected Results	Nondetected Results
%D > 20%	J	UJ
%Drift > 20%	J*	UJ*
* No guidance in NFG, thus professional judgment was used		

Qualified sample results are shown in Table 1. Nonconformances are summarized in Attachment A in Tables A-1a and A-1b.

**Laboratory Blanks/Equipment Blanks/Trip Blanks**

Laboratory method blanks, equipment rinsate and trip 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, equipment rinsate and trip 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.

**Surrogate Spike Recoveries**

The surrogate recoveries (%Rs) were reviewed for conformance with the QC acceptance criteria.

Data qualification on the basis of surrogate recovery nonconformances was as follows:

Nonconformance	Action	
	Detected Compounds	Nondetected Compounds
%R > Upper Limit (UL)	J	No qualification
20% < %R < Lower Limit (LL)	J	UJ
%R < 20%	J	R

Nonconformances are summarized in Attachment A in Table A-2. Qualified sample results are shown in Table 1.

### **MS/MSD Results**

MS/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.

### **Internal Standard Results**

The internal standard (IS) recoveries were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

### **Sample Results/Reporting Issues**

Compounds that were not detected in the sample are reported as undetected (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**

Sample results qualified as a result of validation actions are summarized in Table 1. All actions are described above.

### **ATTACHMENTS**



Attachment A: Nonconformance Summary Tables

Attachment B: Qualifier Codes and Explanations

Attachment C: Reason Codes and Explanations

Table 1 - Data Validation Summary of Qualified Data

Sample ID	Matrix	Compound	Result	LOD	Units	Validation Qualifiers	Validation Reason
VPB152-GW-062714-950-952	WG	1,1,1-TRICHLOROETHANE		1.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	1,1,2,2-TETRACHLOROETHANE		1.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE		1.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	1,1,2-TRICHLOROETHANE		1.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	1,1-DICHLOROETHANE		1.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	1,1-DICHLOROETHENE		1.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	1,2,4-TRICHLOROBENZENE		1.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	1,2-DIBROMO-3-CHLOROPROPANE		1.5	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	1,2-DIBROMOETHANE		1.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	1,2-DICHLOROBENZENE		1.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	1,2-DICHLOROETHANE		1.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	1,2-DICHLOROETHENE, TOTAL		2.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	1,2-DICHLOROPROPANE		1.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	1,3-DICHLOROBENZENE		1.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	1,4-DICHLOROBENZENE		1.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	2-BUTANONE		5.0	UG/L	UJ	c,mc
VPB152-GW-062714-950-952	WG	2-HEXANONE		5.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	4-METHYL-2-PENTANONE		5.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	ACETONE	46	5.0	UG/L	J	c,mc,s
VPB152-GW-062714-950-952	WG	BENZENE		1.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	BROMODICHLOROMETHANE		1.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	BROMOFORM		1.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	BROMOMETHANE		2.0	UG/L	UJ	c,mc
VPB152-GW-062714-950-952	WG	CARBON DISULFIDE		1.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	CARBON TETRACHLORIDE		1.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	CHLOROBENZENE		1.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	CHLOROETHANE		2.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	CHLOROFORM		1.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	CHLOROMETHANE		2.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	CIS-1,2-DICHLOROETHENE		1.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	CIS-1,3-DICHLOROPROPENE		1.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	CYCLOHEXANE		1.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	DIBROMOCHLOROMETHANE		1.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	DICHLORODIFLUOROMETHANE		2.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	ETHYLBENZENE		1.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	ISOPROPYLBENZENE		1.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	M- AND P-XYLENE		2.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	METHYL ACETATE		1.5	UG/L	UJ	c,mc
VPB152-GW-062714-950-952	WG	METHYL CYCLOHEXANE		1.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	METHYL TERT-BUTYL ETHER		1.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	METHYLENE CHLORIDE	6.6	5.0	UG/L	J	mc,s
VPB152-GW-062714-950-952	WG	O-XYLENE		1.0	UG/L	UJ	mc

Sample ID	Matrix	Compound	Result	LOD	Units	Validation Qualifiers	Validation Reason
VPB152-GW-062714-950-952	WG	STYRENE		1.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	TETRACHLOROETHENE		1.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	TOLUENE	5.4	1.0	UG/L	J	mc,s
VPB152-GW-062714-950-952	WG	TRANS-1,2-DICHLOROETHENE		1.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	TRANS-1,3-DICHLOROPROPENE		1.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	TRICHLOROETHENE		1.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	TRICHLOROFLUOROMETHANE		2.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	VINYL CHLORIDE		2.0	UG/L	UJ	mc
VPB152-GW-062714-950-952	WG	XYLENES, TOTAL		3.0	UG/L	UJ	mc
VPB152-TRIP BLANK-063014	WQ	2-BUTANONE	2.5	2.5	UG/L	UJ	c
VPB152-TRIP BLANK-063014	WQ	ACETONE	2.5	2.5	UG/L	UJ	c
VPB152-TRIP BLANK-063014	WQ	BROMOMETHANE	1.0	1.0	UG/L	UJ	c
VPB152-TRIP BLANK-063014	WQ	METHYL ACETATE	0.75	0.75	UG/L	UJ	c

## Attachment A

## Nonconformance Summary Tables

Table A-1 - Initial Calibration Verification

ICV	Compound	% R	Limit
WG145769-7	ACETONE	145.9	80-120%
Associated samples: VPB152-GW-062714-950-952, VPB152-TRIP BLANK-063014			

Table A-1b - Continuing Calibration Verification

CCV ID	Compound	% D	Limit
WG145865 GCMS-C	BROMOMETHANE	20.8	< 20%
	2-BUTANONE	25.3	< 20%
	METHYL ACETATE	22.3	< 20%
Associated samples: VPB152-GW-062714-950-952, VPB152-TRIP BLANK-063014			

Table A-2- Surrogates

Sample ID	Surrogate	% Recovery	Lower Limit	Upper Limit
VPB152-GW-062714-950-952	1,2-DICHLOROETHANE-D4	132	70	120

**Attachment B**  
**Qualifier Codes and Explanations**

<b>Qualifier</b>	<b>Explanation</b>
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.

## Attachment C

### Reason Codes and Explanations

Reason Code	Explanation
be	Equipment blank contamination
bf	Field blank contamination
bl	Laboratory blank contamination
c	Calibration issue
co	Analyte carryover
d	Reporting limit raised due to chromatographic interference
fd	Field duplicate RPDs
h	Holding times
i	Internal standard areas
k	Estimated Maximum Possible Concentration (EMPC)
l	LCS or OPR recoveries
lc	Labeled compound recovery
ld	Laboratory duplicate RPDs
lp	Laboratory control sample/laboratory control sample duplicate RPDs
m	Matrix spike recovery
md	Matrix spike/matrix spike duplicate RPDs
nb	Negative laboratory blank contamination
p	Chemical preservation issue
r	Dual column RPD
q	Quantitation issue
s	Surrogate recovery
su	Ion suppression
t	Temperature preservation issue
x	Percent solids
y	Serial dilution results
z	ICS results

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4831-1DL  
**Client ID:** 152-062714-950-952  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4831  
**Lab File ID:** D9287.D

**Sample Date:** 27-JUN-14  
**Received Date:** 01-JUL-14  
**Extract Date:** 02-JUL-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG145865

**Analysis Date:** 02-JUL-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 03-JUL-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	2.0	ug/L	2	2	4.0	0.48	2.0
Chloromethane	U	2.0	ug/L	2	2	4.0	0.72	2.0
Vinyl Chloride	U	2.0	ug/L	2	2	4.0	0.50	2.0
Bromomethane	U	2.0	ug/L	2	2	4.0	0.98	2.0
Chloroethane	U	2.0	ug/L	2	2	4.0	1.1	2.0
Trichlorofluoromethane	U	2.0	ug/L	2	2	4.0	0.48	2.0
1,1-Dichloroethene	U	1.0	ug/L	2	1	2.0	0.70	1.0
Carbon Disulfide	U	1.0	ug/L	2	1	2.0	0.50	1.0
Freon-113	U	1.0	ug/L	2	1	2.0	0.62	1.0
<b>Methylene Chloride</b>	J	6.6	ug/L	2	5	10.	2.3	5.0
<b>Acetone</b>	J	46	ug/L	2	5	10.	4.4	5.0
trans-1,2-Dichloroethene	U	1.0	ug/L	2	1	2.0	0.50	1.0
Methyl tert-butyl Ether	U	1.0	ug/L	2	1	2.0	0.72	1.0
1,1-Dichloroethane	U	1.0	ug/L	2	1	2.0	0.42	1.0
cis-1,2-Dichloroethene	U	1.0	ug/L	2	1	2.0	0.42	1.0
Chloroform	U	1.0	ug/L	2	1	2.0	0.64	1.0
1,1,1-Trichloroethane	U	1.0	ug/L	2	1	2.0	0.40	1.0
2-Butanone	U	5.0	ug/L	2	5	10.	2.6	5.0
Cyclohexane	U	1.0	ug/L	2	1	2.0	0.62	1.0
Carbon Tetrachloride	U	1.0	ug/L	2	1	2.0	0.44	1.0
Benzene	U	1.0	ug/L	2	1	2.0	0.52	1.0
1,2-Dichloroethane	U	1.0	ug/L	2	1	2.0	0.40	1.0
Trichloroethene	U	1.0	ug/L	2	1	2.0	0.56	1.0
1,2-Dichloropropane	U	1.0	ug/L	2	1	2.0	0.50	1.0
Bromodichloromethane	U	1.0	ug/L	2	1	2.0	0.66	1.0
cis-1,3-Dichloropropene	U	1.0	ug/L	2	1	2.0	0.38	1.0
<b>Toluene</b>	J	5.4	ug/L	2	1	2.0	0.54	1.0
4-Methyl-2-Pentanone	U	5.0	ug/L	2	5	10.	2.6	5.0
trans-1,3-Dichloropropene	U	1.0	ug/L	2	1	2.0	0.40	1.0
1,1,2-Trichloroethane	U	1.0	ug/L	2	1	2.0	0.66	1.0
Tetrachloroethene	U	1.0	ug/L	2	1	2.0	0.80	1.0
Dibromochloromethane	U	1.0	ug/L	2	1	2.0	0.60	1.0
2-Hexanone	U	5.0	ug/L	2	5	10.	3.4	5.0
Chlorobenzene	U	1.0	ug/L	2	1	2.0	0.44	1.0
Ethylbenzene	U	1.0	ug/L	2	1	2.0	0.42	1.0



**Report of Analytical Results**

**Client:** ENSAFE  
**Lab ID:** SH4831-1DL  
**Client ID:** 152-062714-950-952  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4831  
**Lab File ID:** D9287.D

**Sample Date:** 27-JUN-14  
**Received Date:** 01-JUL-14  
**Extract Date:** 02-JUL-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG145865

**Analysis Date:** 02-JUL-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 03-JUL-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U <i>0.5</i>	3.0	ug/L	2	3	6.0	0.50	3.0
Styrene	U	1.0	ug/L	2	1	2.0	0.46	1.0
Bromoform	U	1.0	ug/L	2	1	2.0	0.46	1.0
Isopropylbenzene	U	1.0	ug/L	2	1	2.0	0.46	1.0
1,1,2,2-Tetrachloroethane	U	1.0	ug/L	2	1	2.0	0.76	1.0
1,3-Dichlorobenzene	U	1.0	ug/L	2	1	2.0	0.52	1.0
1,4-Dichlorobenzene	U	1.0	ug/L	2	1	2.0	0.48	1.0
1,2-Dichlorobenzene	U	1.0	ug/L	2	1	2.0	0.30	1.0
1,2,4-Trichlorobenzene	U	1.0	ug/L	2	1	2.0	0.74	1.0
Methyl Acetate	U	1.5	ug/L	2	1	2.0	1.1	1.5
Methylcyclohexane	U	1.0	ug/L	2	1	2.0	0.60	1.0
o-Xylene	U	1.0	ug/L	2	1	2.0	0.50	1.0
M+P-Xylenes	U	2.0	ug/L	2	2	4.0	1.2	2.0
1,2-Dichloroethylene (Total)	U	2.0	ug/L	2	2	4.0	0.42	2.0
1,2-Dibromoethane	U	1.0	ug/L	2	1	2.0	0.44	1.0
1,2-Dibromo-3-Chloropropane	U	1.5	ug/L	2	1	2.0	1.0	1.5
P-Bromofluorobenzene		82.5	%					
Toluene-d8		112.	%					
1,2-Dichloroethane-d4	*	132.	%					
Dibromofluoromethane		112.	%					



## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4831-2  
**Client ID:** VPB152-TB-063014  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4831  
**Lab File ID:** D9284.D

**Sample Date:** 30-JUN-14  
**Received Date:** 01-JUL-14  
**Extract Date:** 02-JUL-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG145865

**Analysis Date:** 02-JUL-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 03-JUL-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	<del>U</del> UJ	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
Acetone	<del>U</del> UJ	2.5	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	<del>U</del> UJ	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50

*PC 7/16/14*

### Report of Analytical Results

Client: ENSAFE  
 Lab ID: SH4831-2  
 Client ID: VPB152-TB-063014  
 Project: Navy Clean WE15-03-06 NW  
 SDG: SH4831  
 Lab File ID: D9284.D

Sample Date: 30-JUN-14  
 Received Date: 01-JUL-14  
 Extract Date: 02-JUL-14  
 Extracted By: DJP  
 Extraction Method: SW846 5030  
 Lab Prep Batch: WG145865

Analysis Date: 02-JUL-14  
 Analyst: DJP  
 Analysis Method: SW846 8260B  
 Matrix: AQ  
 % Solids: NA  
 Report Date: 03-JUL-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	<del>U</del> UJ	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		88.0	%					
Toluene-d8		119.	%					
1,2-Dichloroethane-d4	*	135.	%					
Dibromofluoromethane	*	118.	%					

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## Data Validation Report

Project: Regional Groundwater Investigation - NWIRP Bethpage  
Laboratory: Test-America, South Burlington, Vermont  
Service Request: 200-22958  
Analyses/Method: EPA Method TO-15, VOCs Collected in Canisters - GC/MS  
Validation Level: Limited  
RESCON Project 60266526.SA.DV  
Number:  
Prepared by: Sheena Blair/RESCON Completed on: 09/14/2014  
Reviewed by: Lori Herberich/RESCON File Name: 200-22958\_TO-15

### SUMMARY

The sample listed below was collected by Resolution Consultants from the Regional Groundwater Investigation - NWIRP Bethpage site on June 24, 2014.

Sample ID	Matrix/Sample Type
VPB152-AIR-062414	Ambient Air

Data validation activities were conducted with reference to *Determination Of Volatile Organic Compounds (VOCs) In Air Collected In Specially-Prepared Canisters And Analyzed By Gas Chromatography/Mass Spectrometry (GC/MS)* (USEPA, Method TO-15) and the *USEPA Contract Laboratory Program National Functional Guidelines for Superfund Organic Methods Data Review* (June 2008). 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
- ✓ GC/MS performance checks
- ✓ Initial calibration/continuing calibration verification
- ✓ Laboratory blanks
- NA Matrix duplicate (MD) results
- ✓ Laboratory control sample (LCS) results
- NA Field duplicates
- ✓ Internal standards
- ✓ Sample results/reporting issues

The symbol (✓) indicates that no validation qualifiers were applied based on this parameter. NA indicates that the parameter was not included as part of this data set or was not applicable to this

validation and therefore not reviewed. The symbol (X) 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.

### **Holding Times and Sample Preservation**

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

### **GC/MS Performance Checks**

The data were reviewed to ensure that the 4-bromofluorobenzene (BFB) tuning was performed at the correct frequency and that the method acceptance criteria were met. 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:

- the initial calibration (ICAL) percent relative standard deviation (%RSD), correlation coefficient (r)/coefficient of determination ( $r^2$ ), and/or response factor method acceptance criteria were met;
- the continuing calibration verification standard (CCV) method percent difference or percent drift (%Ds) and RF acceptance criteria were met; and
- the retention time method acceptance criteria were met.

The QC acceptance criteria were met.

### **Laboratory Blanks**

Laboratory method blanks were evaluated as to whether there were contaminants detected above the detection limit (DL). Blank results were reviewed for conformance with the QC acceptance criteria. Data validation qualifications for individual samples are based on the maximum contaminant concentration detected in all associated blanks.

The QC acceptance criteria were met; qualification of the sample results was not required.

### **MS/MSD Results**

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

### **LCS Results**

The LCS recoveries 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.

### **Internal Standard Results**

The internal standard (IS) recoveries were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

### **Sample Results/Reporting Issues**

Compounds that were not detected in the sample are reported as undetected (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: Nonconformance Summary Tables

Attachment B: Qualifier Codes and Explanations

**Attachment A**

**Nonconformance Summary Tables**

No nonconformances were identified during this review.

**Attachment B****Qualifier Codes and Explanations**

<b>Qualifier</b>	<b>Explanation</b>
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.

**TestAmerica Burlington**  
 30 Community Drive  
 Suite 11

South Burlington, VT 05403  
 Phone 802-880-1990 fax 802-880-1919

### Canister Samples Chain of Custody Record

TestAmerica Analytical Testing Corp. assumes no liability with respect to the collection and shipment of these samples.

Client Contact Information Company: Resolution Consultants Address: 100 Red Schoolhouse Rd, City/State/Zip: Chittenden, VT 05747 Phone: 847-435-4180 FAX: Project Name: N111P Bethpage Site: 66266526 PO #			Project Manager: E. Vivandou Phone: 847-435-4180 Email: E.Vivandou@Resolution.com Site Contact: TIA Contact: Analysis Turnaround Time Standard (Specify) Push (Specify)			Samples Collected By: <u>Vivant Vivandou</u> of <u>COCs</u>													
Sample ID	Sample Description	Time Start	Time Stop	Canister Vacuum in Field, % (Bar)	Canister Vacuum in Lab, % (Bar)	Flow Controller ID	Canister ID	TO-15	MA-4PH	EPA 30C	EPA 30C	ASTM D-1946	Other (Please specify in notes section)	Sample Type	Indoor Air	Artificial Air	Soft Gas	Leak Gas	Other (Please specify in notes section)
	VPB152-AIR-062414	6/24/14 800	1555	-29	-7	3695	5052	X											
Special Instructions/QC Requirements & Comments:																			
Temperature (Fahrenheit) Ambient Interior Exterior Pressure (Inches of Hg) Ambient Interior Shop																			
Date/Time: 6/24/14 / 1730 Samples Shipped by: <u>[Signature]</u> Samples Returned by: Date/Time: Date/Time:																			
Date/Time: 6/24/14 / 1730 Samples Received by: <u>LABVZ 62514 1030</u> Received by: Received by:																			
Lab Use Only Shipper Name: Operated by: Conditions:																			



200-22958 COC

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Client: Katahdin Analytical Services

Job Number: 200-22958-1

Sdg Number: 200-22958

Client Sample ID: VPB152-AIR-062414

Lab Sample ID: 200-22958-1

Date Sampled: 06/24/2014 1555

Client Matrix: Air

Date Received: 06/25/2014 1030

## TO-15 Volatile Organic Compounds in Ambient Air

Analysis Method:	TO-15	Analysis Batch:	200-74232	Instrument ID:	CHW.i
Prep Method:	Summa Canister	Prep Batch:	N/A	Lab File ID:	8298_026.d
Dilution:	1.0			Initial Weight/Volume:	200 mL
Analysis Date:	06/27/2014 0919			Final Weight/Volume:	200 mL
Prep Date:	06/27/2014 0919			Injection Volume:	200 mL

Analyte	Result (ppb v/v)	Qualifier	DL	LOQ
1,1,1-Trichloroethane	0.080	U	0.20	0.20
1,1,1,2-Tetrachloroethane	0.030	U	0.20	0.20
1,1,1,2-Trichloro-1,2,2-trifluoroethane	0.030	U	0.20	0.20
1,1,2-Trichloroethane	0.030	U	0.20	0.20
1,1-Dichloroethane	0.080	U	0.20	0.20
1,1-Dichloroethene	0.080	U	0.20	0.20
1,2,4-Trichlorobenzene	0.080	U	0.50	0.50
1,2-Dibromoethane (EDB)	0.080	U	0.20	0.20
1,2-Dichlorobenzene	0.030	U	0.20	0.20
1,2-Dichloroethane	0.030	U	0.20	0.20
1,2-Dichloropropane	0.080	U	0.20	0.20
Acetone	7.5		5.0	5.0
1,3-Dichlorobenzene	0.030	U	0.20	0.20
1,4-Dichlorobenzene	0.030	U	0.20	0.20
2-Butanone (MEK)	0.73		0.50	0.50
2-Hexanone	0.20	U	0.50	0.50
4-Methyl-2-pentanone	0.080	U	0.50	0.50
Benzene	0.030	U	0.20	0.20
Bromoform	0.030	U	0.20	0.20
Bromomethane	0.080	U	0.20	0.20
Carbon disulfide	10		0.50	0.50
Carbon tetrachloride	0.080	U	0.20	0.20
Chlorobenzene	0.030	U	0.20	0.20
Dibromochloromethane	0.030	U	0.20	0.20
Chloroethane	0.080	U	0.50	0.50
Chloroform	0.080	U	0.20	0.20
Chloromethane	0.50		0.50	0.50
cis-1,2-Dichloroethene	0.080	U	0.20	0.20
cis-1,3-Dichloropropene	0.080	U	0.20	0.20
Cyclohexane	0.080	U	0.20	0.20
Bromodichloromethane	0.030	U	0.20	0.20
Dichlorodifluoromethane	0.51		0.50	0.50
Ethylbenzene	0.030	U	0.20	0.20
Isopropylbenzene	0.030	U	0.20	0.20
Methyl tert-butyl ether	0.080	U	0.20	0.20
Methylene Chloride	0.20	U	0.50	0.50
m,p-Xylene	0.080	U	0.50	0.50
Xylene, o-	0.030	U	0.20	0.20
Styrene	0.030	U	0.20	0.20
Tetrachloroethene	0.030	U	0.20	0.20
Toluene	0.030	U	0.20	0.20
trans-1,2-Dichloroethene	0.080	U	0.20	0.20
trans-1,3-Dichloropropene	0.080	U	0.20	0.20
Trichloroethene	0.080	U	0.20	0.20
Trichlorofluoromethane	0.22		0.20	0.20
Vinyl chloride	0.080	U	0.20	0.20

## Analytical Data

Client: Katahdin Analytical Services

Job Number: 200-22958-1

Sdg Number: 200-22958

Client Sample ID: **VPB152-AIR-062414**

Lab Sample ID: 200-22958-1

Date Sampled: 06/24/2014 1555

Client Matrix: Air

Date Received: 06/25/2014 1030

### TO-15 Volatile Organic Compounds in Ambient Air

Analysis Method:	TO-15	Analysis Batch:	200-74232	Instrument ID:	CHW.i
Prep Method:	Summa Canister	Prep Batch:	N/A	Lab File ID:	8298_026.d
Dilution:	1.0			Initial Weight/Volume:	200 mL
Analysis Date:	06/27/2014 0919			Final Weight/Volume:	200 mL
Prep Date:	06/27/2014 0919			Injection Volume:	200 mL

Analyte	Result (ppb v/v)	Qualifier	DL	LOQ
Xylene (total)	0.080	U	0.20	0.20

Analyte	Result (ug/m3)	Qualifier	DL	LOQ
1,1,1-Trichloroethane	0.44	U	1.1	1.1
1,1,2,2-Tetrachloroethane	0.21	U	1.4	1.4
1,1,2-Trichloro-1,2,2-trifluoroethane	0.23	U	1.5	1.5
1,1,2-Trichloroethane	0.16	U	1.1	1.1
1,1-Dichloroethane	0.32	U	0.81	0.81
1,1-Dichloroethene	0.32	U	0.79	0.79
1,2,4-Trichlorobenzene	0.59	U	3.7	3.7
1,2-Dibromoethane (EDB)	0.61	U	1.5	1.5
1,2-Dichlorobenzene	0.18	U	1.2	1.2
1,2-Dichloroethane	0.12	U	0.81	0.81
1,2-Dichloropropane	0.37	U	0.92	0.92
Acetone	18		12	12
1,3-Dichlorobenzene	0.18	U	1.2	1.2
1,4-Dichlorobenzene	0.18	U	1.2	1.2
2-Butanone (MEK)	2.1		1.5	1.5
2-Hexanone	0.82	U	2.0	2.0
4-Methyl-2-pentanone	0.33	U	2.0	2.0
Benzene	0.096	U	0.64	0.64
Bromoform	0.31	U	2.1	2.1
Bromomethane	0.31	U	0.78	0.78
Carbon disulfide	32		1.6	1.6
Carbon tetrachloride	0.50	U	1.3	1.3
Chlorobenzene	0.14	U	0.92	0.92
Dibromochloromethane	0.26	U	1.7	1.7
Chloroethane	0.21	U	1.3	1.3
Chloroform	0.39	U	0.98	0.98
Chloromethane	1.0		1.0	1.0
cis-1,2-Dichloroethene	0.32	U	0.79	0.79
cis-1,3-Dichloropropene	0.36	U	0.91	0.91
Cyclohexane	0.28	U	0.69	0.69
Bromodichloromethane	0.20	U	1.3	1.3
Dichlorodifluoromethane	2.5		2.5	2.5
Ethylbenzene	0.13	U	0.87	0.87
Isopropylbenzene	0.15	U	0.98	0.98
Methyl tert-butyl ether	0.29	U	0.72	0.72
Methylene Chloride	0.69	U	1.7	1.7
m,p-Xylene	0.35	U	2.2	2.2
Xylene, o-	0.13	U	0.87	0.87
Styrene	0.13	U	0.85	0.85
Tetrachloroethene	0.20	U	1.4	1.4
Toluene	0.11	U	0.75	0.75
trans-1,2-Dichloroethene	0.32	U	0.79	0.79
trans-1,3-Dichloropropene	0.36	U	0.91	0.91

# Analytical Data

Client: Katahdin Analytical Services

Job Number: 200-22958-1  
Sdg Number: 200-22958

Client Sample ID: VPB152-AIR-062414

Lab Sample ID: 200-22958-1  
Client Matrix: Air

Date Sampled: 06/24/2014 1555  
Date Received: 06/25/2014 1030

## TO-15 Volatile Organic Compounds in Ambient Air

Analysis Method:	TO-15	Analysis Batch:	200-74232	Instrument ID:	CHW.i
Prep Method:	Summa Canister	Prep Batch:	N/A	Lab File ID:	8298_026.d
Dilution:	1.0			Initial Weight/Volume:	200 mL
Analysis Date:	06/27/2014 0919			Final Weight/Volume:	200 mL
Prep Date:	06/27/2014 0919			Injection Volume:	200 mL

Analyte	Result (ug/m3)	Qualifier	DL	LOQ
Trichloroethene	0.43	U	1.1	1.1
Trichlorofluoromethane	1.2		1.1	1.1
Vinyl chloride	0.20	U	0.51	0.51
Xylene (total)	0.35	U	0.87	0.87



## Data Validation Report

Project: Regional Groundwater Investigation - NWIRP Bethpage

Laboratory: Katahdin Analytical Services, Scarborough, Maine

Service Request: SH4793

Analyses/Method: EPA SW-846 Method 8260B for VOCs (GC/MS) and Standard Method 5310B for Total Organic Carbon by High-Temperature Combustion

Validation Level: Limited

RESCON Project 60266526.SA.DV  
Number:

Prepared by: Sheena Blair/RESCON Completed on: 10/16/2014

Reviewed by: Lori Herberich/RESCON File Name: SH4793\_5310B and 8260B

### SUMMARY

The samples listed below were collected by Resolution Consultants (RESCON) from the Regional Groundwater Investigation - NWIRP Bethpage site on June 24, 25, and 26, 2014.

Sample ID	Matrix/Sample Type
VPB152-EB-062514	Equipment blank
VPB152-GW-062414-870-872	Groundwater
VPB152-GW-062414-880-882	Groundwater
VPB152-GW-062514-910-912	Groundwater
VPB152-GW-062614-930-932	Groundwater
VPB152-TRIPBLANK-062614	Trip Blank

The samples were analyzed in accordance with:

- *Test Methods for Evaluating Solid Waste, Physical/Chemical Methods SW-846, Method 8260B, Volatile Organic Compounds by Gas Chromatography/Mass Spectrometry (USEPA, 1996).*
- *Standard Methods for the Examination of Water and Wastewater, Method SM310B, Total Organic Carbon by High-Temperature Combustion*

Data validation activities were conducted with reference to these methods, *USEPA Contract Laboratory Program National Functional Guidelines for Superfund Organic Methods Data Review* (June 2008), *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):

- X Data completeness (chain-of-custody [COC])/sample integrity
- X Holding times and sample preservation
- ✓ GC/MS performance checks
- X Initial calibration/continuing calibration verification
- X Laboratory blanks/equipment blanks/ trip blanks
- ✓ Surrogate spike recoveries
- ✓ Matrix spike (MS) results
- ✓ Laboratory control sample (LCS) results
- NA Field duplicate results
- ✓ Internal standard results
- ✓ 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 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. Selected data points were negated and/or estimated, due to nonconformances of certain QC criteria (see discussion below). Qualified sample results are presented in Table 1.

## 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 "VPB-" prefix and the "GW" from the sample ID, and truncated the ID for the Trip Blank in the report. The submitted EDD file reflects the full sample ID.

Sample VPB152-GW-062614-930-932 was mostly soil and had very little standing water. The laboratory decanted the water from the individual vials into one vial as a composite. Due to lack of sample volume the composite was analyzed at a 2-fold dilution. Positive and nondetect results for this sample were qualified as estimated (J and UJ) respectively, due to possible loss of sample integrity during the decanting procedure.

### **Holding Times and Sample Preservation**

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

The cooler temperature upon arrival at the laboratory was 17.4°C, which grossly exceeded the QC acceptance limit of 4°C ± 2°C.

All samples were qualified as follows:

#### **Cooler Temperature (all sample matrices)**

Criteria	Actions			
	Aqueous		Solid	
	Detected	Nondetected	Detected	Nondetected
Cooler T < 2°C	No qualification <sup>1</sup>		No qualification <sup>1</sup>	
Cooler T >6°C but <10°C	No qualification <sup>1</sup>		No qualification <sup>1</sup>	
Cooler T >10°C	J <sup>2</sup>	UJ <sup>2</sup>	J <sup>2</sup>	UJ <sup>2</sup>
<sup>1</sup> No guidance listed in Region 2. NFG recommends professional judgment, thus RESCON professional judgment was used. <sup>2</sup> Per Region 2 guidance that states qualify data if cooler temperature >10°C due to no ice/ice melted upon arrival.				

Qualified sample results are shown in Table 1.

### **GC/MS Performance Checks**

The data were reviewed to ensure that the 4-bromofluorobenzene (BFB) tuning was performed at the correct frequency and that the method acceptance criteria were met. All QC acceptance criteria were met.

### **Initial Calibration/Continuing Calibration Verification**

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

- the initial calibration (ICAL) percent relative standard deviation (%RSD), correlation coefficient (r)/coefficient of determination (r<sup>2</sup>), and/or response factor method acceptance criteria were met;
- the initial calibration verification (ICV) percent recovery (%R) criteria were met;
- the continuing calibration verification standard (CCV) method percent difference or percent drift (%Ds), %Rs, and/or RF acceptance criteria were met; and/or
- the retention time method acceptance criteria were met.

Nonconformances are summarized in Attachment A in Table A-1.

Data qualification to the analytes associated with the specific ICAL and/or CCV was as follows:

**ICAL Linearity Nonconformances:**

Nonconformance	Actions	
	Detected Results	Nondetected Results
%RSD > 15% and quantitation based on mean RF	J	UJ
r or r <sup>2</sup> < 0.99 and quantitation based on linear regression	J*	UJ*
* No guidance in NFG, thus RESCON professional judgment was used		

**ICV Recovery Nonconformances:**

Nonconformance	Actions	
	Detected Compounds	Nondetected Compounds
%R > 120%	J	No qualification
20% < %R < 80%	J	UJ
%R < 20% (see note)	J	R*

Notes: Based on NFG 2008 VOC guidance, professional judgment is used to reject (R) non-detects in all associated samples for any analyte with < 20% recovery. Also, professional judgment is used to estimate (UJ) rather the reject (R) sample results previously negated (U) on the basis of blank contamination.

**CCV Linearity Nonconformances:**

Nonconformance	Actions	
	Detected Results	Nondetected Results
%D > 20%	J	UJ
%Drift > 20%	J*	UJ*
* No guidance in NFG, thus professional judgment was used		

Qualified sample results are shown in Table 1.

**Laboratory Blanks/Equipment Blanks/Trip Blanks**

Laboratory method blanks, equipment rinsate and trip 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, equipment rinsate and trip 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.

Nonconformances are summarized in Attachment A in Tables A-2 and A-3.

Sample results were qualified as follows:

For common lab contaminants (methylene chloride, acetone, 2-butanone):

Blank type	Blank result	Sample result	Action for samples
Method, Storage, Field, Trip, or Instrument*	Detects	Not detected	No qualification
	≤ 2x LOQ	< 2x LOQ	Report sample LOQ value with a U
		≥ 2x LOQ and ≤ 4x LOQ	Report the sample result with a U**
		≥ 4x LOQ	No qualifications
	> 2x LOQ	< 2x LOQ	Report sample LOQ value with a U
		≥ 2x LOQ and < blank contamination	Report the sample result with a U
≥ 2x LOQ and ≥ blank contamination		If the result is ≤ 2x blank result, report the sample result U.** If the result is > 2x blank result, no qualification is required.**	
* Qualifications based on instrument blank results affect only the sample analyzed immediately after the sample that has target compounds that exceed the calibration range or non-target compounds that exceed 100 g/L.			
**Based on RESCON professional judgment			

For TOC:

Blank Type	Blank Result	Sample Result	Action for Samples
PB / EB/ FB (Positive)	> LOQ	≥DL but ≤ LOQ	Qualify as nondetect (U) at the LOQ
		>LOQ but < 10x Blank Result	Qualify results as unusable
		≥10x Blank Result	No action
	≥DL but ≤LOQ	Nondetect	No action
		≥DL but ≤LOQ	Qualify as nondetect (U) at the LOQ
		> LOQ	Use RESCON professional judgment

LOQ - Limit of Quantitation.

### Surrogate Spike Recoveries

The surrogate recoveries (%Rs) were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

### MS Results

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

For TOC the MS/MSD %Rs and relative percent differences (RPDs) were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

### LCS Results

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

Acetone had a high %R in one of the LCSs; however, the associated sample results were nondetect. No validation action was required for the high bias.



**Field Duplicate Results**

Field duplicate RPDs were reviewed for conformance with the QC criterion of  $\leq 30\%$  for aqueous matrices. This criterion applies if both results were greater than five times the Limit of Quantitation (LOQ). All QC acceptance criteria were met.

**Internal Standard Results**

The internal standard (IS) recoveries were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

**Sample Results/Reporting Issues**

Compounds that were not detected in the sample are reported as undetected (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**

Sample results qualified as a result of validation actions are summarized in Table 1. All actions are described above.

**ATTACHMENTS**

Attachment A: Nonconformance Summary Tables

Attachment B: Qualifier Codes and Explanations

Attachment C: Reason Codes and Explanations

Sample ID	Matrix	Compound	Result	LOD	Units	Validation Qualifiers	Validation Reason
VPB152-EB-062514	WQ	TOTAL ORGANIC CARBON		1.0*	MG/L	UJ	t,bl
VPB152-EB-062514	WQ	1,1,1-TRICHLOROETHANE		0.50	UG/L	UJ	t
VPB152-EB-062514	WQ	1,1,2,2-TETRACHLOROETHANE		0.50	UG/L	UJ	t
VPB152-EB-062514	WQ	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE		0.50	UG/L	UJ	t,c
VPB152-EB-062514	WQ	1,1,2-TRICHLOROETHANE		0.50	UG/L	UJ	t
VPB152-EB-062514	WQ	1,1-DICHLOROETHANE		0.50	UG/L	UJ	t
VPB152-EB-062514	WQ	1,1-DICHLOROETHENE		0.50	UG/L	UJ	t,c
VPB152-EB-062514	WQ	1,2,4-TRICHLOROBENZENE		0.50	UG/L	UJ	t
VPB152-EB-062514	WQ	1,2-DIBROMO-3-CHLOROPROPANE		0.75	UG/L	UJ	t
VPB152-EB-062514	WQ	1,2-DIBROMOETHANE		0.50	UG/L	UJ	t
VPB152-EB-062514	WQ	1,2-DICHLOROBENZENE		0.50	UG/L	UJ	t
VPB152-EB-062514	WQ	1,2-DICHLOROETHANE		0.50	UG/L	UJ	t
VPB152-EB-062514	WQ	1,2-DICHLOROETHENE, TOTAL		1.0	UG/L	UJ	t
VPB152-EB-062514	WQ	1,2-DICHLOROPROPANE		0.50	UG/L	UJ	t
VPB152-EB-062514	WQ	1,3-DICHLOROBENZENE		0.50	UG/L	UJ	t
VPB152-EB-062514	WQ	1,4-DICHLOROBENZENE		0.50	UG/L	UJ	t
VPB152-EB-062514	WQ	2-BUTANONE		2.5	UG/L	UJ	t
VPB152-EB-062514	WQ	2-HEXANONE		2.5	UG/L	UJ	t
VPB152-EB-062514	WQ	4-METHYL-2-PENTANONE		2.5	UG/L	UJ	t
VPB152-EB-062514	WQ	ACETONE	8.1	2.5	UG/L	J	t,c
VPB152-EB-062514	WQ	BENZENE		0.50	UG/L	UJ	t
VPB152-EB-062514	WQ	BROMODICHLOROMETHANE		0.50	UG/L	UJ	t
VPB152-EB-062514	WQ	BROMOFORM		0.50	UG/L	UJ	t
VPB152-EB-062514	WQ	BROMOMETHANE		1.0	UG/L	UJ	t
VPB152-EB-062514	WQ	CARBON DISULFIDE		0.50	UG/L	UJ	t,c
VPB152-EB-062514	WQ	CARBON TETRACHLORIDE		0.50	UG/L	UJ	t
VPB152-EB-062514	WQ	CHLOROBENZENE		0.50	UG/L	UJ	t
VPB152-EB-062514	WQ	CHLOROETHANE		1.0	UG/L	UJ	t
VPB152-EB-062514	WQ	CHLOROFORM		0.50	UG/L	UJ	t
VPB152-EB-062514	WQ	CHLOROMETHANE		1.0	UG/L	UJ	t
VPB152-EB-062514	WQ	CIS-1,2-DICHLOROETHENE		0.50	UG/L	UJ	t
VPB152-EB-062514	WQ	CIS-1,3-DICHLOROPROPENE		0.50	UG/L	UJ	t
VPB152-EB-062514	WQ	CYCLOHEXANE		0.50	UG/L	UJ	t
VPB152-EB-062514	WQ	DIBROMOCHLOROMETHANE		0.50	UG/L	UJ	t
VPB152-EB-062514	WQ	DICHLORODIFLUOROMETHANE		1.0	UG/L	UJ	t
VPB152-EB-062514	WQ	ETHYLBENZENE		0.50	UG/L	UJ	t
VPB152-EB-062514	WQ	ISOPROPYLBENZENE		0.50	UG/L	UJ	t
VPB152-EB-062514	WQ	M- AND P-XYLENE		1.0	UG/L	UJ	t
VPB152-EB-062514	WQ	METHYL ACETATE		0.75	UG/L	UJ	t

Sample ID	Matrix	Compound	Result	LOD	Units	Validation Qualifiers	Validation Reason
VPB152-EB-062514	WQ	METHYL CYCLOHEXANE		0.50	UG/L	UJ	t,c
VPB152-EB-062514	WQ	METHYL TERT-BUTYL ETHER		0.50	UG/L	UJ	t
VPB152-EB-062514	WQ	METHYLENE CHLORIDE		2.5	UG/L	UJ	t,c
VPB152-EB-062514	WQ	O-XYLENE		0.50	UG/L	UJ	t
VPB152-EB-062514	WQ	STYRENE		0.50	UG/L	UJ	t
VPB152-EB-062514	WQ	TETRACHLOROETHENE		0.50	UG/L	UJ	t
VPB152-EB-062514	WQ	TOLUENE		0.50	UG/L	UJ	t
VPB152-EB-062514	WQ	TRANS-1,2-DICHLOROETHENE		0.50	UG/L	UJ	t
VPB152-EB-062514	WQ	TRANS-1,3-DICHLOROPROPENE		0.50	UG/L	UJ	t
VPB152-EB-062514	WQ	TRICHLOROETHENE		0.50	UG/L	UJ	t
VPB152-EB-062514	WQ	TRICHLOROFLUOROMETHANE		1.0	UG/L	UJ	t
VPB152-EB-062514	WQ	VINYL CHLORIDE		1.0	UG/L	UJ	t
VPB152-EB-062514	WQ	XYLENES, TOTAL		1.5	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	1,1,1-TRICHLOROETHANE		0.50	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	1,1,2,2-TETRACHLOROETHANE		0.50	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE		0.50	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	1,1,2-TRICHLOROETHANE		0.50	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	1,1-DICHLOROETHANE		0.50	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	1,1-DICHLOROETHENE		0.50	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	1,2,4-TRICHLOROBENZENE		0.50	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	1,2-DIBROMO-3-CHLOROPROPANE		0.75	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	1,2-DIBROMOETHANE		0.50	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	1,2-DICHLOROBENZENE		0.50	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	1,2-DICHLOROETHANE		0.50	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	1,2-DICHLOROETHENE, TOTAL		1.0	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	1,2-DICHLOROPROPANE		0.50	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	1,3-DICHLOROBENZENE		0.50	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	1,4-DICHLOROBENZENE		0.50	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	2-BUTANONE		2.5	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	2-HEXANONE		2.5	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	4-METHYL-2-PENTANONE		2.5	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	ACETONE		5.0*	UG/L	UJ	t,be
VPB152-GW-062414-870-872	WG	BENZENE		0.50	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	BROMODICHLOROMETHANE		0.50	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	BROMOFORM		0.50	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	BROMOMETHANE		1.0	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	CARBON DISULFIDE		0.50	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	CARBON TETRACHLORIDE		0.50	UG/L	UJ	t

Sample ID	Matrix	Compound	Result	LOD	Units	Validation Qualifiers	Validation Reason
VPB152-GW-062414-870-872	WG	CHLOROBENZENE		0.50	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	CHLOROETHANE		1.0	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	CHLOROFORM		0.50	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	CHLOROMETHANE		1.0	UG/L	UJ	t,c
VPB152-GW-062414-870-872	WG	CIS-1,2-DICHLOROETHENE		0.50	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	CIS-1,3-DICHLOROPROPENE		0.50	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	CYCLOHEXANE		0.50	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	DIBROMOCHLOROMETHANE		0.50	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	DICHLORODIFLUOROMETHANE		1.0	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	ETHYLBENZENE		0.50	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	ISOPROPYLBENZENE		0.50	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	M- AND P-XYLENE		1.0	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	METHYL ACETATE		0.75	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	METHYL CYCLOHEXANE		0.50	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	METHYL TERT-BUTYL ETHER		0.50	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	METHYLENE CHLORIDE		2.5	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	O-XYLENE		0.50	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	STYRENE		0.50	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	TETRACHLOROETHENE		0.50	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	TOLUENE		0.50	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	TRANS-1,2-DICHLOROETHENE		0.50	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	TRANS-1,3-DICHLOROPROPENE		0.50	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	TRICHLOROETHENE		0.50	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	TRICHLOROFUOROMETHANE		1.0	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	VINYL CHLORIDE		1.0	UG/L	UJ	t
VPB152-GW-062414-870-872	WG	XYLENES, TOTAL		1.5	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	1,1,1-TRICHLOROETHANE		0.50	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	1,1,2,2-TETRACHLOROETHANE		0.50	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE		0.50	UG/L	UJ	t,c
VPB152-GW-062414-880-882	WG	1,1,2-TRICHLOROETHANE		0.50	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	1,1-DICHLOROETHANE		0.50	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	1,1-DICHLOROETHENE		0.50	UG/L	UJ	t,c
VPB152-GW-062414-880-882	WG	1,2,4-TRICHLOROBENZENE		0.50	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	1,2-DIBROMO-3-CHLOROPROPANE		0.75	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	1,2-DIBROMOETHANE		0.50	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	1,2-DICHLOROBENZENE		0.50	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	1,2-DICHLOROETHANE		0.50	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	1,2-DICHLOROETHENE, TOTAL		1.0	UG/L	UJ	t

Sample ID	Matrix	Compound	Result	LOD	Units	Validation Qualifiers	Validation Reason
VPB152-GW-062414-880-882	WG	1,2-DICHLOROPROPANE		0.50	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	1,3-DICHLOROBENZENE		0.50	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	1,4-DICHLOROBENZENE		0.50	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	2-BUTANONE		2.5	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	2-HEXANONE		2.5	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	4-METHYL-2-PENTANONE		2.5	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	ACETONE		13*	UG/L	UJ	t,be
VPB152-GW-062414-880-882	WG	BENZENE		0.50	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	BROMODICHLOROMETHANE		0.50	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	BROMOFORM		0.50	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	BROMOMETHANE		1.0	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	CARBON DISULFIDE		0.50	UG/L	UJ	t,c
VPB152-GW-062414-880-882	WG	CARBON TETRACHLORIDE		0.50	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	CHLOROBENZENE		0.50	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	CHLOROETHANE		1.0	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	CHLOROFORM		0.50	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	CHLOROMETHANE		1.0	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	CIS-1,2-DICHLOROETHENE		0.50	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	CIS-1,3-DICHLOROPROPENE		0.50	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	CYCLOHEXANE		0.50	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	DIBROMOCHLOROMETHANE		0.50	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	DICHLORODIFLUOROMETHANE		1.0	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	ETHYLBENZENE		0.50	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	ISOPROPYLBENZENE		0.50	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	M- AND P-XYLENE		1.0	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	METHYL ACETATE		0.75	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	METHYL CYCLOHEXANE		0.50	UG/L	UJ	t,c
VPB152-GW-062414-880-882	WG	METHYL TERT-BUTYL ETHER		0.50	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	METHYLENE CHLORIDE		2.5	UG/L	UJ	t,c
VPB152-GW-062414-880-882	WG	O-XYLENE		0.50	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	STYRENE		0.50	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	TETRACHLOROETHENE		0.50	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	TOLUENE	1.4	0.50	UG/L	J	t
VPB152-GW-062414-880-882	WG	TRANS-1,2-DICHLOROETHENE		0.50	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	TRANS-1,3-DICHLOROPROPENE		0.50	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	TRICHLOROETHENE		0.50	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	TRICHLOROFLUOROMETHANE		1.0	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	VINYL CHLORIDE		1.0	UG/L	UJ	t
VPB152-GW-062414-880-882	WG	XYLENES, TOTAL		1.5	UG/L	UJ	t

Sample ID	Matrix	Compound	Result	LOD	Units	Validation Qualifiers	Validation Reason
VPB152-GW-062514-910-912	WG	1,1,1-TRICHLOROETHANE		0.50	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	1,1,2,2-TETRACHLOROETHANE		0.50	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE		0.50	UG/L	UJ	t,c
VPB152-GW-062514-910-912	WG	1,1,2-TRICHLOROETHANE		0.50	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	1,1-DICHLOROETHANE		0.50	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	1,1-DICHLOROETHENE		0.50	UG/L	UJ	t,c
VPB152-GW-062514-910-912	WG	1,2,4-TRICHLOROBENZENE		0.50	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	1,2-DIBROMO-3-CHLOROPROPANE		0.75	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	1,2-DIBROMOETHANE		0.50	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	1,2-DICHLOROBENZENE		0.50	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	1,2-DICHLOROETHANE		0.50	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	1,2-DICHLOROETHENE, TOTAL		1.0	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	1,2-DICHLOROPROPANE		0.50	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	1,3-DICHLOROBENZENE		0.50	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	1,4-DICHLOROBENZENE		0.50	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	2-BUTANONE		2.5	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	2-HEXANONE		2.5	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	4-METHYL-2-PENTANONE		2.5	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	ACETONE		5.0*	UG/L	UJ	t,be
VPB152-GW-062514-910-912	WG	BENZENE		0.50	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	BROMODICHLOROMETHANE		0.50	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	BROMOFORM		0.50	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	BROMOMETHANE		1.0	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	CARBON DISULFIDE		0.50	UG/L	UJ	t,c
VPB152-GW-062514-910-912	WG	CARBON TETRACHLORIDE		0.50	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	CHLOROBENZENE		0.50	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	CHLOROETHANE		1.0	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	CHLOROFORM		0.50	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	CHLOROMETHANE		1.0	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	CIS-1,2-DICHLOROETHENE		0.50	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	CIS-1,3-DICHLOROPROPENE		0.50	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	CYCLOHEXANE		0.50	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	DIBROMOCHLOROMETHANE		0.50	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	DICHLORODIFLUOROMETHANE		1.0	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	ETHYLBENZENE		0.50	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	ISOPROPYLBENZENE		0.50	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	M- AND P-XYLENE		1.0	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	METHYL ACETATE		0.75	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	METHYL CYCLOHEXANE		0.50	UG/L	UJ	t,c

VPB152-GW-062514-910-912	WG	METHYL TERT-BUTYL ETHER		0.50	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	METHYLENE CHLORIDE		2.5	UG/L	UJ	t,c
VPB152-GW-062514-910-912	WG	O-XYLENE		0.50	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	STYRENE		0.50	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	TETRACHLOROETHENE		0.50	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	TOLUENE		0.50	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	TRANS-1,2-DICHLOROETHENE		0.50	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	TRANS-1,3-DICHLOROPROPENE		0.50	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	TRICHLOROETHENE		0.50	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	TRICHLOROFLUOROMETHANE		1.0	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	VINYL CHLORIDE		1.0	UG/L	UJ	t
VPB152-GW-062514-910-912	WG	XYLENES, TOTAL		1.5	UG/L	UJ	t
VPB152-GW-062614-930-932	WG	1,1,1-TRICHLOROETHANE		1.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	1,1,2,2-TETRACHLOROETHANE		1.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE		1.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	1,1,2-TRICHLOROETHANE		1.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	1,1-DICHLOROETHANE		1.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	1,1-DICHLOROETHENE		1.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	1,2,4-TRICHLOROBENZENE		1.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	1,2-DIBROMO-3-CHLOROPROPANE		1.5	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	1,2-DIBROMOETHANE		1.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	1,2-DICHLOROBENZENE		1.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	1,2-DICHLOROETHANE		1.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	1,2-DICHLOROETHENE, TOTAL		2.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	1,2-DICHLOROPROPANE		1.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	1,3-DICHLOROBENZENE		1.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	1,4-DICHLOROBENZENE		1.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	2-BUTANONE		5.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	2-HEXANONE		5.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	4-METHYL-2-PENTANONE		5.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	ACETONE		32*	UG/L	UJ	t,be,mc
VPB152-GW-062614-930-932	WG	BENZENE		1.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	BROMODICHLOROMETHANE		1.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	BROMOFORM		1.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	BROMOMETHANE		2.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	CARBON DISULFIDE		1.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	CARBON TETRACHLORIDE		1.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	CHLOROBENZENE		1.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	CHLOROETHANE		2.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	CHLOROFORM		1.0	UG/L	UJ	t,mc

VPB152-GW-062614-930-932	WG	CHLOROMETHANE		2.0	UG/L	UJ	t,mc,c
VPB152-GW-062614-930-932	WG	CIS-1,2-DICHLOROETHENE		1.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	CIS-1,3-DICHLOROPROPENE		1.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	CYCLOHEXANE		1.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	DIBROMOCHLOROMETHANE		1.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	DICHLORODIFLUOROMETHANE		2.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	ETHYLBENZENE		1.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	ISOPROPYLBENZENE		1.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	M- AND P-XYLENE		2.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	METHYL ACETATE		1.5	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	METHYL CYCLOHEXANE		1.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	METHYL TERT-BUTYL ETHER		1.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	METHYLENE CHLORIDE	3.2	5.0	UG/L	J	mc,t
VPB152-GW-062614-930-932	WG	O-XYLENE		1.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	STYRENE		1.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	TETRACHLOROETHENE		1.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	TOLUENE	3.8	1.0	UG/L	J	mc,t
VPB152-GW-062614-930-932	WG	TRANS-1,2-DICHLOROETHENE		1.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	TRANS-1,3-DICHLOROPROPENE		1.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	TRICHLOROETHENE		1.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	TRICHLOROFUOROMETHANE		2.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	VINYL CHLORIDE		2.0	UG/L	UJ	t,mc
VPB152-GW-062614-930-932	WG	XYLENES, TOTAL		3.0	UG/L	UJ	t,mc
VPB152-TRIPBLANK-062614	WQ	1,1,1-TRICHLOROETHANE		0.50	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	1,1,2,2-TETRACHLOROETHANE		0.50	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE		0.50	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	1,1,2-TRICHLOROETHANE		0.50	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	1,1-DICHLOROETHANE		0.50	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	1,1-DICHLOROETHENE		0.50	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	1,2,4-TRICHLOROBENZENE		0.50	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	1,2-DIBROMO-3-CHLOROPROPANE		0.75	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	1,2-DIBROMOETHANE		0.50	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	1,2-DICHLOROBENZENE		0.50	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	1,2-DICHLOROETHANE		0.50	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	1,2-DICHLOROETHENE, TOTAL		1.0	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	1,2-DICHLOROPROPANE		0.50	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	1,3-DICHLOROBENZENE		0.50	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	1,4-DICHLOROBENZENE		0.50	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	2-BUTANONE		2.5	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	2-HEXANONE		2.5	UG/L	UJ	t



VPB152-TRIPBLANK-062614	WQ	4-METHYL-2-PENTANONE		2.5	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	ACETONE		2.5	UG/L	UL	t
VPB152-TRIPBLANK-062614	WQ	BENZENE		0.50	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	BROMODICHLOROMETHANE		0.50	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	BROMOFORM		0.50	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	BROMOMETHANE		1.0	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	CARBON DISULFIDE		0.50	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	CARBON TETRACHLORIDE		0.50	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	CHLOROBENZENE		0.50	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	CHLOROETHANE		1.0	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	CHLOROFORM		0.50	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	CHLOROMETHANE		1.0	UG/L	UJ	t,c
VPB152-TRIPBLANK-062614	WQ	CIS-1,2-DICHLOROETHENE		0.50	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	CIS-1,3-DICHLOROPROPENE		0.50	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	CYCLOHEXANE		0.50	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	DIBROMOCHLOROMETHANE		0.50	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	DICHLORODIFLUOROMETHANE		1.0	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	ETHYLBENZENE		0.50	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	ISOPROPYLBENZENE		0.50	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	M- AND P-XYLENE		1.0	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	METHYL ACETATE		0.75	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	METHYL CYCLOHEXANE		0.50	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	METHYL TERT-BUTYL ETHER		0.50	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	METHYLENE CHLORIDE		2.5	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	O-XYLENE		0.50	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	STYRENE		0.50	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	TETRACHLOROETHENE		0.50	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	TOLUENE		0.50	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	TRANS-1,2-DICHLOROETHENE		0.50	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	TRANS-1,3-DICHLOROPROPENE		0.50	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	TRICHLOROETHENE		0.50	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	TRICHLOROFLUOROMETHANE		1.0	UG/L	UJ	t
VPB152-TRIPBLANK-062614	WQ	VINYL CHLORIDE		1.0	UG/L	UJ	t
*LOQ							

## Attachment A

## Nonconformance Summary Tables

Table A-1a - Initial Calibration

Calibration Date/Time	Compound	% RSD	Limits
01-July-2014 09:21	CHLOROETHANE	16	≤15%
Associated samples: VPB152-GW-062414-870-872, VPB152-TRIP BLANK-062614,VPB152-GW-062614-930-932			

Table A-1b - Initial Calibration Verification

ICV	Compound	% R	Limit
WG145769-7	ACETONE	128	80-120%
Associated samples: VPB152-GW-062414-870-872, VPB152-TRIP BLANK-062614,VPB152-GW-062614-930-932			
WG146050-7	CARBON DISULFIDE	141	80-120%
	ACETONE	163	80-120%
	2-BUTANONE	122	80-120%
Associated samples: VPB152-GW-062414-880-882, VPB152-EB-062514,VPB152-GW-062514-910-912			

Table A-1c - Continuing Calibration Verification

CCV ID	Compound	% D	Limit
WG146127-4	1,1-DICHLOROETHENE	22	≤ 20%
	CARBON DISULFIDE	27	≤ 20%
	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	39	≤ 20%
	METHYLENE CHLORIDE	24	≤ 20%
Associated samples: VPB152-GW-062414-880-882, VPB152-EB-062514,VPB152-GW-062514-910-912			

Table A-3 - Lab Blanks

Blank ID	Compound	Result	LOD	Units	Associated Samples
WG146111-1	TOTAL ORGANIC CARBON	0.26	0.50	MG/L	VPB152-EB-062514

Table A-4 - Field Blanks

Blank ID	Compound	Result	LOD	Units	Associated Samples
VPB152-EB-062514	ACETONE	8.1	2.5	UG/L	Samples in SDG SH4793

**Attachment B****Qualifier Codes and Explanations**

<b>Qualifier</b>	<b>Explanation</b>
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.

## Attachment C

### Reason Codes and Explanations

Reason Code	Explanation
be	Equipment blank contamination
bf	Field blank contamination
bl	Laboratory blank contamination
c	Calibration issue
co	Analyte carryover
d	Reporting limit raised due to chromatographic interference
fd	Field duplicate RPDs
h	Holding times
i	Internal standard areas
k	Estimated Maximum Possible Concentration (EMPC)
l	LCS or OPR recoveries
lc	Labeled compound recovery
ld	Laboratory duplicate RPDs
lp	Laboratory control sample/laboratory control sample duplicate RPDs
m	Matrix spike recovery
md	Matrix spike/matrix spike duplicate RPDs
nb	Negative laboratory blank contamination
p	Chemical preservation issue
r	Dual column RPD
q	Quantitation issue
s	Surrogate recovery
su	Ion suppression
t	Temperature preservation issue
x	Percent solids
y	Serial dilution results
z	ICS results
mc	Method compliance nonconformance



## CHAIN of CUSTODY

PLEASE BEAR DOWN AND  
 PRINT LEGIBLY IN PEN

Client: Resolution Consultants Contact: E. Uioanda Phone #: (845) 425-4180 Fax #: ( )

Address: 100 Red Schoolhouse Rd City: Chastnut Ridge State: NY Zip Code: 10977

Purchase Order #: \_\_\_\_\_ Proj. Name / No.: NR1RP Bathpage / 66866526 Katahdin Quote #: \_\_\_\_\_

Bill (if different than above): \_\_\_\_\_ Address: \_\_\_\_\_

Sampler (Print / Sign): Vincent Viorakio / Nat Viorak Copies To: \_\_\_\_\_

LAB USE ONLY WORK ORDER #: SH4793  
 KATAHDIN PROJECT NUMBER \_\_\_\_\_

ANALYSIS AND CONTAINER TYPE PRESERVATIVES

REMARKS: \_\_\_\_\_

SHIPPING INFO:  FED EX  UPS  CLIENT

AIRBILL NO: \_\_\_\_\_

TEMP °C \_\_\_\_\_  TEMP BLANK  INTACT  NOT INTACT

Filt.	Filt.	Filt.	Filt.	Filt.	Filt.	Filt.	Filt.	Filt.	Filt.	Filt.																		
Y	O	N	Y	O	N	Y	O	N	Y	O	N	Y	O	N	Y	O	N	Y	O	N	Y	O	N					

COMMENTS \_\_\_\_\_

17.46

Relinquished By: (Signature) <u>[Signature]</u>	Date / Time <u>6/26 17:15</u>	Received By: (Signature) <u>FedEx</u>	Relinquished By: (Signature)	Date / Time	Received By: (Signature)
Relinquished By: (Signature)	Date / Time <u>6/30 9:15</u>	Received By: (Signature) <u>Jennifer Obrien</u>	Relinquished By: (Signature)	Date / Time	Received By: (Signature)



## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4793-1  
**Client ID:** 152-062414-870-872  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4793  
**Lab File ID:** D9267.D

**Sample Date:** 24-JUN-14  
**Received Date:** 30-JUN-14  
**Extract Date:** 01-JUL-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG145769

**Analysis Date:** 01-JUL-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260C  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 17-JUL-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	4.5 1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
Acetone	L	<del>7.2</del> 5.0	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50

*Handwritten:* 7/25/14



## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4793-1  
**Client ID:** 152-062414-870-872  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4793  
**Lab File ID:** D9267.D

**Sample Date:** 24-JUN-14  
**Received Date:** 30-JUN-14  
**Extract Date:** 01-JUL-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG145769

**Analysis Date:** 01-JUL-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260C  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 17-JUL-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		85.7	%					
Toluene-d8		111.	%					
1,2-Dichloroethane-d4		115.	%					
Dibromofluoromethane		103.	%					

DJP/75/14

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4793-3RA  
**Client ID:** 152-062414-880-882  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4793  
**Lab File ID:** C8048.D

**Sample Date:** 24-JUN-14  
**Received Date:** 30-JUN-14  
**Extract Date:** 08-JUL-14  
**Extracted By:** REC  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG146127

**Analysis Date:** 08-JUL-14  
**Analyst:** REC  
**Analysis Method:** SW846 8260C  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 17-JUL-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
<b>Acetone</b>		13	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
<b>Toluene</b>		1.4	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50

*Handwritten signature:* F. 11/25/14

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4793-3RA  
**Client ID:** 152-062414-880-882  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4793  
**Lab File ID:** C8048.D

**Sample Date:** 24-JUN-14  
**Received Date:** 30-JUN-14  
**Extract Date:** 08-JUL-14  
**Extracted By:** REC  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG146127

**Analysis Date:** 08-JUL-14  
**Analyst:** REC  
**Analysis Method:** SW846 8260C  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 17-JUL-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U <i>55</i>	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		93.5	%					
Toluene-d8		102.	%					
1,2-Dichloroethane-d4		109.	%					
Dibromofluoromethane		97.7	%					

*Rulzsky*

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4793-4RA  
**Client ID:** VPB152-EB-062514  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4793  
**Lab File ID:** C8049.D

**Sample Date:** 25-JUN-14  
**Received Date:** 30-JUN-14  
**Extract Date:** 08-JUL-14  
**Extracted By:** REC  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG146127

**Analysis Date:** 08-JUL-14  
**Analyst:** REC  
**Analysis Method:** SW846 8260C  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 17-JUL-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
Acetone	U	8.1	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50

JUN 25/14

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4793-4RA  
**Client ID:** VPB152-EB-062514  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4793  
**Lab File ID:** C8049.D

**Sample Date:** 25-JUN-14  
**Received Date:** 30-JUN-14  
**Extract Date:** 08-JUL-14  
**Extracted By:** REC  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG146127

**Analysis Date:** 08-JUL-14  
**Analyst:** REC  
**Analysis Method:** SW846 8260C  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 17-JUL-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		89.8	%					
Toluene-d8		95.8	%					
1,2-Dichloroethane-d4		101.	%					
Dibromofluoromethane		90.7	%					

R 11/25/14

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4793-5  
**Client ID:** VPB152-TB-062614  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4793  
**Lab File ID:** D9263.D

**Sample Date:** 26-JUN-14  
**Received Date:** 30-JUN-14  
**Extract Date:** 01-JUL-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG145769

**Analysis Date:** 01-JUL-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260C  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 17-JUL-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
Acetone	UL	2.5	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50

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### Report of Analytical Results

Client: ENSAFE  
 Lab ID: SH4793-5  
 Client ID: VPB152-TB-062614  
 Project: Navy Clean WE15-03-06 NW  
 SDG: SH4793  
 Lab File ID: D9263.D

Sample Date: 26-JUN-14  
 Received Date: 30-JUN-14  
 Extract Date: 01-JUL-14  
 Extracted By: DJP  
 Extraction Method: SW846 5030  
 Lab Prep Batch: WG145769

Analysis Date: 01-JUL-14  
 Analyst: DJP  
 Analysis Method: SW846 8260C  
 Matrix: AQ  
 % Solids: NA  
 Report Date: 17-JUL-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		85.9	%					
Toluene-d8		110.	%					
1,2-Dichloroethane-d4		110.	%					
Dibromofluoromethane		102.	%					

*Ric/25/17*



## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4793-6RA  
**Client ID:** 152-062514-910-912  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4793  
**Lab File ID:** C8050.D

**Sample Date:** 25-JUN-14  
**Received Date:** 30-JUN-14  
**Extract Date:** 08-JUL-14  
**Extracted By:** REC  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG146127

**Analysis Date:** 08-JUL-14  
**Analyst:** REC  
**Analysis Method:** SW846 8260C  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 17-JUL-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
Acetone	U	<del>9.7</del> 5.0	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50

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### Report of Analytical Results

Client: ENSAFE  
 Lab ID: SH4793-6RA  
 Client ID: 152-062514-910-912  
 Project: Navy Clean WE15-03-06 NW  
 SDG: SH4793  
 Lab File ID: C8050.D

Sample Date: 25-JUN-14  
 Received Date: 30-JUN-14  
 Extract Date: 08-JUL-14  
 Extracted By: REC  
 Extraction Method: SW846 5030  
 Lab Prep Batch: WG146127

Analysis Date: 08-JUL-14  
 Analyst: REC  
 Analysis Method: SW846 8260C  
 Matrix: AQ  
 % Solids: NA  
 Report Date: 17-JUL-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U UJ	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		92.5	%					
Toluene-d8		97.5	%					
1,2-Dichloroethane-d4		104.	%					
Dibromofluoromethane		92.2	%					

*Handwritten signature and date: 6/25/14*

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4793-7DL  
**Client ID:** 152-062614-930-932  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4793  
**Lab File ID:** D9265.D

**Sample Date:** 26-JUN-14  
**Received Date:** 30-JUN-14  
**Extract Date:** 01-JUL-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG145769

**Analysis Date:** 01-JUL-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260C  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 17-JUL-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	2.0	ug/L	2	2	4.0	0.48	2.0
Chloromethane	U	2.0	ug/L	2	2	4.0	0.72	2.0
Vinyl Chloride	U	2.0	ug/L	2	2	4.0	0.50	2.0
Bromomethane	U	2.0	ug/L	2	2	4.0	0.98	2.0
Chloroethane	U	2.0	ug/L	2	2	4.0	1.1	2.0
Trichlorofluoromethane	U	2.0	ug/L	2	2	4.0	0.48	2.0
1,1-Dichloroethene	U	1.0	ug/L	2	1	2.0	0.70	1.0
Carbon Disulfide	U	1.0	ug/L	2	1	2.0	0.50	1.0
Freon-113	U	1.0	ug/L	2	1	2.0	0.62	1.0
<b>Methylene Chloride</b>	J	3.2	ug/L	2	5	10.	2.3	5.0
<b>Acetone</b>	L	32	ug/L	2	5	10.	4.4	5.0
trans-1,2-Dichloroethene	U	1.0	ug/L	2	1	2.0	0.50	1.0
Methyl tert-butyl Ether	U	1.0	ug/L	2	1	2.0	0.72	1.0
1,1-Dichloroethane	U	1.0	ug/L	2	1	2.0	0.42	1.0
cis-1,2-Dichloroethene	U	1.0	ug/L	2	1	2.0	0.42	1.0
Chloroform	U	1.0	ug/L	2	1	2.0	0.64	1.0
1,1,1-Trichloroethane	U	1.0	ug/L	2	1	2.0	0.40	1.0
2-Butanone	U	5.0	ug/L	2	5	10.	2.6	5.0
Cyclohexane	U	1.0	ug/L	2	1	2.0	0.62	1.0
Carbon Tetrachloride	U	1.0	ug/L	2	1	2.0	0.44	1.0
Benzene	U	1.0	ug/L	2	1	2.0	0.52	1.0
1,2-Dichloroethane	U	1.0	ug/L	2	1	2.0	0.40	1.0
Trichloroethene	U	1.0	ug/L	2	1	2.0	0.56	1.0
1,2-Dichloropropane	U	1.0	ug/L	2	1	2.0	0.50	1.0
Bromodichloromethane	U	1.0	ug/L	2	1	2.0	0.66	1.0
cis-1,3-Dichloropropene	U	1.0	ug/L	2	1	2.0	0.38	1.0
<b>Toluene</b>	J	3.8	ug/L	2	1	2.0	0.54	1.0
4-Methyl-2-Pentanone	U	5.0	ug/L	2	5	10.	2.6	5.0
trans-1,3-Dichloropropene	U	1.0	ug/L	2	1	2.0	0.40	1.0
1,1,2-Trichloroethane	U	1.0	ug/L	2	1	2.0	0.66	1.0
Tetrachloroethene	U	1.0	ug/L	2	1	2.0	0.80	1.0
Dibromochloromethane	U	1.0	ug/L	2	1	2.0	0.60	1.0
2-Hexanone	U	5.0	ug/L	2	5	10.	3.4	5.0
Chlorobenzene	U	1.0	ug/L	2	1	2.0	0.44	1.0
Ethylbenzene	U	1.0	ug/L	2	1	2.0	0.42	1.0

*R. 1/25/14*

### Report of Analytical Results

Client: ENSAFE  
 Lab ID: SH4793-7DL  
 Client ID: 152-062614-930-932  
 Project: Navy Clean WE15-03-06 NW  
 SDG: SH4793  
 Lab File ID: D9265.D

Sample Date: 26-JUN-14  
 Received Date: 30-JUN-14  
 Extract Date: 01-JUL-14  
 Extracted By: DJP  
 Extraction Method: SW846 5030  
 Lab Prep Batch: WG145769

Analysis Date: 01-JUL-14  
 Analyst: DJP  
 Analysis Method: SW846 8260C  
 Matrix: AQ  
 % Solids: NA  
 Report Date: 17-JUL-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	3.0	ug/L	2	3	6.0	0.50	3.0
Styrene	U	1.0	ug/L	2	1	2.0	0.46	1.0
Bromoform	U	1.0	ug/L	2	1	2.0	0.46	1.0
Isopropylbenzene	U	1.0	ug/L	2	1	2.0	0.46	1.0
1,1,2,2-Tetrachloroethane	U	1.0	ug/L	2	1	2.0	0.76	1.0
1,3-Dichlorobenzene	U	1.0	ug/L	2	1	2.0	0.52	1.0
1,4-Dichlorobenzene	U	1.0	ug/L	2	1	2.0	0.48	1.0
1,2-Dichlorobenzene	U	1.0	ug/L	2	1	2.0	0.30	1.0
1,2,4-Trichlorobenzene	U	1.0	ug/L	2	1	2.0	0.74	1.0
Methyl Acetate	U	1.5	ug/L	2	1	2.0	1.1	1.5
Methylcyclohexane	U	1.0	ug/L	2	1	2.0	0.60	1.0
o-Xylene	U	1.0	ug/L	2	1	2.0	0.50	1.0
M+P-Xylenes	U	2.0	ug/L	2	2	4.0	1.2	2.0
1,2-Dichloroethylene (Total)	U	2.0	ug/L	2	2	4.0	0.42	2.0
1,2-Dibromoethane	U	1.0	ug/L	2	1	2.0	0.44	1.0
1,2-Dibromo-3-Chloropropane	U	1.5	ug/L	2	1	2.0	1.0	1.5
P-Bromofluorobenzene		85.4	%					
Toluene-d8		110.	%					
1,2-Dichloroethane-d4		110.	%					
Dibromofluoromethane		103.	%					

*R 11/25/14*

## Report of Analytical Results

**Client:** Rick Purdy  
AECOM  
701 Edgewater Drive  
Wakefield, MA 01880

**Lab Sample ID:** SH4793-4  
**Report Date:** 19-JUL-14  
**Client PO:** 16518  
**Project:** Navy Clean WE15-03-0  
**SDG:** SH4793

Sample Description  
VPB152-EB-062514

Matrix AQ      Date Sampled 25-JUN-14      Date Received 30-JUN-14

Parameter	Result	Adj LOQ	Adj MDL	Adj LOD	Anst. Method	QC.Batch	Anal. Date	Prep. Method	Prep. Date	Footnotes
Total Organic Carbon	10.37 mg/L	1.0	0.10	.5	SM5310B	WG146111	30-JUN-14 19:35:58	N/A	N/A	N/A

1.0 U

10.37





## Data Validation Report

Project: Regional Groundwater Investigation - NWIRP Bethpage  
Laboratory: Katahdin Analytical Services, Scarborough, Maine  
Service Request: SH4577  
Analyses/Method: EPA SW-846 Method 8260B for VOCs (GC/MS)  
Validation Level: Limited  
RESCON Project 60266526.SA.DV  
Number: \_\_\_\_\_  
Prepared by: Sheena Blair/RESCON Completed on: 8/30/2014  
Reviewed by: Lori Herberich/RESCON File Name: SH4577\_8260B

### SUMMARY

The samples listed below were collected by Resolution Consultants (RESCON) from the Regional Groundwater Investigation - NWIRP Bethpage site on June 20 and 23, 2014.

Sample ID	Matrix/Sample Type
VPB152-GW-062014-830-832	Groundwater
VPB152-GW-062314-850-852	Groundwater
VPB152-TRIP BLANK-062314	Trip Blank

Data validation activities were conducted with reference to *Test Methods for Evaluating Solid Waste, Physical/Chemical Methods SW846, specifically SW-846 Method 8260B, Volatile Organic Compounds by Gas Chromatography/Mass Spectrometry* (USEPA, 1996), *USEPA Contract Laboratory Program National Functional Guidelines for Superfund Organic Methods Data Review* (June 2008), and *Quality Systems Manual (QSM) for Environmental Laboratories, Version 4.2* (DoD, 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 review elements (where applicable to the method):

- ✓ Data completeness (chain-of-custody [COC])/sample integrity
- ✓ Holding times and sample preservation
- ✓ GC/MS performance checks
- ✗ Initial calibration/continuing calibration verification
- ✓ Laboratory blanks/trip blanks/equipment blanks
- ✓ Surrogate spike recoveries
- NA Matrix spike (MS) and/or matrix spike duplicate (MSD) results
- ✓ Laboratory control sample (LCS) results
- NA Field duplicate results
- ✓ Internal standard results

✓ 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 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. Selected data points were estimated due to nonconformances of certain QC criteria (see discussion below). Qualified sample results are presented in Table 1.

## 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 "VPB" prefix and the "GW" from the sample ID, and truncated the ID for the Trip Blank 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.

### GC/MS Performance Checks

The data were reviewed to ensure that the 4-bromofluorobenzene (BFB) tuning was performed at the correct frequency and that the method acceptance criteria were met. 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:

- the initial calibration (ICAL) percent relative standard deviation (%RSD), correlation coefficient (r)/coefficient of determination ( $r^2$ ), and/or response factor method acceptance criteria were met;
- the initial calibration verification standard (ICV) percent recoveries (%Rs) acceptance criteria were met;

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

Data qualification to the analytes associated with the specific ICAL, ICV and/or CCV was as follows:

**ICAL Linearity Nonconformances:**

Nonconformance	Actions	
	Detected Results	Nondetected Results
%RSD > 15% and quantitation based on mean RF	J	UJ
r or r <sup>2</sup> < 0.99 and quantitation based on linear regression	J*	UJ*
* No guidance in NFG, thus AECOM professional judgment was used		

**ICV Recovery Nonconformances:**

Nonconformance	Actions	
	Detected Compounds	Nondetected Compounds
%R > 120%	J	No qualification
20% < %R < 80%	J	UJ
%R < 20% (see note)	J	R*

Notes: Based on NFG 2008 VOC guidance, professional judgment is used to reject (R) nondetects in all associated samples for any analyte with < 20% recovery. Also, professional judgment is used to estimate (UJ) rather the reject (R) sample results previously negated (U) on the basis of blank contamination.

**CCV Linearity Nonconformances:**

Nonconformance	Actions	
	Detected Results	Nondetected Results
%D > 20%	J	UJ
%Drift > 20%	J*	UJ*
* No guidance in NFG, thus professional judgment was used		

Qualified sample results are shown in Table 1. Nonconformances are summarized in Attachment A in Table A-1a, A-1b, and A-1c.

**Laboratory Blanks/Equipment Blanks/Trip Blanks**

Laboratory method blanks, equipment rinsate and trip 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, equipment rinsate and trip 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.



### **Surrogate Spike Recoveries**

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

### **MS/MSD Results**

MS/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. The 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.

### **Internal Standard Results**

The internal standard (IS) recoveries were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

### **Sample Results/Reporting Issues**

Compounds that were not detected in the sample are reported as undetected (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**

Sample results qualified as a result of validation actions are summarized in Table 1. All actions are described above.

**ATTACHMENTS**

Attachment A: Nonconformance Summary Tables

Attachment B: Qualifier Codes and Explanations

Attachment C: Reason Codes and Explanations

**Table 1 - Data Validation Summary of Qualified Data**

Sample ID	Matrix	Compound	Result	LOD	Units	Validation Qualifiers	Validation Reason
VPB152-GW-062014-830-832	WG	2-BUTANONE		2.5	UG/L	UJ	c
VPB152-GW-062014-830-832	WG	2-HEXANONE		2.5	UG/L	UJ	c
VPB152-GW-062014-830-832	WG	4-METHYL-2-PENTANONE		2.5	UG/L	UJ	c
VPB152-GW-062014-830-832	WG	ACETONE	10	2.5	UG/L	J	c
VPB152-GW-062014-830-832	WG	CHLOROETHANE		1.0	UG/L	UJ	c
VPB152-GW-062014-830-832	WG	CYCLOHEXANE		0.50	UG/L	UJ	c
VPB152-GW-062014-830-832	WG	DICHLORODIFLUOROMETHANE		1.0	UG/L	UJ	c
VPB152-GW-062014-830-832	WG	TRICHLOROFLUOROMETHANE		1.0	UG/L	UJ	c
VPB152-GW-062014-830-832	WG	VINYL CHLORIDE		1.0	UG/L	UJ	c
VPB152-GW-062314-850-852	WG	2-BUTANONE		2.5	UG/L	UJ	c
VPB152-GW-062314-850-852	WG	2-HEXANONE		2.5	UG/L	UJ	c
VPB152-GW-062314-850-852	WG	4-METHYL-2-PENTANONE		2.5	UG/L	UJ	c
VPB152-GW-062314-850-852	WG	ACETONE	7.7	2.5	UG/L	J	c
VPB152-GW-062314-850-852	WG	CHLOROETHANE		1.0	UG/L	UJ	c
VPB152-GW-062314-850-852	WG	CYCLOHEXANE		0.50	UG/L	UJ	c
VPB152-GW-062314-850-852	WG	DICHLORODIFLUOROMETHANE		1.0	UG/L	UJ	c
VPB152-GW-062314-850-852	WG	TRICHLOROFLUOROMETHANE		1.0	UG/L	UJ	c
VPB152-GW-062314-850-852	WG	VINYL CHLORIDE		1.0	UG/L	UJ	c
VPB152-TRIP BLANK-062314	WQ	2-BUTANONE		2.5	UG/L	UJ	c
VPB152-TRIP BLANK-062314	WQ	2-HEXANONE		2.5	UG/L	UJ	c
VPB152-TRIP BLANK-062314	WQ	4-METHYL-2-PENTANONE		2.5	UG/L	UJ	c
VPB152-TRIP BLANK-062314	WQ	CHLOROETHANE		1.0	UG/L	UJ	c
VPB152-TRIP BLANK-062314	WQ	CYCLOHEXANE		0.50	UG/L	UJ	c
VPB152-TRIP BLANK-062314	WQ	DICHLORODIFLUOROMETHANE		1.0	UG/L	UJ	c
VPB152-TRIP BLANK-062314	WQ	TRICHLOROFLUOROMETHANE		1.0	UG/L	UJ	c
VPB152-TRIP BLANK-062314	WQ	VINYL CHLORIDE		1.0	UG/L	UJ	c

**Attachment A****Nonconformance Summary Tables****Table A-1a - Initial Calibration**

ICAL	Compound	% RSD	Limit
WG144357-7 GCMS-C	CHLOROETHANE	36.2	< 15%
Associated samples: All samples in the SDG			

**Table A-1b - Initial Calibration Verification**

ICV	Compound	% R	Limit
WG144357-7 GCMS-C	ACETONE	120.4	80-120%
Associated samples: All samples in the SDG			

**Table A-1c - Continuing Calibration Verification**

CCV ID	Compound	% D	Limit
WG145394 6/25/2014 GCMS-C	DICHLORODIFLUOROMETHANE	38.0	< 20%
	VINYL CHLORIDE	20.4	< 20%
	TRICHLOROFLUOROMETHANE	33.2	< 20%
	2-BUTANONE	40.0	< 20%
	CYCLOHEXANE	20.6	< 20%
	4-METHYL-2-PENTANONE	46.4	< 20%
	2-HEXANONE	27.9	< 20%
Associated samples: All samples in the SDG			

**Attachment B****Qualifier Codes and Explanations**

<b>Qualifier</b>	<b>Explanation</b>
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.

## Attachment C

### Reason Codes and Explanations

Reason Code	Explanation
be	Equipment blank contamination
bf	Field blank contamination
bl	Laboratory blank contamination
c	Calibration issue
co	Analyte carryover
d	Reporting limit raised due to chromatographic interference
fd	Field duplicate RPDs
h	Holding times
i	Internal standard areas
k	Estimated Maximum Possible Concentration (EMPC)
l	LCS or OPR recoveries
lc	Labeled compound recovery
ld	Laboratory duplicate RPDs
lp	Laboratory control sample/laboratory control sample duplicate RPDs
m	Matrix spike recovery
md	Matrix spike/matrix spike duplicate RPDs
nb	Negative laboratory blank contamination
p	Chemical preservation issue
r	Dual column RPD
q	Quantitation issue
s	Surrogate recovery
su	Ion suppression
t	Temperature preservation issue
x	Percent solids
y	Serial dilution results
z	ICS results

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4577-1  
**Client ID:** 152-062014-830-832  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4577  
**Lab File ID:** C7913.D

**Sample Date:** 20-JUN-14  
**Received Date:** 24-JUN-14  
**Extract Date:** 25-JUN-14  
**Extracted By:** REC  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG145394

**Analysis Date:** 25-JUN-14  
**Analyst:** REC  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 26-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	<del>U</del> UJ	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	<del>U</del> UJ	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	<del>U</del> UJ	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	<del>U</del> UJ	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
Acetone	J	10	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	<del>U</del> UJ	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	<del>U</del> UJ	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene		3.5	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	<del>U</del> UJ	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	<del>U</del> UJ	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50

*REC 6/14/14*

### Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4577-1  
**Client ID:** 152-062014-830-832  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4577  
**Lab File ID:** C7913.D

**Sample Date:** 20-JUN-14  
**Received Date:** 24-JUN-14  
**Extract Date:** 25-JUN-14  
**Extracted By:** REC  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG145394

**Analysis Date:** 25-JUN-14  
**Analyst:** REC  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 26-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylenes	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		87.1	%					
Toluene-d8		89.4	%					
1,2-Dichloroethane-d4		112.	%					
Dibromofluoromethane		98.4	%					



### Report of Analytical Results

Client: ENSAFE  
 Lab ID: SH4577-2  
 Client ID: 152-TB-062314  
 Project: Navy Clean WE15-03-06 NW  
 SDG: SH4577  
 Lab File ID: C7904.D

Sample Date: 23-JUN-14  
 Received Date: 24-JUN-14  
 Extract Date: 25-JUN-14  
 Extracted By: REC  
 Extraction Method: SW846 5030  
 Lab Prep Batch: WG145394

Analysis Date: 25-JUN-14  
 Analyst: REC  
 Analysis Method: SW846 8260B  
 Matrix: AQ  
 % Solids: NA  
 Report Date: 26-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	<del>U</del> UJ	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	<del>U</del> UJ	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	<del>U</del> UJ	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	<del>U</del> UJ	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
Acetone	U	2.5	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	<del>U</del> UJ	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	<del>U</del> UJ	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	<del>U</del> UJ	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	<del>U</del> UJ	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50



### Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4577-2  
**Client ID:** 152-TB-062314  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4577  
**Lab File ID:** C7904.D

**Sample Date:** 23-JUN-14  
**Received Date:** 24-JUN-14  
**Extract Date:** 25-JUN-14  
**Extracted By:** REC  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG145394

**Analysis Date:** 25-JUN-14  
**Analyst:** REC  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 26-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		87.9	%					
Toluene-d8		90.5	%					
1,2-Dichloroethane-d4		117.	%					
Dibromofluoromethane		100.	%					

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4577-3  
**Client ID:** 152-062314-850-852  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4577  
**Lab File ID:** C7914.D

**Sample Date:** 23-JUN-14  
**Received Date:** 24-JUN-14  
**Extract Date:** 25-JUN-14  
**Extracted By:** REC  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG145394

**Analysis Date:** 25-JUN-14  
**Analyst:** REC  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 26-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	<del>U</del> UJ	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	<del>U</del> UJ	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	<del>U</del> UJ	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	<del>U</del> UJ	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
Acetone	J	7.7	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	<del>U</del> UJ	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	<del>U</del> UJ	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	<del>U</del> UJ	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	<del>U</del> UJ	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4577-3  
**Client ID:** 152-062314-850-852  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4577  
**Lab File ID:** C7914.D

**Sample Date:** 23-JUN-14  
**Received Date:** 24-JUN-14  
**Extract Date:** 25-JUN-14  
**Extracted By:** REC  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG145394

**Analysis Date:** 25-JUN-14  
**Analyst:** REC  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 26-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		87.3	%					
Toluene-d8		88.1	%					
1,2-Dichloroethane-d4		117.	%					
Dibromofluoromethane		98.7	%					



## Data Validation Report

Project: Regional Groundwater Investigation - NWIRP Bethpage  
Laboratory: Katahdin Analytical Services, Scarborough, Maine  
Service Request: SH4477  
Analyses/Method: EPA SW-846 Method 8260B for VOCs (GC/MS)  
Validation Level: Limited  
RESCON Project 60266526.SA.DV  
Number:  
Prepared by: Sheena Blair/RESCON Completed on: 9/10/2014  
Reviewed by: Lori Herberich/RESCON File Name: SH4477\_8260B

### SUMMARY

The samples listed below were collected by Resolution Consultants (RESCON) from the Regional Groundwater Investigation - NWIRP Bethpage site on June 17 and 19, 2014.

Sample ID	Matrix/Sample Type
VPB152-GW-061714-725-727	Groundwater
VPB152-GW-061714-740-742	Groundwater
VPB152-GW-061914-780-782	Groundwater
VPB152-GW-061914-800-802	Groundwater
VPB152-TRIP BLANK-061914	Trip Blank

Data validation activities were conducted with reference to *Test Methods for Evaluating Solid Waste, Physical/Chemical Methods SW-846, specifically SW-846 Method 8260B, Volatile Organic Compounds by Gas Chromatography/Mass Spectrometry* (USEPA, 1996), *USEPA Contract Laboratory Program National Functional Guidelines for Superfund Organic Methods Data Review* (June 2008), and *Quality Systems Manual (QSM) for Environmental Laboratories, Version 4.2* (DoD, 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 review elements (where applicable to the method):

- X Data completeness (chain-of-custody [COC])/sample integrity
- ✓ Holding times and sample preservation
- ✓ GC/MS performance checks
- X Initial calibration/continuing calibration verification
- X Laboratory blanks/equipment blanks/trip blanks
- ✓ Surrogate spike recoveries
- NA Matrix spike (MS) and/or matrix spike duplicate (MSD) results
- ✓ Laboratory control sample (LCS) results

- NA Field duplicate results
- ✓ Internal standard results
- ✓ 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 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. Selected data points were estimated or negated due to nonconformances of certain QC criteria (see discussion below). Qualified sample results are presented in Table 1.

## 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 "VPB" prefix and the "GW" from the sample ID, and truncated the ID for the Trip Blank in the report. The submitted EDD file reflects the full sample ID.

Sample VPB152-GW-061714-725-727 was extremely silty and had very little standing water. The laboratory decanted the water from all three individual vials and made a composite of the sample. As a result sample VPB152-GW-061714-725-727 was analyzed at a 50-fold dilution. Positive and nondetect results for this sample were qualified as estimated (J and UJ) respectively, due to possible loss of sample integrity during the compositing procedure. Qualified sample results are presented in Table 1.

### 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.

### GC/MS Performance Checks

The data were reviewed to ensure that the 4-bromofluorobenzene (BFB) tuning was performed at the correct frequency and that the method acceptance criteria were met. 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:

- the initial calibration (ICAL) percent relative standard deviation (%RSD), correlation coefficient (r)/coefficient of determination ( $r^2$ ), and/or response factor method acceptance criteria were met;
- the initial calibration verification standard (ICV) percent recoveries (%Rs) acceptance criteria were met;
- the continuing calibration verification standard (CCV) method percent difference or percent drift (%Ds) and RF acceptance criteria were met; and
- the retention time method acceptance criteria were met.

Data qualification to the analytes associated with the specific ICAL, ICV and/or CCV was as follows:

#### ICAL Linearity Nonconformances:

Nonconformance	Actions	
	Detected	Nondetected Results
%RSD > 15% and quantitation based on mean RF	J	UJ
r or $r^2$ < 0.99 and quantitation based on linear	J*	UJ*
* No guidance in NFG, thus professional judgment was used		

#### ICV Recovery Nonconformances:

Nonconformance	Actions	
	Detected Compounds	Nondetected Compounds
%R > 120%	J	No qualification
20% < %R < 80%	J	UJ
%R < 20% (see note)	J	R*

Notes: Based on NFG 2008 VOC guidance, professional judgment is used to reject (R) nondetects in all associated samples for any analyte with < 20% recovery. Also, professional judgment is used to estimate (UJ) rather the reject (R) sample results previously negated (U) on the basis of blank contamination.

#### CCV Linearity Nonconformances:

Nonconformance	Actions	
	Detected Results	Nondetected Results
%D > 20%	J	UJ
%Drift > 20%	J*	UJ*
* No guidance in NFG, thus professional judgment was used		

Qualified sample results are shown in Table 1. Nonconformances are summarized in Attachment A in Table A-1a, A-1b, and A-1c.

#### Laboratory Blanks/Equipment Blanks/Trip Blanks

Laboratory method blanks, equipment rinsate and trip 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, equipment rinsate and trip 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.

Sample results were qualified as follows:

Blank type	Blank result	Sample result	Action for samples
Method, Storage, Field, Trip, or Instrument*	Detects	Not detected	No qualification
	≤ LOQ	< LOQ	Report sample LOQ value with a U
		≥ LOQ and < 2x LOQ	Report the sample result with a U**
		≥ 2x the LOQ	No qualifications
	> LOQ	< LOQ	Report sample LOQ value with a U
		≥ LOQ and < blank contamination	Report the sample result with a U or reject the sample result as unusable R
		≥ LOQ and ≥ blank contamination	If the result is ≤ 2x blank result, report the sample result U.** If the result is > 2x blank result, no qualification is required.**
TIC detected	Detects	If the result is ≤ 2x blank result, report the sample result U.** If the result is > 2x blank result, no qualification is required.**	
* Qualifications based on instrument blank results affect only the sample analyzed immediately after the sample that has target compounds that exceed the calibration range or non-target compounds that exceed 100 g/L.			
**Based on professional judgment.			

LOQ - Limit of Quantitation

Nonconformances are summarized in Attachment A in Table A-2. Qualified sample results are shown in Table 1.

### **Surrogate Spike Recoveries**

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

### **MS/MSD Results**

MS/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. The 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.

**Internal Standard Results**

The internal standard (IS) recoveries were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

**Sample Results/Reporting Issues**

Compounds that were not detected in the sample are reported as undetected (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**

Sample results qualified as a result of validation actions are summarized in Table 1. All actions are described above.

**ATTACHMENTS**

Attachment A: Nonconformance Summary Tables

Attachment B: Qualifier Codes and Explanations

Attachment C: Reason Codes and Explanations

**Table 1 - Data Validation Summary of Qualified Data**

Sample ID	Matrix	Compound	Result	LOD	Units	Validation Qualifiers	Validation Reason
VPB152-GW-061714-725-727	WG	1,1,1-TRICHLOROETHANE		25	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	1,1,2,2-TETRACHLOROETHANE		25	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE		25	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	1,1,2-TRICHLOROETHANE		25	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	1,1-DICHLOROETHANE		25	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	1,1-DICHLOROETHENE		25	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	1,2,4-TRICHLOROBENZENE		25	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	1,2-DIBROMO-3-CHLOROPROPANE		38	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	1,2-DIBROMOETHANE		25	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	1,2-DICHLOROBENZENE		25	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	1,2-DICHLOROETHANE		25	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	1,2-DICHLOROETHENE, TOTAL		50	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	1,2-DICHLOROPROPANE		25	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	1,3-DICHLOROBENZENE		25	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	1,4-DICHLOROBENZENE		25	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	2-BUTANONE		120	UG/L	UJ	c,mc
VPB152-GW-061714-725-727	WG	2-HEXANONE		120	UG/L	UJ	c,mc
VPB152-GW-061714-725-727	WG	4-METHYL-2-PENTANONE		120	UG/L	UJ	c,mc
VPB152-GW-061714-725-727	WG	ACETONE		120	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	BENZENE		25	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	BROMODICHLOROMETHANE		25	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	BROMOFORM		25	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	BROMOMETHANE		50	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	CARBON DISULFIDE		25	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	CARBON TETRACHLORIDE		25	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	CHLOROBENZENE		25	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	CHLOROETHANE		50	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	CHLOROFORM		25	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	CHLOROMETHANE		50	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	CIS-1,2-DICHLOROETHENE		25	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	CIS-1,3-DICHLOROPROPENE		25	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	CYCLOHEXANE		25	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	DIBROMOCHLOROMETHANE		25	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	DICHLORODIFLUOROMETHANE		50	UG/L	UJ	c,mc
VPB152-GW-061714-725-727	WG	ETHYLBENZENE		25	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	ISOPROPYLBENZENE		25	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	M- AND P-XYLENE		50	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	METHYL ACETATE		38	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	METHYL CYCLOHEXANE		25	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	METHYL TERT-BUTYL ETHER		25	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	METHYLENE CHLORIDE	67	120	UG/L	J	mc
VPB152-GW-061714-725-727	WG	O-XYLENE		25	UG/L	UJ	mc

Sample ID	Matrix	Compound	Result	LOD	Units	Validation Qualifiers	Validation Reason
VPB152-GW-061714-725-727	WG	STYRENE		25	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	TETRACHLOROETHENE		25	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	TOLUENE		25	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	TRANS-1,2-DICHLOROETHENE		25	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	TRANS-1,3-DICHLOROPROPENE		25	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	TRICHLOROETHENE		25	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	TRICHLOROFLUOROMETHANE		50	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	VINYL CHLORIDE		50	UG/L	UJ	mc
VPB152-GW-061714-725-727	WG	XYLENES, TOTAL		75	UG/L	UJ	mc
VPB152-GW-061714-740-742	WG	2-BUTANONE	2.1	2.5	UG/L	J	c
VPB152-GW-061714-740-742	WG	2-HEXANONE		2.5	UG/L	UJ	c
VPB152-GW-061714-740-742	WG	4-METHYL-2-PENTANONE		2.5	UG/L	UJ	c
VPB152-GW-061714-740-742	WG	ACETONE	11	2.5	UG/L	J	c
VPB152-GW-061714-740-742	WG	DICHLORODIFLUOROMETHANE		1.0	UG/L	UJ	c
VPB152-GW-061914-780-782	WG	2-BUTANONE		2.5	UG/L	UJ	c
VPB152-GW-061914-780-782	WG	2-HEXANONE		2.5	UG/L	UJ	c
VPB152-GW-061914-780-782	WG	4-METHYL-2-PENTANONE		2.5	UG/L	UJ	c
VPB152-GW-061914-780-782	WG	ACETONE	7.5	2.5	UG/L	J	c
VPB152-GW-061914-780-782	WG	DICHLORODIFLUOROMETHANE		1.0	UG/L	UJ	c
VPB152-GW-061914-800-802	WG	2-BUTANONE		2.5	UG/L	UJ	c
VPB152-GW-061914-800-802	WG	2-HEXANONE		2.5	UG/L	UJ	c
VPB152-GW-061914-800-802	WG	4-METHYL-2-PENTANONE		2.5	UG/L	UJ	c
VPB152-GW-061914-800-802	WG	ACETONE	11	2.5	UG/L	J	c
VPB152-GW-061914-800-802	WG	CARBON DISULFIDE		1.0*	UG/L	U	bf
VPB152-GW-061914-800-802	WG	DICHLORODIFLUOROMETHANE		1.0	UG/L	UJ	c
VPB152-TRIP BLANK-061914	WQ	2-BUTANONE		2.5	UG/L	UJ	c
VPB152-TRIP BLANK-061914	WQ	2-HEXANONE		2.5	UG/L	UJ	c
VPB152-TRIP BLANK-061914	WQ	4-METHYL-2-PENTANONE		2.5	UG/L	UJ	c
VPB152-TRIP BLANK-061914	WQ	DICHLORODIFLUOROMETHANE		1.0	UG/L	UJ	c
*LOQ							

## Attachment A

## Nonconformance Summary Tables

Table A-1a - Initial Calibration

ICAL	Compound	% RSD	Limit
WG144357-7 GCMS-C	CHLOROETHANE	36.2	< 15%
Associated samples: All samples in the SDG			

Table A-1b - Initial Calibration Verification

ICV	Compound	% R	Limit
WG144357-7 GCMS-C	ACETONE	120.4	80-120%
Associated samples: All samples in the SDG			

Table A-1c - Continuing Calibration Verification

CCV ID	Compound	% D	Limit
WG145087 6/20/2014 GCMS-C	DICHLORODIFLUOROMETHANE	60.6	< 20%
	2-BUTANONE	23.9	< 20%
	4-METHYL-2-PENTANONE	40.2	< 20%
	2-HEXANONE	21.1	< 20%
Associated samples: All samples in the SDG			

**Attachment B****Qualifier Codes and Explanations**

<b>Qualifier</b>	<b>Explanation</b>
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.

## Attachment C

## Reason Codes and Explanations

Reason Code	Explanation
be	Equipment blank contamination
bf	Field blank contamination
bl	Laboratory blank contamination
c	Calibration issue
co	Analyte carryover
d	Reporting limit raised due to chromatographic interference
fd	Field duplicate RPDs
h	Holding times
i	Internal standard areas
k	Estimated Maximum Possible Concentration (EMPC)
l	LCS or OPR recoveries
lc	Labeled compound recovery
ld	Laboratory duplicate RPDs
lp	Laboratory control sample/laboratory control sample duplicate RPDs
m	Matrix spike recovery
md	Matrix spike/matrix spike duplicate RPDs
nb	Negative laboratory blank contamination
p	Chemical preservation issue
r	Dual column RPD
q	Quantitation Issue
s	Surrogate recovery
su	Ion suppression
t	Temperature preservation issue
x	Percent solids
y	Serial dilution results
z	ICS results

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4477-1DL  
**Client ID:** 152-061714-725-727  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4477  
**Lab File ID:** C7780.D

**Sample Date:** 17-JUN-14  
**Received Date:** 20-JUN-14  
**Extract Date:** 20-JUN-14  
**Extracted By:** REC  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG145087

**Analysis Date:** 20-JUN-14  
**Analyst:** REC  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 23-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	50	ug/L	50	2	100	12.	50.
Chloromethane	U	50	ug/L	50	2	100	18.	50.
Vinyl Chloride	U	50	ug/L	50	2	100	12.	50.
Bromomethane	U	50	ug/L	50	2	100	24.	50.
Chloroethane	U	50	ug/L	50	2	100	28.	50.
Trichlorofluoromethane	U	50	ug/L	50	2	100	12.	50.
1,1-Dichloroethene	U	25	ug/L	50	1	50.	18.	25.
Carbon Disulfide	U	25	ug/L	50	1	50.	12.	25.
Freon-113	U	25	ug/L	50	1	50.	16.	25.
<b>Methylene Chloride</b>	J	67	ug/L	50	5	250	56.	120
Acetone	U	120	ug/L	50	5	250	110	120
trans-1,2-Dichloroethene	U	25	ug/L	50	1	50.	12.	25.
Methyl tert-butyl Ether	U	25	ug/L	50	1	50.	18.	25.
1,1-Dichloroethane	U	25	ug/L	50	1	50.	10.	25.
cis-1,2-Dichloroethene	U	25	ug/L	50	1	50.	10.	25.
Chloroform	U	25	ug/L	50	1	50.	16.	25.
1,1,1-Trichloroethane	U	25	ug/L	50	1	50.	10.	25.
2-Butanone	U	120	ug/L	50	5	250	66.	120
Cyclohexane	U	25	ug/L	50	1	50.	16.	25.
Carbon Tetrachloride	U	25	ug/L	50	1	50.	11.	25.
Benzene	U	25	ug/L	50	1	50.	13.	25.
1,2-Dichloroethane	U	25	ug/L	50	1	50.	10.	25.
Trichloroethene	U	25	ug/L	50	1	50.	14.	25.
1,2-Dichloropropane	U	25	ug/L	50	1	50.	12.	25.
Bromodichloromethane	U	25	ug/L	50	1	50.	16.	25.
cis-1,3-Dichloropropene	U	25	ug/L	50	1	50.	9.5	25.
Toluene	U	25	ug/L	50	1	50.	14.	25.
4-Methyl-2-Pentanone	U	120	ug/L	50	5	250	66.	120
trans-1,3-Dichloropropene	U	25	ug/L	50	1	50.	10.	25.
1,1,2-Trichloroethane	U	25	ug/L	50	1	50.	16.	25.
Tetrachloroethene	U	25	ug/L	50	1	50.	20.	25.
Dibromochloromethane	U	25	ug/L	50	1	50.	15.	25.
2-Hexanone	U	120	ug/L	50	5	250	85.	120
Chlorobenzene	U	25	ug/L	50	1	50.	11.	25.
Ethylbenzene	U	25	ug/L	50	1	50.	10.	25.

*Handwritten signature and date: J. 4/13/14*

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4477-1DL  
**Client ID:** 152-061714-725-727  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4477  
**Lab File ID:** C7780.D

**Sample Date:** 17-JUN-14  
**Received Date:** 20-JUN-14  
**Extract Date:** 20-JUN-14  
**Extracted By:** REC  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG145087

**Analysis Date:** 20-JUN-14  
**Analyst:** REC  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 23-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U UJ	75	ug/L	50	3	150	12.	75.
Styrene	U	25	ug/L	50	1	50.	12.	25.
Bromoform	U	25	ug/L	50	1	50.	12.	25.
Isopropylbenzene	U	25	ug/L	50	1	50.	12.	25.
1,1,2,2-Tetrachloroethane	U	25	ug/L	50	1	50.	19.	25.
1,3-Dichlorobenzene	U	25	ug/L	50	1	50.	13.	25.
1,4-Dichlorobenzene	U	25	ug/L	50	1	50.	12.	25.
1,2-Dichlorobenzene	U	25	ug/L	50	1	50.	7.5	25.
1,2,4-Trichlorobenzene	U	25	ug/L	50	1	50.	18.	25.
Methyl Acetate	U	38	ug/L	50	1	50.	26.	38.
Methylcyclohexane	U	25	ug/L	50	1	50.	15.	25.
o-Xylene	U	25	ug/L	50	1	50.	12.	25.
M+P-Xylenes	U	50	ug/L	50	2	100	30.	50.
1,2-Dichloroethylene (Total)	U	50	ug/L	50	2	100	10.	50.
1,2-Dibromoethane	U	25	ug/L	50	1	50.	11.	25.
1,2-Dibromo-3-Chloropropane	U	38	ug/L	50	1	50.	25.	38.
P-Bromofluorobenzene		94.6	%					
Toluene-d8		96.1	%					
1,2-Dichloroethane-d4		111.	%					
Dibromofluoromethane		97.9	%					

*R-11/13/14*



## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4477-2  
**Client ID:** 152-061714-740-742  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4477  
**Lab File ID:** C7787.D

**Sample Date:** 17-JUN-14  
**Received Date:** 20-JUN-14  
**Extract Date:** 20-JUN-14  
**Extracted By:** REC  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG145087

**Analysis Date:** 20-JUN-14  
**Analyst:** REC  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 23-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U <del>U</del> UJ	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
<b>Acetone</b>	J	11	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
<b>2-Butanone</b>	J <del>J</del> J	2.1	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
<b>Toluene</b>		5.6	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	U <del>U</del> UJ	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U <del>U</del> UJ	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50



## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4477-2  
**Client ID:** 152-061714-740-742  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4477  
**Lab File ID:** C7787.D

**Sample Date:** 17-JUN-14  
**Received Date:** 20-JUN-14  
**Extract Date:** 20-JUN-14  
**Extracted By:** REC  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG145087

**Analysis Date:** 20-JUN-14  
**Analyst:** REC  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 23-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		99.3	%					
Toluene-d8		98.3	%					
1,2-Dichloroethane-d4		113.	%					
Dibromofluoromethane		101.	%					

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4477-3  
**Client ID:** 152-061914-780-782  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4477  
**Lab File ID:** C7788.D

**Sample Date:** 19-JUN-14  
**Received Date:** 20-JUN-14  
**Extract Date:** 20-JUN-14  
**Extracted By:** REC  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG145087

**Analysis Date:** 20-JUN-14  
**Analyst:** REC  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 23-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	<del>U</del> UJ	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
<b>Acetone</b>	J	7.5	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	<del>U</del> UJ	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	<del>U</del> UJ	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	<del>U</del> UJ	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50

REC 11/23/14

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4477-3  
**Client ID:** 152-061914-780-782  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4477  
**Lab File ID:** C7788.D

**Sample Date:** 19-JUN-14  
**Received Date:** 20-JUN-14  
**Extract Date:** 20-JUN-14  
**Extracted By:** REC  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG145087

**Analysis Date:** 20-JUN-14  
**Analyst:** REC  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 23-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		101.	%					
Toluene-d8		99.4	%					
1,2-Dichloroethane-d4		117.	%					
Dibromofluoromethane		103.	%					

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4477-4  
**Client ID:** 152-061914-800-802  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4477  
**Lab File ID:** C7789.D

**Sample Date:** 19-JUN-14  
**Received Date:** 20-JUN-14  
**Extract Date:** 20-JUN-14  
**Extracted By:** REC  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG145087

**Analysis Date:** 20-JUN-14  
**Analyst:** REC  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 23-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	<del>U</del> UJ	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
<b>Carbon Disulfide</b>	<del>J</del> U	<del>0.32</del> 1.0	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
<b>Acetone</b>	J	11	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	<del>U</del> UJ	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	<del>U</del> UJ	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	<del>U</del> UJ	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50

REC 6/23/14

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4477-4  
**Client ID:** 152-061914-800-802  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4477  
**Lab File ID:** C7789.D

**Sample Date:** 19-JUN-14  
**Received Date:** 20-JUN-14  
**Extract Date:** 20-JUN-14  
**Extracted By:** REC  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG145087

**Analysis Date:** 20-JUN-14  
**Analyst:** REC  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 23-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		99.2	%					
Toluene-d8		98.3	%					
1,2-Dichloroethane-d4		117.	%					
Dibromofluoromethane		102.	%					

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4477-5  
**Client ID:** VPB152-TB-061914  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4477  
**Lab File ID:** C7779.D

**Sample Date:** 19-JUN-14  
**Received Date:** 20-JUN-14  
**Extract Date:** 20-JUN-14  
**Extracted By:** REC  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG145087

**Analysis Date:** 20-JUN-14  
**Analyst:** REC  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 23-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	<del>U</del> UJ	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
Acetone	U	2.5	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	<del>U</del> UJ	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	<del>U</del> UJ	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	<del>U</del> UJ	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50

*Handwritten signature and date: R 4/13/14*

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4477-5  
**Client ID:** VPB152-TB-061914  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4477  
**Lab File ID:** C7779.D

**Sample Date:** 19-JUN-14  
**Received Date:** 20-JUN-14  
**Extract Date:** 20-JUN-14  
**Extracted By:** REC  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG145087

**Analysis Date:** 20-JUN-14  
**Analyst:** REC  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 23-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		95.0	%					
Toluene-d8		95.4	%					
1,2-Dichloroethane-d4		115.	%					
Dibromofluoromethane		99.6	%					





## Data Validation Report

Project: Regional Groundwater Investigation - NWIRP Bethpage  
Laboratory: Katahdin Analytical Services, Scarborough, Maine  
Service Request: SH4301  
Analyses/Method: EPA SW-846 Method 8260B for VOCs (GC/MS)  
Validation Level: Limited  
RESCON Project 60266526.SA.DV  
Number: \_\_\_\_\_  
Prepared by: Sheena Blair/RESCON Completed on: 9/10/2014  
Reviewed by: Lori Herberich/RESCON File Name: SH4301\_8260B

### SUMMARY

The samples listed below were collected by Resolution Consultants (RESCON) from the Regional Groundwater Investigation - NWIRP Bethpage site on June 13 and 16, 2014.

Sample ID	Matrix/Sample Type
VPB152-GW-061314-620-622	Groundwater
VPB152-GW-061314-640-642	Groundwater
VPB152-GWD-061314	Field Duplicate of VPB152-GW-061314-640-642
VPB152-GW-061614-660-662	Groundwater
VPB152-GW-061614-680-682	Groundwater
VPB152-GW-061614-700-702	Groundwater
VPB152-TRIP BLANK-061614	Trip Blank

Data validation activities were conducted with reference to *Test Methods for Evaluating Solid Waste, Physical/Chemical Methods SW-846, specifically SW-846 Method 8260B, Volatile Organic Compounds by Gas Chromatography/Mass Spectrometry* (USEPA, 1996), *USEPA Contract Laboratory Program National Functional Guidelines for Superfund Organic Methods Data Review* (June 2008), and *Quality Systems Manual (QSM) for Environmental Laboratories, Version 4.2* (DoD, 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 review elements (where applicable to the method):

- X Data completeness (chain-of-custody [COC])/sample integrity
- ✓ Holding times and sample preservation
- ✓ GC/MS performance checks
- X Initial calibration/continuing calibration verification
- X Laboratory blanks/equipment blanks/trip blanks

- ✓ Surrogate spike recoveries
- ✓ Matrix spike (MS) and/or matrix spike duplicate (MSD) results
- ✓ Laboratory control sample (LCS) results
- ✓ Field duplicate results
- ✓ Internal standard results
- ✓ 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 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. Selected data points were estimated or negated due to nonconformances of certain QC criteria (see discussion below). Qualified sample results are presented in Table 1.

## 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 "VPB" prefix and the "GW" from the sample ID, and truncated the ID for the Trip Blank in the report. The submitted EDD file reflects the full sample ID.

Selected samples were mostly soil and had very little standing water.

- For samples VPB152-GW-061314-620-622 and VPB152-GW-061614-700-702 the laboratory decanted the liquid from three vials into one vial as a composite. Sample VPB152-GW-061614-700-702 was analyzed at a 100-fold dilution due to the limited sample volume available.

Positive and nondetect results for these sample were qualified as estimated (J and UJ) respectively, due to possible loss of sample integrity during the decanting procedure. Qualified sample results are presented in Table 1.

### **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.

### **GC/MS Performance Checks**

The data were reviewed to ensure that the 4-bromofluorobenzene (BFB) tuning was performed at the correct frequency and that the method acceptance criteria were met. 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:

- the initial calibration (ICAL) percent relative standard deviation (%RSD), correlation coefficient (r)/coefficient of determination ( $r^2$ ), and/or response factor method acceptance criteria were met;
- the initial calibration verification standard (ICV) percent recoveries (%Rs) acceptance criteria were met;
- the continuing calibration verification standard (CCV) method percent difference or percent drift (%Ds) and RF acceptance criteria were met; and
- the retention time method acceptance criteria were met.

Data qualification to the analytes associated with the specific ICAL, ICV, and/or CCV was as follows:

#### **ICV Recovery Nonconformances:**

Nonconformance	Actions	
	Detected Compounds	Nondetected Compounds
%R > 120%	J	No qualification
20% < %R < 80%	J	UJ
%R < 20% (see note)	J	R*

Notes: Based on NFG 2008 VOC guidance, professional judgment is used to reject (R) nondetects in all associated samples for any analyte with < 20% recovery. Also, professional judgment is used to estimate (UJ) rather the reject (R) sample results previously negated (U) on the basis of blank contamination.

#### **CCV Linearity Nonconformances:**

Nonconformance	Actions	
	Detected Results	Nondetected Results
%D > 20%	J	UJ
%Drift > 20%	J*	UJ*
* No guidance in NFG, thus professional judgment was used		

Qualified sample results are shown in Table 1. Nonconformances are summarized in Attachment A in Table A-1a, A-1b, A-1c, and A-1d.

### **Laboratory Blanks/Equipment Blanks/Trip Blanks**

Laboratory method blanks, equipment rinsate and trip 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, equipment rinsate and trip 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.

Sample results were qualified as follows:

For common lab contaminants (methylene chloride, acetone, 2-butanone):

Blank type	Blank result	Sample result	Action for samples
Method, Storage, Field, Trip, or Instrument*	Detects	Not detected	No qualification
	≤ 2x LOQ	< 2x LOQ	Report sample LOQ value with a U
		≥ 2x LOQ and ≤ 4x LOQ	Report the sample result with a U**
		≥ 4x LOQ	No qualifications
	> 2x LOQ	< 2x LOQ	Report sample LOQ value with a U
		≥ 2x LOQ and < blank contamination	Report the sample result with a U
		≥ 2x LOQ and ≥ blank contamination	If the result is ≤ 2x blank result, report the sample result U.** If the result is > 2x blank result, no qualification is required.**
* Qualifications based on instrument blank results affect only the sample analyzed immediately after the sample that has target compounds that exceed the calibration range or non-target compounds that exceed 100 µg/L.			
**Based on professional judgment			

For all other compounds:

Blank type	Blank result	Sample result	Action for samples
Method, Storage, Field, Trip, or Instrument*	Detects	Not detected	No qualification
	≤ LOQ	< LOQ	Report sample LOQ value with a U
		≥ LOQ and ≤ 2x LOQ	Report the sample result with a U**
		≥ 2x the LOQ	No qualifications
	> LOQ	< LOQ	Report sample LOQ value with a U
		≥ LOQ and < blank contamination	Report the sample result with a U or reject the sample result as unusable R
		≥ LOQ and ≥ blank contamination	If the result is ≤ 2x blank result, report the sample result U.** If the result is > 2x blank result, no qualification is required.**
	TIC detected	Detects	If the result is ≤ 2x blank result, report the sample result U.** If the result is > 2x blank result, no qualification is required.**
* Qualifications based on instrument blank results affect only the sample analyzed immediately after the sample that has target compounds that exceed the calibration range or non-target compounds that exceed 100 g/L.			
**Based on professional judgment.			

LOQ - Limit of Quantitation

Nonconformances are summarized in Attachment A in Table A-2. Qualified sample results are shown in Table 1.

### **Surrogate Spike Recoveries**

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

### **MS/MSD Results**

The MS/MSD %Rs and relative percent differences (RPDs) were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

### **LCS Results**

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

Although selected compounds had high %Rs, no validation action was required because the associated results were nondetect.

### **Field Duplicate Results**

Field duplicate RPDs were reviewed for conformance with the QC criterion of  $\leq 30\%$  for aqueous matrices. This criterion applies if both results were greater than five times the Limit of Quantitation (LOQ). All QC acceptance criteria were met.

### **Internal Standard Results**

The internal standard (IS) recoveries were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

### **Sample Results/Reporting Issues**

Compounds that were not detected in the sample are reported as undetected (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**

Sample results qualified as a result of validation actions are summarized in Table 1. All actions are described above.

**ATTACHMENTS**

Attachment A: Nonconformance Summary Tables

Attachment B: Qualifier Codes and Explanations

Attachment C: Reason Codes and Explanations

**Table 1 - Data Validation Summary of Qualified Data**

Sample ID	Matrix	Compound	Result	LOD	Units	Validation Qualifiers	Validation Reason
VPB152-GW-061314-620-622	WG	1,1,1-TRICHLOROETHANE		0.50	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	1,1,2,2-TETRACHLOROETHANE		0.50	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE		0.50	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	1,1,2-TRICHLOROETHANE		0.50	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	1,1-DICHLOROETHANE		0.50	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	1,1-DICHLOROETHENE		0.50	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	1,2,4-TRICHLOROBENZENE		0.50	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	1,2-DIBROMO-3-CHLOROPROPANE		0.75	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	1,2-DIBROMOETHANE		0.50	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	1,2-DICHLOROBENZENE		0.50	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	1,2-DICHLOROETHANE		0.50	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	1,2-DICHLOROETHENE, TOTAL		1.0	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	1,2-DICHLOROPROPANE		0.50	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	1,3-DICHLOROBENZENE		0.50	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	1,4-DICHLOROBENZENE		0.50	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	2-BUTANONE		2.5	UG/L	UJ	c,mc
VPB152-GW-061314-620-622	WG	2-HEXANONE		2.5	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	4-METHYL-2-PENTANONE		2.5	UG/L	UJ	c,mc
VPB152-GW-061314-620-622	WG	ACETONE	14	2.5	UG/L	J	c,mc
VPB152-GW-061314-620-622	WG	BENZENE		0.50	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	BROMODICHLOROMETHANE		0.50	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	BROMOFORM		0.50	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	BROMOMETHANE		1.0	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	CARBON DISULFIDE	0.40	0.50	UG/L	J	mc
VPB152-GW-061314-620-622	WG	CARBON TETRACHLORIDE		0.50	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	CHLOROBENZENE		0.50	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	CHLOROETHANE		1.0	UG/L	UJ	c,mc
VPB152-GW-061314-620-622	WG	CHLOROFORM		0.50	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	CHLOROMETHANE		1.0	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	CIS-1,2-DICHLOROETHENE		0.50	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	CIS-1,3-DICHLOROPROPENE		0.50	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	CYCLOHEXANE		0.50	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	DIBROMOCHLOROMETHANE		0.50	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	DICHLORODIFLUOROMETHANE		1.0	UG/L	UJ	c,mc
VPB152-GW-061314-620-622	WG	ETHYLBENZENE		0.50	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	ISOPROPYLBENZENE		0.50	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	M- AND P-XYLENE		1.0	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	METHYL ACETATE		0.75	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	METHYL CYCLOHEXANE		0.50	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	METHYL TERT-BUTYL ETHER		0.50	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	METHYLENE CHLORIDE		2.5	UG/L	UJ	mc

Sample ID	Matrix	Compound	Result	LOD	Units	Validation Qualifiers	Validation Reason
VPB152-GW-061314-620-622	WG	O-XYLENE		0.50	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	STYRENE		0.50	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	TETRACHLOROETHENE		0.50	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	TOLUENE	3.7	0.50	UG/L	J	mc
VPB152-GW-061314-620-622	WG	TRANS-1,2-DICHLOROETHENE		0.50	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	TRANS-1,3-DICHLOROPROPENE		0.50	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	TRICHLOROETHENE		0.50	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	TRICHLOROFLUOROMETHANE		1.0	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	VINYL CHLORIDE		1.0	UG/L	UJ	mc
VPB152-GW-061314-620-622	WG	XYLENES, TOTAL		1.5	UG/L	UJ	mc
VPB152-GW-061314-640-642	WG	2-BUTANONE		2.5	UG/L	UJ	c
VPB152-GW-061314-640-642	WG	4-METHYL-2-PENTANONE		2.5	UG/L	UJ	c
VPB152-GW-061314-640-642	WG	ACETONE	7.0	2.5	UG/L	J	c
VPB152-GW-061314-640-642	WG	CARBON DISULFIDE		1.0*	UG/L	U	bl
VPB152-GW-061314-640-642	WG	CHLOROETHANE		1.0	UG/L	UJ	c
VPB152-GW-061314-640-642	WG	DICHLORODIFLUOROMETHANE		1.0	UG/L	UJ	c
VPB152-GW-061614-660-662	WG	2-BUTANONE		2.5	UG/L	UJ	c
VPB152-GW-061614-660-662	WG	4-METHYL-2-PENTANONE		2.5	UG/L	UJ	c
VPB152-GW-061614-660-662	WG	ACETONE	8.4	2.5	UG/L	J	c
VPB152-GW-061614-660-662	WG	CHLOROETHANE		1.0	UG/L	UJ	c
VPB152-GW-061614-660-662	WG	DICHLORODIFLUOROMETHANE		1.0	UG/L	UJ	c
VPB152-GW-061614-680-682	WG	2-BUTANONE	4.8	2.5	UG/L	J	c
VPB152-GW-061614-680-682	WG	4-METHYL-2-PENTANONE		2.5	UG/L	UJ	c
VPB152-GW-061614-680-682	WG	ACETONE	22	2.5	UG/L	J	c
VPB152-GW-061614-680-682	WG	CARBON DISULFIDE		1.0*	UG/L	U	bl
VPB152-GW-061614-680-682	WG	CHLOROETHANE		1.0	UG/L	UJ	c
VPB152-GW-061614-680-682	WG	DICHLORODIFLUOROMETHANE		1.0	UG/L	UJ	c
VPB152-GW-061614-700-702	WG	1,1,1-TRICHLOROETHANE		50	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	1,1,2,2-TETRACHLOROETHANE		50	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE		50	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	1,1,2-TRICHLOROETHANE		50	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	1,1-DICHLOROETHANE		50	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	1,1-DICHLOROETHENE		50	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	1,2,4-TRICHLOROBENZENE		50	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	1,2-DIBROMO-3-CHLOROPROPANE		75	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	1,2-DIBROMOETHANE		50	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	1,2-DICHLOROBENZENE		50	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	1,2-DICHLOROETHANE		50	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	1,2-DICHLOROETHENE, TOTAL		100	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	1,2-DICHLOROPROPANE		50	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	1,3-DICHLOROBENZENE		50	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	1,4-DICHLOROBENZENE		50	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	2-BUTANONE		250	UG/L	UJ	c,mc
VPB152-GW-061614-700-702	WG	2-HEXANONE		250	UG/L	UJ	mc



Sample ID	Matrix	Compound	Result	LOD	Units	Validation Qualifiers	Validation Reason
VPB152-GW-061614-700-702	WG	4-METHYL-2-PENTANONE		250	UG/L	UJ	c,mc
VPB152-GW-061614-700-702	WG	ACETONE		250	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	BENZENE		50	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	BROMODICHLOROMETHANE		50	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	BROMOFORM		50	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	BROMOMETHANE		100	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	CARBON DISULFIDE		50	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	CARBON TETRACHLORIDE		50	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	CHLOROBENZENE		50	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	CHLOROETHANE		100	UG/L	UJ	c,mc
VPB152-GW-061614-700-702	WG	CHLOROFORM		50	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	CHLOROMETHANE		100	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	CIS-1,2-DICHLOROETHENE		50	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	CIS-1,3-DICHLOROPROPENE		50	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	CYCLOHEXANE		50	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	DIBROMOCHLOROMETHANE		50	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	DICHLORODIFLUOROMETHANE		100	UG/L	UJ	c,mc
VPB152-GW-061614-700-702	WG	ETHYLBENZENE		50	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	ISOPROPYLBENZENE		50	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	M- AND P-XYLENE		100	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	METHYL ACETATE		75	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	METHYL CYCLOHEXANE		50	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	METHYL TERT-BUTYL ETHER		50	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	METHYLENE CHLORIDE		500*	UG/L	UJ	bl,mc
VPB152-GW-061614-700-702	WG	O-XYLENE		50	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	STYRENE		50	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	TETRACHLOROETHENE		50	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	TOLUENE		50	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	TRANS-1,2-DICHLOROETHENE		50	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	TRANS-1,3-DICHLOROPROPENE		50	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	TRICHLOROETHENE		50	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	TRICHLOROFLUOROMETHANE		100	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	VINYL CHLORIDE		100	UG/L	UJ	mc
VPB152-GW-061614-700-702	WG	XYLENES, TOTAL		150	UG/L	UJ	mc
VPB152-GWD-061314	WG	2-BUTANONE		2.5	UG/L	UJ	c
VPB152-GWD-061314	WG	4-METHYL-2-PENTANONE		2.5	UG/L	UJ	c
VPB152-GWD-061314	WG	ACETONE	9.5	2.5	UG/L	J	c
VPB152-GWD-061314	WG	CARBON DISULFIDE		1.0*	UG/L	U	bl
VPB152-GWD-061314	WG	CHLOROETHANE		1.0	UG/L	UJ	c
VPB152-GWD-061314	WG	DICHLORODIFLUOROMETHANE		1.0	UG/L	UJ	c
VPB152-TRIP BLANK-061614	WQ	2-BUTANONE		2.5	UG/L	UJ	c
VPB152-TRIP BLANK-061614	WQ	4-METHYL-2-PENTANONE		2.5	UG/L	UJ	c
VPB152-TRIP BLANK-061614	WQ	CHLOROETHANE		1.0	UG/L	UJ	c
VPB152-TRIP BLANK-061614	WQ	DICHLORODIFLUOROMETHANE		1.0	UG/L	UJ	c

\*LOQ

## Attachment A

### Nonconformance Summary Tables

**Table A-1a - Initial Calibration**

ICAL	Compound	% RSD	Limit
WG144357-7 GCMS-C	CHLOROETHANE	36.2	< 15%
Associated samples: All samples in the SDG			

**Table A-1b - Initial Calibration Verification**

ICV	Compound	% R	Limit
WG144357-7 GCMS-C	ACETONE	120.4	80-120%
Associated samples: All samples in the SDG			

**Table A-1c - Continuing Calibration Verification**

CCV ID	Compound	% D	Limit
WG144955 6/18/2014 GCMS-C	DICHLORODIFLUOROMETHANE	56.1	< 20%
	2-BUTANONE	20.7	< 20%
	4-METHYL-2-PENTANONE	37.9	< 20%
Associated samples: Samples VPB152-GW-061314-620-622, VPB152-GW-061614-700-702, VPB152-TRIP BLANK-061614			

**Table A-1d - Continuing Calibration Verification**

CCV ID	Compound	% D	Limit
WG145023 6/19/2014 GCMS-C	DICHLORODIFLUOROMETHANE	55.0	< 20%
	2-BUTANONE	29.0	< 20%
	4-METHYL-2-PENTANONE	37.4	< 20%
Associated samples: Samples VPB152-GW-061314-640-642, VPB152-GW-061614-660-662, VPB152-GW-061614-680-682, VPB152-GWD-061314			

**Table A-2 - Lab Blanks**

Blank ID	Compound	Result	LOD	Units	Associated Samples
WG144955-2	METHYLENE CHLORIDE	1.3	2.5	UG/L	VPB152-GW-061614-700-702
WG145023-2	CARBON DISULFIDE	0.32	0.50	UG/L	VPB152-GW-061314-640-642, VPB152-GW-061614-680-682, VPB152-GWD-061314

**Attachment B**  
**Qualifier Codes and Explanations**

<b>Qualifier</b>	<b>Explanation</b>
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.

### Attachment C

#### Reason Codes and Explanations

Reason Code	Explanation
be	Equipment blank contamination
bf	Field blank contamination
bl	Laboratory blank contamination
c	Calibration issue
co	Analyte carryover
d	Reporting limit raised due to chromatographic interference
fd	Field duplicate RPDs
h	Holding times
i	Internal standard areas
k	Estimated Maximum Possible Concentration (EMPC)
l	LCS or OPR recoveries
lc	Labeled compound recovery
ld	Laboratory duplicate RPDs
lp	Laboratory control sample/laboratory control sample duplicate RPDs
m	Matrix spike recovery
md	Matrix spike/matrix spike duplicate RPDs
nb	Negative laboratory blank contamination
p	Chemical preservation issue
r	Dual column RPD
q	Quantitation issue
s	Surrogate recovery
su	Ion suppression
t	Temperature preservation issue
x	Percent solids
y	Serial dilution results
z	ICS results

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4301-1  
**Client ID:** 152-061314-620-622  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4301  
**Lab File ID:** C7731.D

**Sample Date:** 13-JUN-14  
**Received Date:** 17-JUN-14  
**Extract Date:** 18-JUN-14  
**Extracted By:** REC  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144955

**Analysis Date:** 18-JUN-14  
**Analyst:** REC  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 20-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
<b>Carbon Disulfide</b>	U	0.40	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
<b>Acetone</b>	U	14	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
<b>Toluene</b>	U	3.7	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50



## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4301-1  
**Client ID:** 152-061314-620-622  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4301  
**Lab File ID:** C7731.D

**Sample Date:** 13-JUN-14  
**Received Date:** 17-JUN-14  
**Extract Date:** 18-JUN-14  
**Extracted By:** REC  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144955

**Analysis Date:** 18-JUN-14  
**Analyst:** REC  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 20-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U <b>JS</b>	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		93.6	%					
Toluene-d8		95.3	%					
1,2-Dichloroethane-d4		113.	%					
Dibromofluoromethane		101.	%					

*Handwritten signature and date: 6/20/14*

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4301-2RA  
**Client ID:** 152-061314-640-642  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4301  
**Lab File ID:** C7753.D

**Sample Date:** 13-JUN-14  
**Received Date:** 17-JUN-14  
**Extract Date:** 19-JUN-14  
**Extracted By:** REC  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG145023

**Analysis Date:** 19-JUN-14  
**Analyst:** REC  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 20-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U <del>UL</del> JS	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U <del>UL</del> JS	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
<b>Carbon Disulfide</b>	U <del>UL</del> JS	<del>0.35</del> 1.0	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
<b>Acetone</b>	U JS	7.0	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U <del>UL</del> JS	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	UL	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
<b>Toluene</b>	U JS	0.72	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	U <del>UL</del> JS	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50



## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4301-2RA  
**Client ID:** 152-061314-640-642  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4301  
**Lab File ID:** C7753.D

**Sample Date:** 13-JUN-14  
**Received Date:** 17-JUN-14  
**Extract Date:** 19-JUN-14  
**Extracted By:** REC  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG145023

**Analysis Date:** 19-JUN-14  
**Analyst:** REC  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 20-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	UL	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	UL	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	UL	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		96.4	%					
Toluene-d8		96.6	%					
1,2-Dichloroethane-d4		114.	%					
Dibromofluoromethane		102.	%					



### Report of Analytical Results

Client: ENSAFE  
 Lab ID: SH4301-3RA  
 Client ID: VPB152-GWD-061314  
 Project: Navy Clean WE15-03-06 NW  
 SDG: SH4301  
 Lab File ID: C7754.D

Sample Date: 13-JUN-14  
 Received Date: 17-JUN-14  
 Extract Date: 19-JUN-14  
 Extracted By: REC  
 Extraction Method: SW846 5030  
 Lab Prep Batch: WG145023

Analysis Date: 19-JUN-14  
 Analyst: REC  
 Analysis Method: SW846 8260B  
 Matrix: AQ  
 % Solids: NA  
 Report Date: 20-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	UL JS	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U JS	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	J U	0.38 1.0	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
Acetone	J	9.5	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U JS	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	UL	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	J	0.79	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	U JS	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4301-3RA  
**Client ID:** VPB152-GWD-061314  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4301  
**Lab File ID:** C7754.D

**Sample Date:** 13-JUN-14  
**Received Date:** 17-JUN-14  
**Extract Date:** 19-JUN-14  
**Extracted By:** REC  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG145023

**Analysis Date:** 19-JUN-14  
**Analyst:** REC  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 20-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	UL	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	UL	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	UL	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		96.3	%					
Toluene-d8		98.2	%					
1,2-Dichloroethane-d4		114.	%					
Dibromofluoromethane		102.	%					

### Report of Analytical Results

Client: ENSAFE  
 Lab ID: SH4301-4RA  
 Client ID: 152-061614-660-662  
 Project: Navy Clean WE15-03-06 NW  
 SDG: SH4301  
 Lab File ID: C7755.D

Sample Date: 16-JUN-14  
 Received Date: 17-JUN-14  
 Extract Date: 19-JUN-14  
 Extracted By: REC  
 Extraction Method: SW846 5030  
 Lab Prep Batch: WG145023

Analysis Date: 19-JUN-14  
 Analyst: REC  
 Analysis Method: SW846 8260B  
 Matrix: AQ  
 % Solids: NA  
 Report Date: 24-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	<del>UL</del> <b>UJ</b>	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	<del>U</del> <b>UJ</b>	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
<b>Acetone</b>	<b>J</b>	8.4	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	<del>U</del> <b>UJ</b>	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	ULM	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
<b>Toluene</b>	<b>J</b>	0.52	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	<del>U</del> <b>UJ</b>	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50



## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4301-4RA  
**Client ID:** 152-061614-660-662  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4301  
**Lab File ID:** C7755.D

**Sample Date:** 16-JUN-14  
**Received Date:** 17-JUN-14  
**Extract Date:** 19-JUN-14  
**Extracted By:** REC  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG145023

**Analysis Date:** 19-JUN-14  
**Analyst:** REC  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 24-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	ULM	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	UL	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	UL	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		93.0	%					
Toluene-d8		94.6	%					
1,2-Dichloroethane-d4		112.	%					
Dibromofluoromethane		97.4	%					

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4301-5RA  
**Client ID:** 152-061614-680-682  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4301  
**Lab File ID:** C7756.D

**Sample Date:** 16-JUN-14  
**Received Date:** 17-JUN-14  
**Extract Date:** 19-JUN-14  
**Extracted By:** REC  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG145023

**Analysis Date:** 19-JUN-14  
**Analyst:** REC  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 20-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U UJ	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U UJ	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
<b>Carbon Disulfide</b>	U J UJ	<del>0.47</del> 1.0	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
<b>Acetone</b>	J	22	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
<b>2-Butanone</b>	J J	4.8	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	UL	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
<b>Toluene</b>	J	0.69	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50

*Handwritten signature/initials*

### Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4301-5RA  
**Client ID:** 152-061614-680-682  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4301  
**Lab File ID:** C7756.D

**Sample Date:** 16-JUN-14  
**Received Date:** 17-JUN-14  
**Extract Date:** 19-JUN-14  
**Extracted By:** REC  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG145023

**Analysis Date:** 19-JUN-14  
**Analyst:** REC  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 20-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	UL	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	UL	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	UL	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		101.	%					
Toluene-d8		101.	%					
1,2-Dichloroethane-d4		119.	%					
Dibromofluoromethane		102.	%					

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4301-7DL  
**Client ID:** 152-061614-700-702  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4301  
**Lab File ID:** C7728.D

**Sample Date:** 16-JUN-14  
**Received Date:** 17-JUN-14  
**Extract Date:** 18-JUN-14  
**Extracted By:** REC  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144955

**Analysis Date:** 18-JUN-14  
**Analyst:** REC  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 20-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	100	ug/L	100	2	200	24.	100
Chloromethane	U	100	ug/L	100	2	200	36.	100
Vinyl Chloride	U	100	ug/L	100	2	200	25.	100
Bromomethane	U	100	ug/L	100	2	200	49.	100
Chloroethane	U	100	ug/L	100	2	200	55.	100
Trichlorofluoromethane	U	100	ug/L	100	2	200	24.	100
1,1-Dichloroethene	U	50	ug/L	100	1	100	35.	50.
Carbon Disulfide	U	50	ug/L	100	1	100	25.	50.
Freon-113	U	50	ug/L	100	1	100	31.	50.
<b>Methylene Chloride</b>	U	<del>180</del> 500	ug/L	100	5	500	110	250
Acetone	U	250	ug/L	100	5	500	220	250
trans-1,2-Dichloroethene	U	50	ug/L	100	1	100	25.	50.
Methyl tert-butyl Ether	U	50	ug/L	100	1	100	36.	50.
1,1-Dichloroethane	U	50	ug/L	100	1	100	21.	50.
cis-1,2-Dichloroethene	U	50	ug/L	100	1	100	21.	50.
Chloroform	U	50	ug/L	100	1	100	32.	50.
1,1,1-Trichloroethane	U	50	ug/L	100	1	100	20.	50.
2-Butanone	U	250	ug/L	100	5	500	130	250
Cyclohexane	U	50	ug/L	100	1	100	31.	50.
Carbon Tetrachloride	U	50	ug/L	100	1	100	22.	50.
Benzene	U	50	ug/L	100	1	100	26.	50.
1,2-Dichloroethane	U	50	ug/L	100	1	100	20.	50.
Trichloroethene	U	50	ug/L	100	1	100	28.	50.
1,2-Dichloropropane	U	50	ug/L	100	1	100	25.	50.
Bromodichloromethane	U	50	ug/L	100	1	100	33.	50.
cis-1,3-Dichloropropene	U	50	ug/L	100	1	100	19.	50.
Toluene	U	50	ug/L	100	1	100	27.	50.
4-Methyl-2-Pentanone	U	250	ug/L	100	5	500	130	250
trans-1,3-Dichloropropene	U	50	ug/L	100	1	100	20.	50.
1,1,2-Trichloroethane	U	50	ug/L	100	1	100	33.	50.
Tetrachloroethene	U	50	ug/L	100	1	100	40.	50.
Dibromochloromethane	U	50	ug/L	100	1	100	30.	50.
2-Hexanone	U	250	ug/L	100	5	500	170	250
Chlorobenzene	U	50	ug/L	100	1	100	22.	50.
Ethylbenzene	U	50	ug/L	100	1	100	21.	50.



## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4301-7DL  
**Client ID:** 152-061614-700-702  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4301  
**Lab File ID:** C7728.D

**Sample Date:** 16-JUN-14  
**Received Date:** 17-JUN-14  
**Extract Date:** 18-JUN-14  
**Extracted By:** REC  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144955

**Analysis Date:** 18-JUN-14  
**Analyst:** REC  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 20-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	150	ug/L	100	3	300	25.	150
Styrene	U	50	ug/L	100	1	100	23.	50.
Bromoform	U	50	ug/L	100	1	100	23.	50.
Isopropylbenzene	U	50	ug/L	100	1	100	23.	50.
1,1,2,2-Tetrachloroethane	U	50	ug/L	100	1	100	38.	50.
1,3-Dichlorobenzene	U	50	ug/L	100	1	100	26.	50.
1,4-Dichlorobenzene	U	50	ug/L	100	1	100	24.	50.
1,2-Dichlorobenzene	U	50	ug/L	100	1	100	15.	50.
1,2,4-Trichlorobenzene	U	50	ug/L	100	1	100	37.	50.
Methyl Acetate	U	75	ug/L	100	1	100	53.	75.
Methylcyclohexane	U	50	ug/L	100	1	100	30.	50.
o-Xylene	U	50	ug/L	100	1	100	25.	50.
M+P-Xylenes	U	100	ug/L	100	2	200	59.	100
1,2-Dichloroethylene (Total)	U	100	ug/L	100	2	200	21.	100
1,2-Dibromoethane	U	50	ug/L	100	1	100	22.	50.
1,2-Dibromo-3-Chloropropane	U	75	ug/L	100	1	100	50.	75.
P-Bromofluorobenzene		90.8	%					
Toluene-d8		95.3	%					
1,2-Dichloroethane-d4		113.	%					
Dibromofluoromethane		101.	%					

*Handwritten signature: REC 6/20/14*



## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4301-6  
**Client ID:** VPB152-TB-061614  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4301  
**Lab File ID:** C7725.D

**Sample Date:** 16-JUN-14  
**Received Date:** 17-JUN-14  
**Extract Date:** 18-JUN-14  
**Extracted By:** REC  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144955

**Analysis Date:** 18-JUN-14  
**Analyst:** REC  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 20-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
Acetone	U	2.5	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4301-6  
**Client ID:** VPB152-TB-061614  
**Project:** Navy Clean WE15-03-06\NW  
**SDG:** SH4301  
**Lab File ID:** C7725.D

**Sample Date:** 16-JUN-14  
**Received Date:** 17-JUN-14  
**Extract Date:** 18-JUN-14  
**Extracted By:** REC  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144955

**Analysis Date:** 18-JUN-14  
**Analyst:** REC  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 20-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		92.4	%					
Toluene-d8		95.4	%					
1,2-Dichloroethane-d4		113.	%					
Dibromofluoromethane		98.3	%					



## Data Validation Report

Project: Regional Groundwater Investigation - NWIRP Bethpage  
Laboratory: Katahdin Analytical Services, Scarborough, Maine  
Service Request: SH4222  
Analyses/Method: EPA SW-846 Method 8260B for VOCs (GC/MS) and Standard Method 5310B for Total Organic Carbon by High-Temperature Combustion  
Validation Level: Limited  
RESCON Project 60266526.SA.DV  
Number:  
Prepared by: Sheena Blair/RESCON Completed on: 9/15/2014  
Reviewed by: Lori Herberich/RESCON File Name: SH4222\_5310B and 8260B

### SUMMARY

The samples listed below were collected by Resolution Consultants (RESCON) from the Regional Groundwater Investigation - NWIRP Bethpage site on June 10,11, and 12, 2014.

Sample ID	Matrix/Sample Type
VPB152-EB-061114	Equipment blank
VPB152-GW-061014-440-442	Groundwater
VPB152-GW-061014-460-462	Groundwater
VPB152-GW-061014-480-482	Groundwater
VPB152-GW-061014-500-502	Groundwater
VPB152-GW-061114-520-522	Groundwater
VPB152-GW-061114-540-542	Groundwater
VPB152-GW-061114-560-562	Groundwater
VPB152-GW-061214-580-582	Groundwater
VPB152-GW-061214-605-607	Groundwater
VPB152-TRIP BLANK-061214	Trip Blank

The samples were analyzed in accordance with:

- *Test Methods for Evaluating Solid Waste, Physical/Chemical Methods SW-846, Method 8260B, Volatile Organic Compounds by Gas Chromatography/Mass Spectrometry (USEPA, 1996).*
- *Standard Methods for the Examination of Water and Wastewater, Method SM310B, Total Organic Carbon by High-Temperature Combustion*

Data validation activities were conducted with reference to these methods, *USEPA Contract Laboratory Program National Functional Guidelines for Superfund Organic Methods Data Review (June 2008)*, *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
- ✓ GC/MS performance checks
- X Initial calibration/continuing calibration verification
- X Laboratory blanks/equipment blanks/trip blanks
- ✓ Surrogate spike recoveries
- NA Matrix spike (MS) results
- ✓ Laboratory control sample (LCS) results
- ✓ Field duplicate results
- ✓ Internal standard results
- ✓ 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 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. Selected data points were estimated and/or negated due to nonconformances of certain QC criteria (see discussion below). Qualified sample results are presented in Table 1.

## 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 "VPB" prefix from the sample ID, and truncated IDs for GW and Trip Blank 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. All QC acceptance criteria were met.

### **GC/MS Performance Checks**

The data were reviewed to ensure that the 4-bromofluorobenzene (BFB) tuning was performed at the correct frequency and that the method acceptance criteria were met. All QC acceptance criteria were met.

### **Initial Calibration/Continuing Calibration Verification**

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

- the initial calibration (ICAL) percent relative standard deviation (%RSD), correlation coefficient (r)/coefficient of determination ( $r^2$ ), and/or response factor method acceptance criteria were met;
- the initial calibration verification (ICV) percent recovery (%R) criteria were met; and
- the continuing calibration verification standard (CCV) method percent difference or percent drift (%Ds), %Rs, and/or RF acceptance criteria were met; and/or
- the retention time method acceptance criteria were met.

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

#### **ICV Recovery Nonconformances:**

Nonconformance	Actions	
	Detected Compounds	Nondetected Compounds
%R > 120%	J	No qualification
20% < %R < 80%	J	UJ
%R < 20% (see note)	J	R*

Notes: Based on NFG 2008 VOC guidance, professional judgment is used to reject (R) nondetects in all associated samples for any analyte with < 20% recovery. Also, professional judgment is used to estimate (UJ) rather the reject (R) sample results previously negated (U) on the basis of blank contamination.

#### **CCV Linearity Nonconformances:**

Nonconformance	Actions	
	Detected Results	Nondetected Results
%D > 20%	J	UJ
%Drift > 20%	J*	UJ*
* No guidance in NFG, thus professional judgment was used		

Qualified sample results are shown in Table 1. Nonconformances are summarized in Attachment A in Table A-1a and A-1b.

### **Laboratory Blanks/Equipment Blanks/Trip Blanks**

Laboratory method blanks, equipment rinsate and trip 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, equipment rinsate and trip 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.

Sample results were qualified as follows:

For common lab contaminants (methylene chloride, acetone, 2-butanone):

Blank type	Blank result	Sample result	Action for samples	
Method, Storage, Field, Trip, or Instrument*	Detects	Not detected	No qualification	
		< 2x LOQ	Report sample LOQ value with a U	
		$\geq 2x$ LOQ and $\leq 4x$ LOQ	Report the sample result with a U**	
	$\leq 2x$ LOQ	$\geq 4x$ LOQ	$\geq 4x$ LOQ	No qualifications
			< 2x LOQ	Report sample LOQ value with a U
			$\geq 2x$ LOQ and < blank contamination	Report the sample result with a U
		> 2x LOQ	$\geq 2x$ LOQ and $\geq$ blank contamination	If the result is $\leq 2x$ blank result, report the sample result U.** If the result is > 2x blank result, no qualification is required.**
			< 2x LOQ	Report sample LOQ value with a U
			$\geq 2x$ LOQ and < blank contamination	Report the sample result with a U
* Qualifications based on instrument blank results affect only the sample analyzed immediately after the sample that has target compounds that exceed the calibration range or non-target compounds that exceed 100 g/L.				
**Based on professional judgment				

For TOC:

Blank Type	Blank Result	Sample Result	Action for Samples
PB / EB/ FB (Positive)	> LOQ	$\geq$ DL but $\leq$ LOQ	Qualify as nondetect (U) at the LOQ
		>LOQ but < 10x Blank Result	Qualify results as unusable
		$\geq 10x$ Blank Result	No action
	$\geq$ DL but $\leq$ LOQ	Nondetect	No action
		$\geq$ DL but $\leq$ LOQ	Qualify as nondetect (U) at the LOQ
		> LOQ	Use professional judgment (see below [1])

LOQ - Limit of Quantitation.

Nonconformances are summarized in Attachment A in Tables A-2 and A-3. Qualified sample results are shown in Table 1.

**Surrogate Spike Recoveries**

The surrogate recoveries (%Rs) were reviewed for conformance with the QC acceptance criteria.

Although the %Rs were high for surrogates in two samples, qualification of the data was not required because nondetects were reported for the samples.

**MS Results**

MS/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**

Field duplicate RPDs were reviewed for conformance with the QC criterion of  $\leq 30\%$  for aqueous matrices. This criterion applies if both results were  $>5x$  the LOQ. The criterion is doubled if the results were  $<5x$  the LOQ. All QC acceptance criteria were met.

**Internal Standard Results**

The internal standard (IS) recoveries were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

**Sample Results/Reporting Issues**

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

Compounds detected at concentrations less than the LOQ but greater than the 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**

Sample results qualified as a result of validation actions are summarized in Table 1. All actions are described above.

**ATTACHMENTS**

Attachment A: Nonconformance Summary Tables

Attachment B: Qualifier Codes and Explanations

Attachment C: Reason Codes and Explanations



**Table 1 - Data Validation Summary of Qualified Data**

Sample ID	Matrix	Compound	Result	LOD	Units	Validation Qualifiers	Validation Reason
VPB152-EB-061114	WQ	TOTAL ORGANIC CARBON		1.0*	MG/L	U	bl
VPB152-EB-061114	WQ	ACETONE	4.1	2.5	UG/L	J	c
VPB152-GW-061014-440-442	WG	ACETONE		5.0*	UG/L	U	be
VPB152-GW-061014-480-482	WG	ACETONE		5.0*	UG/L	U	be
VPB152-GW-061014-500-502	WG	1,2-DIBROMO-3-CHLOROPROPANE		0.75	UG/L	UJ	c
VPB152-GW-061014-500-502	WG	2-HEXANONE		2.5	UG/L	UJ	c
VPB152-GW-061014-500-502	WG	4-METHYL-2-PENTANONE		2.5	UG/L	UJ	c
VPB152-GW-061014-500-502	WG	ACETONE		5.0*	UG/L	U	be
VPB152-GW-061114-520-522	WG	ACETONE		5.0*	UG/L	U	be
VPB152-GW-061114-560-562	WG	ACETONE		5.0*	UG/L	U	be
VPB152-GW-061214-580-582	WG	ACETONE		5.0*	UG/L	U	be
*LOQ							

## Attachment A

### Non Conformance Summary Tables

**Table A-1 - Initial Calibration Verification**

ICV	Compound	% R	Limit
WG144515 GC-MS-D	ACETONE	125.4	80-120%
Associated samples: All samples in the SDG			

**Table A-1b - Continuing Calibration Verification**

CCV ID	Compound	% D	Limit
WG144718-4 GCMS-D 6/14/2014	4-METHYL-2-PENTANONE	21.2	< 20%
	2-HEXANONE	27.7	< 20%
	1,2-DIBROMO-3-CHLOROPROPANE	22.5	< 20%
Associated sample: VPB152-GW-061014-500-502			

**Table A-2 - Lab Blanks**

Blank ID	Compound	Result	LOQ	Units	Associated Samples
WG145514-1	TOTAL ORGANIC CARBON	0.21	1.0	MG/L	VPB152-EB-061114

**Table A-3- Field Blanks**

Blank ID	Compound	Result	LOQ	Units	Associated Samples
VPB152-EB-061114	ACETONE	4.1	5.0	UG/L	All field samples in this SDG

**Attachment B**  
**Qualifier Codes and Explanations**

<b>Qualifier</b>	<b>Explanation</b>
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.

### Attachment C

#### Reason Codes and Explanations

Reason Code	Explanation
be	Equipment blank contamination
bf	Field blank contamination
bl	Laboratory blank contamination
c	Calibration issue
co	Analyte carryover
d	Reporting limit raised due to chromatographic interference
fd	Field duplicate RPDs
h	Holding times
i	Internal standard areas
k	Estimated Maximum Possible Concentration (EMPC)
l	LCS or OPR recoveries
lc	Labeled compound recovery
ld	Laboratory duplicate RPDs
lp	Laboratory control sample/laboratory control sample duplicate RPDs
m	Matrix spike recovery
md	Matrix spike/matrix spike duplicate RPDs
nb	Negative laboratory blank contamination
p	Chemical preservation issue
r	Dual column RPD
q	Quantitation issue
s	Surrogate recovery
su	Ion suppression
t	Temperature preservation issue
x	Percent solids
y	Serial dilution results
z	ICS results
mc	Method compliance nonconformance

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4222-1  
**Client ID:** 152-061014-440-442  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4222  
**Lab File ID:** D8915.D

**Sample Date:** 10-JUN-14  
**Received Date:** 13-JUN-14  
**Extract Date:** 13-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144658

**Analysis Date:** 13-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 16-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
Acetone	U	<del>4.8</del> 5.0	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50



## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4222-1  
**Client ID:** 152-061014-440-442  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4222  
**Lab File ID:** D8915.D

**Sample Date:** 10-JUN-14  
**Received Date:** 13-JUN-14  
**Extract Date:** 13-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144658

**Analysis Date:** 13-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 16-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		104.	%					
Toluene-d8		111.	%					
1,2-Dichloroethane-d4		112.	%					
Dibromofluoromethane		106.	%					

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4222-2  
**Client ID:** 152-061014-460-462  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4222  
**Lab File ID:** D8916.D

**Sample Date:** 10-JUN-14  
**Received Date:** 13-JUN-14  
**Extract Date:** 13-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144658

**Analysis Date:** 13-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 16-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
Acetone	U	2.5	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4222-2  
**Client ID:** 152-061014-460-462  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SII4222  
**Lab File ID:** D8916.D

**Sample Date:** 10-JUN-14  
**Received Date:** 13-JUN-14  
**Extract Date:** 13-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144658

**Analysis Date:** 13-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 16-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		105.	%					
Toluene-d8		117.	%					
1,2-Dichloroethane-d4		118.	%					
Dibromofluoromethane		110.	%					



## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4222-3  
**Client ID:** 152-061014-480-482  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4222  
**Lab File ID:** D8917.D

**Sample Date:** 10-JUN-14  
**Received Date:** 13-JUN-14  
**Extract Date:** 13-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144658

**Analysis Date:** 13-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 16-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
Acetone	U	<del>3.2</del> 5.0	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50



## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4222-3  
**Client ID:** 152-061014-480-482  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4222  
**Lab File ID:** D8917.D

**Sample Date:** 10-JUN-14  
**Received Date:** 13-JUN-14  
**Extract Date:** 13-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144658

**Analysis Date:** 13-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 16-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		99.9	%					
Toluene-d8		112.	%					
1,2-Dichloroethane-d4		114.	%					
Dibromofluoromethane		105.	%					

### Report of Analytical Results

Client: ENSAFE  
 Lab ID: SH4222-4RA  
 Client ID: 152-061014-500-502  
 Project: Navy Clean WE15-03-06 NW  
 SDG: SH4222  
 Lab File ID: D8937.D

Sample Date: 10-JUN-14  
 Received Date: 13-JUN-14  
 Extract Date: 14-JUN-14  
 Extracted By: DJP  
 Extraction Method: SW846 5030  
 Lab Prep Batch: WG144718

Analysis Date: 14-JUN-14  
 Analyst: DJP  
 Analysis Method: SW846 8260B  
 Matrix: AQ  
 % Solids: NA  
 Report Date: 16-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
Acetone	<del>U</del> U	<del>3.8</del> 5.0	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	<del>U</del> U	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	<del>U</del> U	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50



## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4222-4RA  
**Client ID:** 152-061014-500-502  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4222  
**Lab File ID:** D8937.D

**Sample Date:** 10-JUN-14  
**Received Date:** 13-JUN-14  
**Extract Date:** 14-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144718

**Analysis Date:** 14-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 16-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	<del>U</del> UJ	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		103.	%					
Toluene-d8		113.	%					
1,2-Dichloroethane-d4		112.	%					
Dibromofluoromethane		107.	%					

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### Report of Analytical Results

Client: ENSAFE  
 Lab ID: SH4222-5  
 Client ID: 152-061114-520-522  
 Project: Navy Clean WE15-03-06 NW  
 SDG: SH4222  
 Lab File ID: D8914.D

Sample Date: 11-JUN-14  
 Received Date: 13-JUN-14  
 Extract Date: 13-JUN-14  
 Extracted By: DJP  
 Extraction Method: SW846 5030  
 Lab Prep Batch: WG144658

Analysis Date: 13-JUN-14  
 Analyst: DJP  
 Analysis Method: SW846 8260B  
 Matrix: AQ  
 % Solids: NA  
 Report Date: 16-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
Acetone	U	<del>9.0</del> 5.0	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50



## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4222-5  
**Client ID:** 152-061114-520-522  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4222  
**Lab File ID:** D8914.D

**Sample Date:** 11-JUN-14  
**Received Date:** 13-JUN-14  
**Extract Date:** 13-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144658

**Analysis Date:** 13-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 16-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		102.	%					
Toluene-d8		112.	%					
1,2-Dichloroethane-d4		112.	%					
Dibromofluoromethane		103.	%					

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4222-6  
**Client ID:** 152-061114-540-542  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4222  
**Lab File ID:** D8919.D

**Sample Date:** 11-JUN-14  
**Received Date:** 13-JUN-14  
**Extract Date:** 13-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144658

**Analysis Date:** 13-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 16-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
Acetone	U	2.5	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4222-6  
**Client ID:** 152-061114-540-542  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4222  
**Lab File ID:** D8919.D

**Sample Date:** 11-JUN-14  
**Received Date:** 13-JUN-14  
**Extract Date:** 13-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144658

**Analysis Date:** 13-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 16-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		101.	%					
Toluene-d8		113.	%					
1,2-Dichloroethane-d4		116.	%					
Dibromofluoromethane		109.	%					



### Report of Analytical Results

Client: ENSAFE  
 Lab ID: SH4222-7  
 Client ID: VPB152-EB-061114  
 Project: Navy Clean WE15-03-06 NW  
 SDG: SH4222  
 Lab File ID: D8920.D

Sample Date: 11-JUN-14  
 Received Date: 13-JUN-14  
 Extract Date: 13-JUN-14  
 Extracted By: DJP  
 Extraction Method: SW846 5030  
 Lab Prep Batch: WG144658

Analysis Date: 13-JUN-14  
 Analyst: DJP  
 Analysis Method: SW846 8260B  
 Matrix: AQ  
 % Solids: NA  
 Report Date: 16-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
Acetone	<del>U</del> U	<del>4.1</del> 5.0	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50



## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4222-7  
**Client ID:** VPB152-EB-061114  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4222  
**Lab File ID:** D8920.D

**Sample Date:** 11-JUN-14  
**Received Date:** 13-JUN-14  
**Extract Date:** 13-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144658

**Analysis Date:** 13-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 16-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		97.7	%					
Toluene-d8		112.	%					
1,2-Dichloroethane-d4		118.	%					
Dibromofluoromethane		110.	%					

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4222-8  
**Client ID:** 152-061114-560-562  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4222  
**Lab File ID:** D8921.D

**Sample Date:** 11-JUN-14  
**Received Date:** 13-JUN-14  
**Extract Date:** 13-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144658

**Analysis Date:** 13-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 16-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
<b>Acetone</b>	<b>U</b>	<del>1.0</del> <b>5.0</b>	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50



### Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SHI4222-8  
**Client ID:** 152-061114-560-562  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4222  
**Lab File ID:** D8921.D

**Sample Date:** 11-JUN-14  
**Received Date:** 13-JUN-14  
**Extract Date:** 13-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144658

**Analysis Date:** 13-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 16-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		99.0	%					
Toluene-d8		112.	%					
1,2-Dichloroethane-d4		117.	%					
Dibromofluoromethane		108.	%					

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4222-9  
**Client ID:** VPB152-TB-061214  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4222  
**Lab File ID:** D8911.D

**Sample Date:** 12-JUN-14  
**Received Date:** 13-JUN-14  
**Extract Date:** 13-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144658

**Analysis Date:** 13-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 16-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
Acetone	U	2.5	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4222-9  
**Client ID:** VPB152-TB-061214  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4222  
**Lab File ID:** D8911.D

**Sample Date:** 12-JUN-14  
**Received Date:** 13-JUN-14  
**Extract Date:** 13-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144658

**Analysis Date:** 13-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 16-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		99.7	%					
Toluene-d8		110.	%					
1,2-Dichloroethane-d4		109.	%					
Dibromofluoromethane		106.	%					

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4222-10  
**Client ID:** 152-061214-580-582  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4222  
**Lab File ID:** D8922.D

**Sample Date:** 12-JUN-14  
**Received Date:** 13-JUN-14  
**Extract Date:** 13-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144658

**Analysis Date:** 13-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 16-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
Acetone	U	6.45.0	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50



## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4222-10  
**Client ID:** 152-061214-580-582  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4222  
**Lab File ID:** D8922.D

**Sample Date:** 12-JUN-14  
**Received Date:** 13-JUN-14  
**Extract Date:** 13-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144658

**Analysis Date:** 13-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 16-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		103.	%					
Toluene-d8		116.	%					
1,2-Dichloroethane-d4	*	120.	%					
Dibromofluoromethane		113.	%					



## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4222-11  
**Client ID:** 152-061214-605-607  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4222  
**Lab File ID:** D8923.D

**Sample Date:** 12-JUN-14  
**Received Date:** 13-JUN-14  
**Extract Date:** 13-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144658

**Analysis Date:** 13-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 16-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
Acetone	U	2.5	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4222-11  
**Client ID:** 152-061214-605-607  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4222  
**Lab File ID:** D8923.D

**Sample Date:** 12-JUN-14  
**Received Date:** 13-JUN-14  
**Extract Date:** 13-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144658

**Analysis Date:** 13-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 16-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropanoic acid	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		108.	%					
Toluene-d8	*	120.	%					
1,2-Dichloroethane-d4	*	125.	%					
Dibromofluoromethane	*	117.	%					

## Report of Analytical Results

**Client:** Rick Purdy  
 AECOM  
 701 Edgewater Drive  
 Wakefield, MA 01880

**Lab Sample ID:** SH4222-7  
**Report Date:** 01-JUL-14  
**Client PO:** 16518  
**Project:** Navy Clean WE15-03-0  
**SDG:** SH4222

**Sample Description**  
 VPB152-EB-061114

**Matrix** AQ  
**Date Sampled** 11-JUN-14  
**Date Received** 13-JUN-14

Parameter	Result	Adj LOQ	Adj MDL	Adj LOD	Anal. Method	QC Batch	Anal. Date	Prep. Method	Prep. Date	Footnotes
Total Organic Carbon	10.20 mg/L	1.0	0.10	.5	SM5310B	WG145514	26-JUN-14 19:24:35	N/A	N/A	N/A

1.0 U

*Result*





## Data Validation Report

Project:	Regional Groundwater Investigation - NWIRP Bethpage	
Laboratory:	Katahdin Analytical Services, Scarborough, Maine	
Service Request:	SH4056	
Analyses/Method:	EPA SW-846, Method 8260B for VOCs (GC/MS), Total Organic Carbon by EPA SW-846 Method 9060 for TOC (Carbonaceous analyzer, IR or FID), and Standard Methods 5310 for Total Organic Carbon by High-Temperature Combustion	
Validation Level:	Limited	
RESCON Project Number:	60266526.SA.DV	
Prepared by:	Sheena Blair/RESCON	Completed on: 9/14/2014
Reviewed by:	Lori Herberich/RESCON	File Name: SH4056__5310B, 8260B and 9060

### SUMMARY

The samples listed below were collected by Resolution Consultants (RESCON) from the Regional Groundwater Investigation - NWIRP Bethpage site on June 6 and 9, 2014.

Sample ID	Matrix/Sample Type
VPB152-FB-060614	Field blank
VPB152-GW-060614-320-322	Groundwater
VPB152-GW-060614-345-347	Groundwater
VPB152-GW-060914-360-362	Groundwater
VPB152-GW-060914-380-382	Groundwater
VPB152-GW-060914-400-402	Groundwater
VPB152-GW-060914-420-422	Groundwater
VPB152-SOIL-060614-325-327	Soil
VPB152-SOIL-D-060614	Field Duplicate of VPB152-SOIL-060614-325-327
VPB152-TRIP BLANK-060914	Trip Blank

The samples were analyzed in accordance with:

- *Test Methods for Evaluating Solid Waste, Physical/Chemical Methods SW-846, Method 8260B, Volatile Organic Compounds by Gas Chromatography/Mass Spectrometry (USEPA, 1996).*
- *Test Methods for Evaluating Solid Waste, Physical/Chemical Methods SW-846, Method 9060A, Total Organic Carbon (USEPA 2004);*
- *Standard Methods for the Examination of Water and Wastewater, Method SM5310B, Total Organic Carbon by High-Temperature Combustion*

Data validation activities were conducted with reference to these methods, *USEPA Contract Laboratory Program National Functional Guidelines for Superfund Organic Methods Data Review* (June 2008), *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
- ✓ GC/MS performance checks
- ✗ Initial calibration/continuing calibration verification
- ✗ Laboratory blanks/equipment blanks/trip blanks
- ✗ Surrogate spike recoveries
- ✓ Matrix spike (MS)/matrix spike duplicate (MSD) results
- ✓ Laboratory control sample (LCS) results
- ✓ Field duplicate results
- ✓ Internal standard results
- ✓ 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. Selected data points were estimated or negated due to nonconformances of certain QC criteria (see discussion below). Qualified sample results are presented in Table 1.

## 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 "VPB-" prefix and the "GW" from the sample ID, and truncated the ID for the Trip Blank 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. All QC acceptance criteria were met.

#### **GC/MS Performance Checks**

The data were reviewed to ensure that the 4-bromofluorobenzene (BFB) tuning was performed at the correct frequency and that the method acceptance criteria were met. All QC acceptance criteria were met.

#### **Initial Calibration/Continuing Calibration Verification**

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

- the initial calibration (ICAL) percent relative standard deviation (%RSD), correlation coefficient (r)/coefficient of determination ( $r^2$ ), and/or response factor method acceptance criteria were met;
- the initial calibration verification (ICV) percent recovery (%R) criteria were met; and
- the continuing calibration verification standard (CCV) method percent difference or percent drift (%Ds), %Rs, and/or RF acceptance criteria were met; and/or
- the retention time method acceptance criteria were met.

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

#### **ICV Recovery Nonconformances:**

Nonconformance	Actions	
	Detected Compounds	Nondetected Compounds
%R > 120%	J	No qualification
20% < %R < 80%	J	UJ
%R < 20% (see note)	J	R*

Notes: Based on NFG 2008 VOC guidance, professional judgment is used to reject (R) nondetects in all associated samples for any analyte with < 20% recovery. Also, professional judgment is used to estimate (UJ) rather the reject (R) sample results previously negated (U) on the basis of blank contamination.

Qualified sample results are shown in Table 1. Nonconformances are summarized in Attachment A in Table A-1.

#### **Laboratory Blanks/Equipment Blanks/Trip Blanks**

Laboratory method blanks, equipment rinsate and trip 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, equipment rinsate and trip 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.

TOC sample results were qualified as follows:

Blank Type	Blank Result	Sample Result	Action for Samples
PB / EB/ FB (Positive)	> LOQ	$\geq$ DL but $\leq$ LOQ	Qualify as nondetect (U) at the LOQ
		>LOQ but < 10x Blank Result	Qualify results as unusable
		$\geq$ 10x Blank Result	No action
	$\geq$ DL but $\leq$ LOQ	Nondetect	No action
		$\geq$ DL but $\leq$ LOQ	Qualify as nondetect (U) at the LOQ
		> LOQ	professional judgment (see below [1])

LOQ - Limit of Quantitation.

Nonconformances are summarized in Attachment A in Table A-2. Qualified sample results are shown in Table 1.

### Surrogate Spike Recoveries

The surrogate recoveries (%Rs) were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

### MS/MSD Results

MS/MSD analyses for VOCs were not performed on a sample reported in this SDG. There were no validation actions taken on this basis.

The MS/MSD %Rs and relative percent differences (RPDs) for TOC were reviewed for conformance with the QC acceptance criteria.

Data qualification to the analytes associated with the specific MS/MSD nonconformances was as follows:

Criteria	Action	
	Detected Compounds	Nondetected Compounds
%R > UL	J	No qualification
$30\% \leq \%R < LL$	J	UJ
%R < 30% (see note 1)	J	R*
%RPD > UL (see note 2)	J	No qualification

Note: Actions are applied to all samples with the same matrix. \*When the native sample concentration is >4X the concentration of the spike added (based on Region I criteria), evaluate the MS, MSD, and native sample with regards to %RSD rather than %R (RESCON professional judgment)

Notes:

- Based on NFG 2008 VOC guidance, professional judgment is used to reject (R) non-detects in all associated samples for any analyte with < 30% recovery. Also, professional judgment is used to estimate (UJ) rather the reject (R) sample results previously negated (U) on the basis of blank contamination.
- In the absence of Region 2 guidance, RPD actions are based on professional judgment.
- If a field duplicate sample was also collected for the native sample chosen for MS/MSD analysis, professional judgment is used to apply MS/MSD actions to the corresponding field duplicate sample as well as the native sample.



Nonconformances are summarized in Attachment A in Table A-3. Qualified sample results are shown in Table 1.

### **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  $\leq 60\%$  for solid matrices. This criterion applies if both results were greater than five times the Limit of Quantitation (LOQ). All QC acceptance criteria were met.

### **Internal Standard Results**

The internal standard (IS) recoveries were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

### **Sample Results/Reporting Issues**

Compounds that were not detected in the sample are reported as undetected (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**

Sample results qualified as a result of validation actions are summarized in Table 1. All actions are described above.

## **ATTACHMENTS**

Attachment A: Nonconformance Summary Tables

Attachment B: Qualifier Codes and Explanations

Attachment C: Reason Codes and Explanations

**Table 1 - Data Validation Summary of Qualified Data**

Sample ID	Matrix	Compound	Result	LOD	Units	Validation Qualifiers	Validation Reason
VPB152-GW-060614-320-322	WG	ACETONE	5.3	2.5	UG/L	J	c
VPB152-GW-060614-345-347	WG	ACETONE	4.4	2.5	UG/L	J	c
VPB152-GW-060914-360-362	WG	ACETONE	9.9	2.5	UG/L	J	c
VPB152-GW-060914-380-382	WG	ACETONE	6.9	2.5	UG/L	J	c
VPB152-GW-060914-420-422	WG	ACETONE	3.1	2.5	UG/L	J	c
VPB152-SOIL-060614-325-327	SO	TOTAL ORGANIC CARBON	27000	620	UG/G	J	m
VPB152-SOIL-D-060614	SO	TOTAL ORGANIC CARBON	16000	570	UG/G	J	m
VPB152-FB-060614	WQ	TOTAL ORGANIC CARBON		1.0*	MG/L	U	bl
*LOQ							

## Attachment A

### Nonconformance Summary Tables

**Table A-1 - Initial Calibration Verification**

ICV	Compound	% R	Limit
WG144515-7 GCMS-D	ACETONE	125.4	80-120%
Associated samples: All samples in the SDG			

**Table A-2 - Lab Blanks**

Blank ID	Compound	Result	QL	Units	Associated Samples
WG144547-1	TOTAL ORGANIC CARBON	0.12	0.50	MG/L	VPB152-FB-060614

**Table A-3 - Matrix Spikes**

Sample ID	Compound	MS % Recovery	MSD % Recovery	Lower Limit	Upper Limit	RPD	RPD Limit
VPB152-SOIL-060614-325-327	TOTAL ORGANIC CARBON	141	126	75	125	11.2	30

**Attachment B****Qualifier Codes and Explanations**

<b>Qualifier</b>	<b>Explanation</b>
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.

## Attachment C

### Reason Codes and Explanations

Reason Code	Explanation
be	Equipment blank contamination
bf	Field blank contamination
bl	Laboratory blank contamination
c	Calibration issue
co	Analyte carryover
d	Reporting limit raised due to chromatographic interference
fd	Field duplicate RPDs
h	Holding times
i	Internal standard areas
k	Estimated Maximum Possible Concentration (EMPC)
l	LCS or OPR recoveries
lc	Labeled compound recovery
ld	Laboratory duplicate RPDs
lp	Laboratory control sample/laboratory control sample duplicate RPDs
m	Matrix spike recovery
md	Matrix spike/matrix spike duplicate RPDs
nb	Negative laboratory blank contamination
p	Chemical preservation issue
r	Dual column RPD
q	Quantitation issue
s	Surrogate recovery
su	Ion suppression
t	Temperature preservation issue
x	Percent solids
y	Serial dilution results
z	ICS results
mc	Method compliance nonconformance

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4056-1  
**Client ID:** 152-060614-320-322  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4056  
**Lab File ID:** D8874.D

**Sample Date:** 06-JUN-14  
**Received Date:** 10-JUN-14  
**Extract Date:** 11-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144515

**Analysis Date:** 11-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 12-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
<b>Carbon Disulfide</b>	J	0.26	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
<b>Acetone</b>	J	5.3	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
<b>Methyl tert-butyl Ether</b>		11	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50



### Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4056-1  
**Client ID:** 152-060614-320-322  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4056  
**Lab File ID:** D8874.D

**Sample Date:** 06-JUN-14  
**Received Date:** 10-JUN-14  
**Extract Date:** 11-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144515

**Analysis Date:** 11-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 12-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		107.	%					
Toluene-d8		115.	%					
1,2-Dichloroethane-d4		105.	%					
Dibromofluoromethane		105.	%					

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4056-4  
**Client ID:** VPB152-FB-060614  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4056  
**Lab File ID:** D8875.D

**Sample Date:** 06-JUN-14  
**Received Date:** 10-JUN-14  
**Extract Date:** 11-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144515

**Analysis Date:** 11-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 12-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
Acetone	U	2.5	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
<b>Dibromochloromethane</b>	J	0.51	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50



### Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4056-4  
**Client ID:** VPB152-FB-060614  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4056  
**Lab File ID:** D8875.D

**Sample Date:** 06-JUN-14  
**Received Date:** 10-JUN-14  
**Extract Date:** 11-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144515

**Analysis Date:** 11-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 12-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		107.	%					
Toluene-d8		115.	%					
1,2-Dichloroethane-d4		112.	%					
Dibromofluoromethane		106.	%					

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4056-5  
**Client ID:** 152-060614-345-347  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4056  
**Lab File ID:** D8876.D

**Sample Date:** 06-JUN-14  
**Received Date:** 10-JUN-14  
**Extract Date:** 11-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144515

**Analysis Date:** 11-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 12-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
<b>Carbon Disulfide</b>	J	0.42	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
<b>Acetone</b>	J J	4.4	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50



### Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4056-5  
**Client ID:** 152-060614-345-347  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4056  
**Lab File ID:** D8876.D

**Sample Date:** 06-JUN-14  
**Received Date:** 10-JUN-14  
**Extract Date:** 11-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144515

**Analysis Date:** 11-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 12-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		109.	%					
Toluene-d8		119.	%					
1,2-Dichloroethane-d4		116.	%					
Dibromofluoromethane		110.	%					

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4056-6  
**Client ID:** 152-060914-360-362  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4056  
**Lab File ID:** D8877.D

**Sample Date:** 06-JUN-14  
**Received Date:** 10-JUN-14  
**Extract Date:** 11-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144515

**Analysis Date:** 11-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 12-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
Acetone	J	9.9	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50



## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4056-6  
**Client ID:** 152-060914-360-362  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4056  
**Lab File ID:** D8877.D

**Sample Date:** 06-JUN-14  
**Received Date:** 10-JUN-14  
**Extract Date:** 11-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144515

**Analysis Date:** 11-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 12-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		108.	%					
Toluene-d8		116.	%					
1,2-Dichloroethane-d4		111.	%					
Dibromofluoromethane		107.	%					

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4056-7  
**Client ID:** 152-060914-380-382  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4056  
**Lab File ID:** D8878.D

**Sample Date:** 09-JUN-14  
**Received Date:** 10-JUN-14  
**Extract Date:** 11-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144515

**Analysis Date:** 11-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 12-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
<b>Acetone</b>	<b>J</b>	6.9	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50



### Report of Analytical Results

Client: ENSAFE  
 Lab ID: SH4056-7  
 Client ID: 152-060914-380-382  
 Project: Navy Clean WE15-03-06 NW  
 SDG: SH4056  
 Lab File ID: D8878.D

Sample Date: 09-JUN-14  
 Received Date: 10-JUN-14  
 Extract Date: 11-JUN-14  
 Extracted By: DJP  
 Extraction Method: SW846 5030  
 Lab Prep Batch: WG144515

Analysis Date: 11-JUN-14  
 Analyst: DJP  
 Analysis Method: SW846 8260B  
 Matrix: AQ  
 % Solids: NA  
 Report Date: 12-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		108.	%					
Toluene-d8		117.	%					
1,2-Dichloroethane-d4		114.	%					
Dibromofluoromethane		104.	%					

### Report of Analytical Results

Client: ENSAFE  
 Lab ID: SH4056-8  
 Client ID: VPB152-TB-060914  
 Project: Navy Clean WE15-03-06 NW  
 SDG: SH4056  
 Lab File ID: D8873.D

Sample Date: 09-JUN-14  
 Received Date: 10-JUN-14  
 Extract Date: 11-JUN-14  
 Extracted By: DJP  
 Extraction Method: SW846 5030  
 Lab Prep Batch: WG144515

Analysis Date: 11-JUN-14  
 Analyst: DJP  
 Analysis Method: SW846 8260B  
 Matrix: AQ  
 % Solids: NA  
 Report Date: 12-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
Acetone	U	2.5	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50



### Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4056-8  
**Client ID:** VPB152-TB-060914  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4056  
**Lab File ID:** D8873.D

**Sample Date:** 09-JUN-14  
**Received Date:** 10-JUN-14  
**Extract Date:** 11-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144515

**Analysis Date:** 11-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 12-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		106.	%					
Toluene-d8		114.	%					
1,2-Dichloroethane-d4		105.	%					
Dibromofluoromethane		102.	%					

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4056-9  
**Client ID:** 152-060914-400-402  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4056  
**Lab File ID:** D8879.D

**Sample Date:** 09-JUN-14  
**Received Date:** 10-JUN-14  
**Extract Date:** 11-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144515

**Analysis Date:** 11-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 12-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
Acetone	U	2.5	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50

### Report of Analytical Results

Client: ENSAFE  
 Lab ID: SH4056-9  
 Client ID: 152-060914-400-402  
 Project: Navy Clean WE15-03-06 NW  
 SDG: SH4056  
 Lab File ID: D8879.D

Sample Date: 09-JUN-14  
 Received Date: 10-JUN-14  
 Extract Date: 11-JUN-14  
 Extracted By: DJP  
 Extraction Method: SW846 5030  
 Lab Prep Batch: WG144515

Analysis Date: 11-JUN-14  
 Analyst: DJP  
 Analysis Method: SW846 8260B  
 Matrix: AQ  
 % Solids: NA  
 Report Date: 12-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		106.	%					
Toluene-d8		115.	%					
1,2-Dichloroethane-d4		117.	%					
Dibromofluoromethane		108.	%					

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4056-10  
**Client ID:** 152-060914-420-422  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4056  
**Lab File ID:** D8880.D

**Sample Date:** 09-JUN-14  
**Received Date:** 10-JUN-14  
**Extract Date:** 12-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144515

**Analysis Date:** 12-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 12-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
Acetone	J	J	3.1	ug/L	1	5	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50



### Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH4056-10  
**Client ID:** 152-060914-420-422  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH4056  
**Lab File ID:** D8880.D

**Sample Date:** 09-JUN-14  
**Received Date:** 10-JUN-14  
**Extract Date:** 12-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144515

**Analysis Date:** 12-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 12-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		109.	%					
Toluene-d8		117.	%					
1,2-Dichloroethane-d4		119.	%					
Dibromofluoromethane		114.	%					

## Report of Analytical Results

**Client:** Rick Purdy  
AECOM  
701 Edgewater Drive  
Wakefield, MA 01880

**Lab Sample ID:** SH4056-2  
**Report Date:** 10-JUL-14  
**Client PO:** 16518  
**Project:** Navy Clean WE15-03-0  
**SDG:** SH4056

Sample Description  
152-060614-325-327

Matrix SL  
Date Sampled 06-JUN-14  
Date Received 10-JUN-14

Parameter	Result	Adj LOQ	Adj MDL	Adj LOD	Anal. Method	QC.Batch	Anal. Date	Prep. Method	Prep. Date	Footnotes
TOC In Soil	27000 ug/gdrywt	620	130	N/A	SW846 9060A Mod.	WG144438	10-JUN-14 16:16:17	N/A	N/A	N/A
Total Solids	64. %	1		N/A	SM2540G	WG144445	11-JUN-14 14:24:36	SM2540G	10-JUN-14	

*Handwritten signature/initials*

### Report of Analytical Results

**Client:** Rick Purdy  
AECOM  
701 Edgewater Drive  
Wakefield, MA 01880

**Lab Sample ID:** SH4056-3  
**Report Date:** 10-JUL-14  
**Client PO:** 16518  
**Project:** Navy Clean WE15-03-0  
**SDG:** SH4056

Sample Description  
152-SOIL-D-060614

Matrix      Date Sampled      Date Received  
SL              06-JUN-14              10-JUN-14

Parameter	Result	Adj LOQ	Adj MDL	Adj LOD	Anal. Method	QC.Batch	Anal. Date	Prep. Method	Prep. Date	Footnotes
TOC In Soil	16000 ug/gdrywt	570	120	N/A	SW846 9060A Mod.	WG144438	10-JUN-14 17:15:33	N/A	N/A	
Total Solids	70. %	1		N/A	SM2540G	WG144445	11-JUN-14 14:24:45	SM2540G	10-JUN-14	

*Rick Purdy*

**Report of Analytical Results**

**Client:** Rick Purdy  
AECOM  
701 Edgewater Drive  
Wakefield, MA 01880

**Lab Sample ID:** SH4056-4  
**Report Date:** 10-JUL-14  
**Client PO:** 16518  
**Project:** Navy Clean WE15-03-0  
**SDG:** SH4056

Sample Description  
VPB152-FB-060614

Matrix AQ      Date Sampled 06-JUN-14      Date Received 10-JUN-14

Parameter	Result	Adj LOQ	Adj MDL	Adj LOD	Anal. Method	QC.Batch	Anal. Date	Prep. Method	Prep. Date	Footnotes
Total Organic Carbon	10.50 mg/L	1.0	0.10	.5	SM5310B	WG144547	11-JUN-14 07:50:02	N/A	N/A	N/A

1.0 U  
Purdy





## Data Validation Report

Project:	Regional Groundwater Investigation - NWIRP Bethpage	
Laboratory:	Katahdin Analytical Services, Scarborough, Maine	
Service Request:	SH3956	
Analyses/Method:	EPA SW-846 Method 8260B for VOCs (GC/MS)	
Validation Level:	Limited	
RESCON Project Number:	60266526.SA.DV	
Prepared by:	Sheena Blair/RESCON	Completed on: 8/30/2014
Reviewed by:	Lori Herberich/RESCON	File Name: SH3956_8260B

### SUMMARY

The samples listed below were collected by Resolution Consultants from the Regional Groundwater Investigation – NWIRP Bethpage site on June 3, 4, and 5, 2014.

Sample ID	Matrix/Sample Type
VPB152-GWD-060314	Field Duplicate of VPB152-GW-060314-150-152
VPB152-GW-060314-150-152	Groundwater
VPB152-GW-060314-58-60	Groundwater
VPB152-GW-060314-98-100	Groundwater
VPB152-GW-060414-200-202	Groundwater
VPB152-GW-060414-220-222	Groundwater
VPB152-GW-060514-240-242	Groundwater
VPB152-GW-060514-260-262	Groundwater
VPB152-GW-060514-280-282	Groundwater
VPB152-GW-060514-300-302	Groundwater
VPB152-TRIP BLANK-060514	Trip Blank

Data validation activities were conducted with reference to *Test Methods for Evaluating Solid Waste, Physical/Chemical Methods SW-846, specifically SW-846 Method 8260B, Volatile Organic Compounds by Gas Chromatography/Mass Spectrometry* (USEPA, 1996), *USEPA Contract Laboratory Program National Functional Guidelines for Superfund Organic Methods Data Review* (June 2008), and *Quality Systems Manual (QSM) for Environmental Laboratories, Version 4.2* (DoD, 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 review elements (where applicable to the method):

- X Data completeness (chain-of-custody [COC])/sample integrity

- ✓ Holding times and sample preservation
- ✓ GC/MS performance checks
- ✗ Initial calibration/continuing calibration verification
- ✓ Laboratory blanks/ equipment blanks/trip blanks
- ✗ Surrogate spike recoveries
- ✗ Matrix spike (MS) and/or matrix spike duplicate (MSD) results
- ✓ Laboratory control sample (LCS) results
- ✓ Field duplicate results
- ✓ Internal standard results
- ✓ 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. Selected data points were estimated due to nonconformances of certain QC criteria (see discussion below). Qualified sample results are presented in Table 1.

## 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 "VPB" prefix and the "GW" from the sample ID, and truncated the ID for the Trip Blank in the report. The submitted EDD file reflects the full sample ID.

Sample VPB152-GW-060314-150-152 was mostly soil and had very little standing water. The laboratory decanted the water from three individual vials into one vial as a composite. Positive and nondetect results for this sample were qualified as estimated (J and UJ, respectively), due to possible loss of sample integrity during the decanting procedure. Qualified sample results are presented in Table 1.

### 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.

**GC/MS Performance Checks**

The data were reviewed to ensure that the 4-bromofluorobenzene (BFB) tuning was performed at the correct frequency and that the method acceptance criteria were met. 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:

- the initial calibration (ICAL) percent relative standard deviation (%RSD), correlation coefficient (r)/coefficient of determination ( $r^2$ ), and/or response factor method acceptance criteria were met;
- the initial calibration verification standard (ICV) percent recoveries (%Rs) acceptance criteria were met;
- the continuing calibration verification standard (CCV) method percent difference or percent drift (%Ds) and RF acceptance criteria were met; and
- the retention time method acceptance criteria were met.

Data qualification to the analytes associated with the specific ICAL,ICV and/or CCV was as follows:

**ICAL Linearity Nonconformances:**

Nonconformance	Actions	
	Detected Results	Nondetected Results
%RSD > 15% and quantitation based on mean RF	J	UJ
r or $r^2$ < 0.99 and quantitation based on linear regression	J*	UJ*
* No guidance in NFG, professional judgment was used		

**ICV Recovery Nonconformances:**

Nonconformance	Actions	
	Detected Compounds	Nondetected Compounds
%R > 120%	J	No qualification
20% < %R < 80%	J	UJ
%R < 20% (see note)	J	R*

Notes: Based on NFG 2008 VOC guidance, professional judgment is used to reject (R) nondetects in all associated samples for any analyte with < 20% recovery. Also, professional judgment is used to estimate (UJ) rather the reject (R) sample results previously negated (U) on the basis of blank contamination.

**CCV Linearity Nonconformances:**

Nonconformance	Actions	
	Detected Results	Nondetected Results
%D > 20%	J	UJ
%Drift > 20%	J*	UJ*
* No guidance in NFG, thus professional judgment was used		

Qualified sample results are shown in Table 1. Nonconformances are summarized in Attachment A in Tables A-1a, A-1b, A-1c, A-1d and A-1e.

### **Laboratory Blanks/Equipment Blanks/Trip Blanks**

Laboratory method blanks, equipment rinsate and trip 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, equipment rinsate and trip 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.

### **Surrogate Spike Recoveries**

The surrogate recoveries (%Rs) were reviewed for conformance with the QC acceptance criteria.

Data qualification on the basis of surrogate recovery nonconformances was as follows:

Nonconformance	Action	
	Detected Compounds	Nondetected Compounds
%R > Upper Limit (UL)	J	No qualification
20% < %R < Lower Limit (LL)	J	UJ
%R < 20%	J	R

Nonconformances are summarized in Attachment A in Table A-2. Qualified sample results are shown in Table 1.

Sample VPB152-GW-060314-150-152 had two surrogates out, one with a high bias and one with a low bias. Positive and nondetect compounds for this sample were qualified, as estimated (J and UJ, respectively), as the bias is undetermined.

### **MS/MSD Results**

The MS/MSD %Rs and relative percent differences (RPDs) were reviewed for conformance with the QC acceptance criteria.

For the MS/MSD performed on sample VPB152-GW-060414-200-202, over 50% of the compounds did not meet the QC acceptance criteria. The validator chose to qualify all compounds reported from this sample. Positive and nondetects were qualified, as estimated (J and UJ, respectively).

Nonconformances are summarized in Attachment A in Table A-3. Qualified sample results are shown in Table 1.

### **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  $\leq 30\%$  for aqueous matrices. This criteria applies if both results were greater than five times the Limit of Quantitation (LOQ). All QC acceptance criteria were met.

**Internal Standard Results**

The internal standard (IS) recoveries were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

**Sample Results/Reporting Issues**

Compounds that were not detected in the sample are reported as undetected (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**

Sample results qualified as a result of validation actions are summarized in Table 1. All actions are described above.

**ATTACHMENTS**

Attachment A: Nonconformance Summary Tables

Attachment B: Qualifier Codes and Explanations

Attachment C: Reason Codes and Explanations

**Table 1 - Data Validation Summary of Qualified Data**

Sample ID	Matrix	Compound	Result	LOD	Units	Validation Qualifiers	Validation Reason
VPB152-GW-060314-150-152	WG	1,1,1-TRICHLOROETHANE	2.2	0.50	UG/L	J	s
VPB152-GW-060314-150-152	WG	1,1,2,2-TETRACHLOROETHANE		0.50	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE		0.50	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	1,1,2-TRICHLOROETHANE		0.50	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	1,1-DICHLOROETHANE	2.0	0.50	UG/L	J	s
VPB152-GW-060314-150-152	WG	1,1-DICHLOROETHENE	0.61	0.50	UG/L	J	s
VPB152-GW-060314-150-152	WG	1,2,4-TRICHLOROBENZENE		0.50	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	1,2-DIBROMO-3-CHLOROPROPANE		0.75	UG/L	UJ	c,s
VPB152-GW-060314-150-152	WG	1,2-DIBROMOETHANE		0.50	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	1,2-DICHLOROBENZENE		0.50	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	1,2-DICHLOROETHANE		0.50	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	1,2-DICHLOROETHENE, TOTAL		1.0	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	1,2-DICHLOROPROPANE		0.50	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	1,3-DICHLOROBENZENE		0.50	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	1,4-DICHLOROBENZENE		0.50	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	2-BUTANONE		2.5	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	2-HEXANONE		2.5	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	4-METHYL-2-PENTANONE		2.5	UG/L	UJ	c,s
VPB152-GW-060314-150-152	WG	ACETONE	2.7	2.5	UG/L	J	c,s
VPB152-GW-060314-150-152	WG	BENZENE		0.50	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	BROMODICHLOROMETHANE		0.50	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	BROMOFORM		0.50	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	BROMOMETHANE		1.0	UG/L	UJ	c,s
VPB152-GW-060314-150-152	WG	CARBON DISULFIDE		0.50	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	CARBON TETRACHLORIDE		0.50	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	CHLOROBENZENE		0.50	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	CHLOROETHANE		1.0	UG/L	UJ	c,s
VPB152-GW-060314-150-152	WG	CHLOROFORM		0.50	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	CHLOROMETHANE		1.0	UG/L	UJ	c,s
VPB152-GW-060314-150-152	WG	CIS-1,2-DICHLOROETHENE		0.50	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	CIS-1,3-DICHLOROPROPENE		0.50	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	CYCLOHEXANE		0.50	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	DIBROMOCHLOROMETHANE		0.50	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	DICHLORODIFLUOROMETHANE		1.0	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	ETHYLBENZENE		0.50	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	ISOPROPYLBENZENE		0.50	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	M- AND P-XYLENE		1.0	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	METHYL ACETATE		0.75	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	METHYL CYCLOHEXANE		0.50	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	METHYL TERT-BUTYL ETHER		0.50	UG/L	UJ	c,s
VPB152-GW-060314-150-152	WG	METHYLENE CHLORIDE		2.5	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	O-XYLENE		0.50	UG/L	UJ	s

Sample ID	Matrix	Compound	Result	LOD	Units	Validation Qualifiers	Validation Reason
VPB152-GW-060314-150-152	WG	STYRENE		0.50	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	TETRACHLOROETHENE		0.50	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	TOLUENE		0.50	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	TRANS-1,2-DICHLOROETHENE		0.50	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	TRANS-1,3-DICHLOROPROPENE		0.50	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	TRICHLOROETHENE		0.50	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	TRICHLOROFLUOROMETHANE		1.0	UG/L	UJ	c,s
VPB152-GW-060314-150-152	WG	VINYL CHLORIDE		1.0	UG/L	UJ	s
VPB152-GW-060314-150-152	WG	XYLENES, TOTAL		1.5	UG/L	UJ	s
VPB152-GW-060314-58-60	WG	1,1,1-TRICHLOROETHANE		0.50	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	1,1,2-TETRACHLOROETHANE		0.50	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE		0.50	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	1,1,2-TRICHLOROETHANE		0.50	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	1,1-DICHLOROETHANE		0.50	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	1,1-DICHLOROETHENE		0.50	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	1,2,4-TRICHLOROBENZENE		0.50	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	1,2-DIBROMO-3-CHLOROPROPANE		0.75	UG/L	UJ	c,mc
VPB152-GW-060314-58-60	WG	1,2-DIBROMOETHANE		0.50	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	1,2-DICHLOROBENZENE		0.50	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	1,2-DICHLOROETHANE		0.50	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	1,2-DICHLOROETHENE, TOTAL		1.0	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	1,2-DICHLOROPROPANE		0.50	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	1,3-DICHLOROBENZENE		0.50	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	1,4-DICHLOROBENZENE		0.50	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	2-BUTANONE		2.5	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	2-HEXANONE		2.5	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	4-METHYL-2-PENTANONE		2.5	UG/L	UJ	c,mc
VPB152-GW-060314-58-60	WG	ACETONE	16	2.5	UG/L	J	c,mc,s
VPB152-GW-060314-58-60	WG	BENZENE		0.50	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	BROMODICHLOROMETHANE		0.50	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	BROMOFORM		0.50	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	BROMOMETHANE		1.0	UG/L	UJ	c,mc
VPB152-GW-060314-58-60	WG	CARBON DISULFIDE		0.50	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	CARBON TETRACHLORIDE		0.50	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	CHLOROBENZENE		0.50	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	CHLOROETHANE		1.0	UG/L	UJ	c,mc
VPB152-GW-060314-58-60	WG	CHLOROFORM		0.50	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	CHLOROMETHANE		1.0	UG/L	UJ	c,mc
VPB152-GW-060314-58-60	WG	CIS-1,2-DICHLOROETHENE		0.50	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	CIS-1,3-DICHLOROPROPENE		0.50	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	CYCLOHEXANE		0.50	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	DIBROMOCHLOROMETHANE	0.34	0.50	UG/L	J	mc,s
VPB152-GW-060314-58-60	WG	DICHLORODIFLUOROMETHANE		1.0	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	ETHYLBENZENE		0.50	UG/L	UJ	mc

Sample ID	Matrix	Compound	Result	LOD	Units	Validation Qualifiers	Validation Reason
VPB152-GW-060314-58-60	WG	ISOPROPYLBENZENE		0.50	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	M- AND P-XYLENE		1.0	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	METHYL ACETATE		0.75	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	METHYL CYCLOHEXANE		0.50	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	METHYL TERT-BUTYL ETHER		0.50	UG/L	UJ	c,mc
VPB152-GW-060314-58-60	WG	METHYLENE CHLORIDE		2.5	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	O-XYLENE		0.50	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	STYRENE		0.50	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	TETRACHLOROETHENE		0.50	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	TOLUENE		0.50	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	TRANS-1,2-DICHLOROETHENE		0.50	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	TRANS-1,3-DICHLOROPROPENE		0.50	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	TRICHLOROETHENE		0.50	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	TRICHLOROFUOROMETHANE		1.0	UG/L	UJ	c,mc
VPB152-GW-060314-58-60	WG	VINYL CHLORIDE		1.0	UG/L	UJ	mc
VPB152-GW-060314-58-60	WG	XYLENES, TOTAL		1.5	UG/L	UJ	mc
VPB152-GW-060314-98-100	WG	1,2-DIBROMO-3-CHLOROPROPANE		0.75	UG/L	UJ	c
VPB152-GW-060314-98-100	WG	2-HEXANONE		2.5	UG/L	UJ	c
VPB152-GW-060314-98-100	WG	4-METHYL-2-PENTANONE		2.5	UG/L	UJ	c
VPB152-GW-060314-98-100	WG	ACETONE	4.4	2.5	UG/L	J	c
VPB152-GW-060414-200-202	WG	1,1,1-TRICHLOROETHANE		0.50	UG/L	UJ	m
VPB152-GW-060414-200-202	WG	1,1,2,2-TETRACHLOROETHANE		0.50	UG/L	UJ	m,md
VPB152-GW-060414-200-202	WG	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE		0.50	UG/L	UJ	m
VPB152-GW-060414-200-202	WG	1,1,2-TRICHLOROETHANE		0.50	UG/L	UJ	m
VPB152-GW-060414-200-202	WG	1,1-DICHLOROETHANE	0.93	0.50	UG/L	J	m
VPB152-GW-060414-200-202	WG	1,1-DICHLOROETHENE		0.50	UG/L	UJ	m
VPB152-GW-060414-200-202	WG	1,2,4-TRICHLOROBENZENE		0.50	UG/L	UJ	m
VPB152-GW-060414-200-202	WG	1,2-DIBROMO-3-CHLOROPROPANE		0.75	UG/L	U	m
VPB152-GW-060414-200-202	WG	1,2-DIBROMOETHANE		0.50	UG/L	UJ	m
VPB152-GW-060414-200-202	WG	1,2-DICHLOROBENZENE		0.50	UG/L	UJ	m
VPB152-GW-060414-200-202	WG	1,2-DICHLOROETHANE		0.50	UG/L	UJ	m
VPB152-GW-060414-200-202	WG	1,2-DICHLOROETHENE, TOTAL		1.0	UG/L	UJ	m
VPB152-GW-060414-200-202	WG	1,2-DICHLOROPROPANE		0.50	UG/L	UJ	m
VPB152-GW-060414-200-202	WG	1,3-DICHLOROBENZENE		0.50	UG/L	UJ	m
VPB152-GW-060414-200-202	WG	1,4-DICHLOROBENZENE		0.50	UG/L	UJ	m
VPB152-GW-060414-200-202	WG	2-BUTANONE		2.5	UG/L	U	m
VPB152-GW-060414-200-202	WG	2-HEXANONE		2.5	UG/L	UJ	m
VPB152-GW-060414-200-202	WG	4-METHYL-2-PENTANONE		2.5	UG/L	UJ	m
VPB152-GW-060414-200-202	WG	ACETONE	8.2	2.5	UG/L	J	m
VPB152-GW-060414-200-202	WG	BENZENE		0.50	UG/L	UJ	m
VPB152-GW-060414-200-202	WG	BROMODICHLOROMETHANE		0.50	UG/L	UJ	m
VPB152-GW-060414-200-202	WG	BROMOFORM		0.50	UG/L	UJ	m
VPB152-GW-060414-200-202	WG	BROMOMETHANE		1.0	UG/L	U	m



Sample ID	Matrix	Compound	Result	LOD	Units	Validation Qualifiers	Validation Reason
VPB152-GW-060414-200-202	WG	CARBON DISULFIDE		0.50	UG/L	U	m
VPB152-GW-060414-200-202	WG	CARBON TETRACHLORIDE		0.50	UG/L	UJ	m
VPB152-GW-060414-200-202	WG	CHLOROBENZENE		0.50	UG/L	UJ	m
VPB152-GW-060414-200-202	WG	CHLOROETHANE		1.0	UG/L	U	m
VPB152-GW-060414-200-202	WG	CHLOROFORM		0.50	UG/L	UJ	m
VPB152-GW-060414-200-202	WG	CHLOROMETHANE		1.0	UG/L	U	m
VPB152-GW-060414-200-202	WG	CIS-1,2-DICHLOROETHENE		0.50	UG/L	UJ	m
VPB152-GW-060414-200-202	WG	CIS-1,3-DICHLOROPROPENE		0.50	UG/L	UJ	m
VPB152-GW-060414-200-202	WG	CYCLOHEXANE		0.50	UG/L	UJ	m
VPB152-GW-060414-200-202	WG	DIBROMOCHLOROMETHANE		0.50	UG/L	UJ	m
VPB152-GW-060414-200-202	WG	DICHLORODIFLUOROMETHANE		1.0	UG/L	U	m
VPB152-GW-060414-200-202	WG	ETHYLBENZENE		0.50	UG/L	UJ	m
VPB152-GW-060414-200-202	WG	ISOPROPYLBENZENE		0.50	UG/L	UJ	m
VPB152-GW-060414-200-202	WG	M- AND P-XYLENE		1.0	UG/L	UJ	m
VPB152-GW-060414-200-202	WG	METHYL ACETATE		0.75	UG/L	UJ	m
VPB152-GW-060414-200-202	WG	METHYL CYCLOHEXANE		0.50	UG/L	UJ	m,md
VPB152-GW-060414-200-202	WG	METHYL TERT-BUTYL ETHER	6.0	0.50	UG/L	J	m
VPB152-GW-060414-200-202	WG	METHYLENE CHLORIDE		2.5	UG/L	U	m
VPB152-GW-060414-200-202	WG	O-XYLENE		0.50	UG/L	UJ	m
VPB152-GW-060414-200-202	WG	STYRENE		0.50	UG/L	UJ	m
VPB152-GW-060414-200-202	WG	TETRACHLOROETHENE		0.50	UG/L	U	m
VPB152-GW-060414-200-202	WG	TOLUENE		0.50	UG/L	UJ	m
VPB152-GW-060414-200-202	WG	TRANS-1,2-DICHLOROETHENE		0.50	UG/L	UJ	m
VPB152-GW-060414-200-202	WG	TRANS-1,3-DICHLOROPROPENE		0.50	UG/L	UJ	m
VPB152-GW-060414-200-202	WG	TRICHLOROETHENE		0.50	UG/L	UJ	m
VPB152-GW-060414-200-202	WG	TRICHLOROFLUOROMETHANE		1.0	UG/L	UJ	m
VPB152-GW-060414-200-202	WG	VINYL CHLORIDE		1.0	UG/L	U	m
VPB152-GW-060414-200-202	WG	XYLENES, TOTAL		1.5	UG/L	UJ	m
VPB152-GW-060414-220-222	WG	1,2-DIBROMO-3-CHLOROPROPANE		0.75	UG/L	UJ	c
VPB152-GW-060414-220-222	WG	4-METHYL-2-PENTANONE		2.5	UG/L	UJ	c
VPB152-GW-060414-220-222	WG	ACETONE		2.5	UG/L	UJ	c
VPB152-GW-060414-220-222	WG	BROMOMETHANE		1.0	UG/L	UJ	c
VPB152-GW-060414-220-222	WG	CHLOROETHANE		1.0	UG/L	UJ	c
VPB152-GW-060414-220-222	WG	CHLOROMETHANE		1.0	UG/L	UJ	c
VPB152-GW-060414-220-222	WG	METHYL TERT-BUTYL ETHER	0.71	0.50	UG/L	J	c,s
VPB152-GW-060414-220-222	WG	TRICHLOROFLUOROMETHANE		1.0	UG/L	UJ	c
VPB152-GW-060514-240-242	WG	1,2-DIBROMO-3-CHLOROPROPANE		0.75	UG/L	UJ	c
VPB152-GW-060514-240-242	WG	2-HEXANONE		2.5	UG/L	UJ	c
VPB152-GW-060514-240-242	WG	4-METHYL-2-PENTANONE		2.5	UG/L	UJ	c
VPB152-GW-060514-240-242	WG	ACETONE	7.8	2.5	UG/L	J	c
VPB152-GW-060514-260-262	WG	1,2-DIBROMO-3-CHLOROPROPANE		0.75	UG/L	UJ	c
VPB152-GW-060514-260-262	WG	2-HEXANONE		2.5	UG/L	UJ	c
VPB152-GW-060514-260-262	WG	4-METHYL-2-PENTANONE		2.5	UG/L	UJ	c

Sample ID	Matrix	Compound	Result	LOD	Units	Validation Qualifiers	Validation Reason
VPB152-GW-060514-260-262	WG	ACETONE	12	2.5	UG/L	J	c
VPB152-GW-060514-280-282	WG	1,2-DIBROMO-3-CHLOROPROPANE		0.75	UG/L	UJ	c
VPB152-GW-060514-280-282	WG	2-HEXANONE		2.5	UG/L	UJ	c
VPB152-GW-060514-280-282	WG	4-METHYL-2-PENTANONE		2.5	UG/L	UJ	c
VPB152-GW-060514-280-282	WG	ACETONE	4.0	2.5	UG/L	J	c
VPB152-GW-060514-300-302	WG	1,2-DIBROMO-3-CHLOROPROPANE		0.75	UG/L	UJ	c
VPB152-GW-060514-300-302	WG	2-HEXANONE		2.5	UG/L	UJ	c
VPB152-GW-060514-300-302	WG	4-METHYL-2-PENTANONE		2.5	UG/L	UJ	c
VPB152-GW-060514-300-302	WG	ACETONE	3.2	2.5	UG/L	J	c

## Attachment A

## Nonconformance Summary Tables

Table A-1a - Initial Calibration Verification

ICAL	Compound	% RSD	Limit
WG143501 5/27/2014	CHLOROETHANE	17.2	< 15%
Associated samples: VPB152-GW-060314-58-60, VPB152-GW-060314-150-152, VPB152-GW-060414-220-222			

Table A-1b - Initial Calibration Verification

ICV	Compound	% R	Limit
WG143501-7 5/27/2014 GCMS-D	4-METHYL-2-PENTANONE	79.8	80-120%
	1,2-DIBROMO-3-CHLOROPROPANE	78.5	80-120%
Associated samples: VPB152-GW-060314-58-60, VPB152-GW-060314-150-152, VPB152-GW-060414-220-222			

Table A-1c - Initial Calibration Verification

ICV	Compound	% R	Limit
WG144515-7 GCMS-D 6/11/2014	ACETONE	125.4	80-120%
Associated samples: VPB152-GW-060514-300-302, VPB152-GW-060314-98-100, VPB152-GW-060514-260-262, VPB152-GW-060514-240-242, VPB152-GW-060514-280-282			

Table A-1d - Continuing Calibration Verification

CCV ID	Compound	% D	Limit
WG144345 6/9/2014 GCMS-D	CHLOROMETHANE	21.3	< 20%
	BROMOMETHANE	43.6	< 20%
	CHLOROETHANE	33.7	< 20%
	TRICHLOROFLUOROMETHANE	30.2	< 20%
	ACETONE	37.0	< 20%
	MTBE	23.3	< 20%
Associated samples: VPB152-GW-060314-58-60, VPB152-GW-060314-150-152, VPB152-GW-060414-220-222			

**Table A-1e - Continuing Calibration Verification**

CCV ID	Compound	% D	Limit
WG145718 GCMS-D 6/14/2014	4-METHYL-2-PENTANONE	21.2	< 20%
	2-HEXNONE	27.7	< 20%
	1,2-DIBROMO-3-CHLOROPROPANE	22.5	< 20%

Associated samples: VPB152-GW-060514-300-302, VPB152-GW-060314-98-100, VPB152-GW-060514-260-262, VPB152-GW-060514-240-242, VPB152-GW-060514-280-282

**Table A-2 - Surrogates**

Sample ID	Surrogate	% Recovery	Lower Limit	Upper Limit
VPB152-GW-060314-150-152	1,2-DICHLOROETHANE-D4	125	70	120
VPB152-GW-060314-150-152	4-BROMOFLUOROBENZENE	72.7	75	120
VPB152-GW-060314-58-60	1,2-DICHLOROETHANE-D4	127	70	120
VPB152-GW-060414-220-222	1,2-DICHLOROETHANE-D4	134	70	120

**Table A-3 - Matrix Spikes**

Sample ID	Compound	MS % Recovery	MSD % Recovery	Lower Limit	Upper Limit	RPD	RPD Limit
VPB152-GW-060414-200-202	ETHYLBENZENE	51.2	59.6	75	125	15	30
VPB152-GW-060414-200-202	STYRENE	51.6	58.4	65	135	12	30
VPB152-GW-060414-200-202	CIS-1,3-DICHLOROPROPENE	46.8	54.2	70	130	15	30
VPB152-GW-060414-200-202	TRANS-1,3-DICHLOROPROPENE	50	57.8	55	140	14	30
VPB152-GW-060414-200-202	1,4-DICHLOROBENZENE	48	56.8	75	125	17	30
VPB152-GW-060414-200-202	1,2-DIBROMOETHANE	55	62.6	80	120	13	30
VPB152-GW-060414-200-202	1,2-DICHLOROETHANE	57	64	70	130	12	30
VPB152-GW-060414-200-202	4-METHYL-2-PENTANONE	44.8	54.8	60	135	20	30
VPB152-GW-060414-200-202	M- AND P-XYLENE	56.2	66.2	75	130	16	30
VPB152-GW-060414-200-202	METHYL CYCLOHEXANE	97.4	71.2	73	125	31	30
VPB152-GW-060414-200-202	TOLUENE	55	62.8	75	120	13	30
VPB152-GW-060414-200-202	CHLOROBENZENE	56.4	63.6	80	120	12	30
VPB152-GW-060414-200-202	CYCLOHEXANE	53	64.4	71	133	19	30
VPB152-GW-060414-200-202	1,2,4-TRICHLOROBENZENE	45	57.8	65	135	25	30
VPB152-GW-060414-200-202	DIBROMOCHLOROMETHANE	55.8	63.2	60	135	12	30
VPB152-GW-060414-200-202	XYLENES, TOTAL	55	64.7	89	116	16	30
VPB152-GW-060414-200-202	CIS-1,2-DICHLOROETHENE	54.8	63	70	125	14	30
VPB152-GW-060414-200-202	TRANS-1,2-DICHLOROETHENE	53.6	61.8	60	140	14	30
VPB152-GW-060414-200-202	METHYL TERT-BUTYL ETHER	65.1	59.5	65	125	8	30
VPB152-GW-060414-200-202	1,2-DICHLOROETHENE, TOTAL	54.2	62.5	84	121	14	30
VPB152-GW-060414-200-202	1,3-DICHLOROBENZENE	53.8	66.8	75	125	22	30
VPB152-GW-060414-200-202	CARBON TETRACHLORIDE	51.2	63	65	140	21	30
VPB152-GW-060414-200-202	2-HEXANONE	37.6	47.8	55	130	24	30

Sample ID	Compound	MS % Recovery	MSD % Recovery	Lower Limit	Upper Limit	RPD	RPD Limit
VPB152-GW-060414-200-202	CHLOROFORM	54	60	65	135	10	30
VPB152-GW-060414-200-202	BENZENE	54.8	62.6	80	120	13	30
VPB152-GW-060414-200-202	1,1,1-TRICHLOROETHANE	52.4	64.4	65	130	20	30
VPB152-GW-060414-200-202	BROMOFORM	48	54.6	70	130	13	30
VPB152-GW-060414-200-202	BROMODICHLOROMETHANE	58	65.8	75	120	13	30
VPB152-GW-060414-200-202	1,1-DICHLOROETHANE	56.5	62.9	70	135	10	30
VPB152-GW-060414-200-202	1,1-DICHLOROETHENE	54.6	62.8	70	130	14	30
VPB152-GW-060414-200-202	TRICHLOROFLUOROMETHANE	56.8	67.6	60	145	17	30
VPB152-GW-060414-200-202	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	88.8	65	73	126	31	30
VPB152-GW-060414-200-202	1,2-DICHLOROPROPANE	56.4	64.2	75	125	13	30
VPB152-GW-060414-200-202	1,1,2-TRICHLOROETHANE	56.2	64.4	75	125	14	30
VPB152-GW-060414-200-202	TRICHLOROETHENE	55.4	61.8	70	125	11	30
VPB152-GW-060414-200-202	METHYL ACETATE	94	69.6	70	132	30	30
VPB152-GW-060414-200-202	1,1,2,2-TETRACHLOROETHANE	56	64.8	65	130	14	30
VPB152-GW-060414-200-202	O-XYLENE	52.6	61.8	80	120	16	30
VPB152-GW-060414-200-202	1,2-DICHLOROBENZENE	55.6	67.2	70	120	19	30
VPB152-GW-060414-200-202	ISOPROPYLBENZENE	58.4	73	75	125	22	30

**Attachment B****Qualifier Codes and Explanations**

<b>Qualifier</b>	<b>Explanation</b>
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.

### Attachment C

#### Reason Codes and Explanations

Reason Code	Explanation
be	Equipment blank contamination
bf	Field blank contamination
bl	Laboratory blank contamination
c	Calibration issue
co	Analyte carryover
d	Reporting limit raised due to chromatographic interference
fd	Field duplicate RPDs
h	Holding times
i	Internal standard areas
k	Estimated Maximum Possible Concentration (EMPC)
l	LCS or OPR recoveries
lc	Labeled compound recovery
ld	Laboratory duplicate RPDs
lp	Laboratory control sample/laboratory control sample duplicate RPDs
m	Matrix spike recovery
md	Matrix spike/matrix spike duplicate RPDs
nb	Negative laboratory blank contamination
p	Chemical preservation issue
r	Dual column RPD
q	Quantitation issue
s	Surrogate recovery
su	Ion suppression
t	Temperature preservation issue
x	Percent solids
y	Serial dilution results
z	ICS results
mc	Method compliance nonconformance

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH3956-1  
**Client ID:** 152-060314-58-60  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH3956  
**Lab File ID:** D8824.D

**Sample Date:** 03-JUN-14  
**Received Date:** 06-JUN-14  
**Extract Date:** 09-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144345

**Analysis Date:** 09-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 21-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U <b>UJ</b>	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
<b>Acetone</b>	<b>J</b>	16	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U <b>UJ</b>	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
<b>Dibromochloromethane</b>	<b>J</b> <b>J</b>	0.34	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U <b>UJ</b>	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U <b>I</b>	0.50	ug/L	1	1	1.0	0.21	0.50





## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH3956-1  
**Client ID:** 152-060314-58-60  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH3956  
**Lab File ID:** D8824.D

**Sample Date:** 03-JUN-14  
**Received Date:** 06-JUN-14  
**Extract Date:** 09-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144345

**Analysis Date:** 09-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 21-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U <i>UJ</i>	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		81.5	%					
Toluene-d8		96.6	%					
1,2-Dichloroethane-d4	*	127.	%					
Dibromofluoromethane		101.	%					

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH3956-2RA  
**Client ID:** 152-060314-98-100  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH3956  
**Lab File ID:** D8935.D

**Sample Date:** 03-JUN-14  
**Received Date:** 06-JUN-14  
**Extract Date:** 14-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144718

**Analysis Date:** 14-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 21-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
<b>Acetone</b>	<del>U</del> J	4.4	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
<b>1,1-Dichloroethane</b>		1.8	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	<del>U</del> UJ	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	<del>U</del> UJ	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH3956-2RA  
**Client ID:** 152-060314-98-100  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH3956  
**Lab File ID:** D8935.D

**Sample Date:** 03-JUN-14  
**Received Date:** 06-JUN-14  
**Extract Date:** 14-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144718

**Analysis Date:** 14-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 21-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropanoic acid	U UJ	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		100.	%					
Toluene-d8		113.	%					
1,2-Dichloroethane-d4		111.	%					
Dibromofluoromethane		106.	%					

G. 6/14/14

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH3956-3  
**Client ID:** 152-060314-150-152  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH3956  
**Lab File ID:** D8826.D

**Sample Date:** 03-JUN-14  
**Received Date:** 06-JUN-14  
**Extract Date:** 09-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144345

**Analysis Date:** 09-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 21-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
<b>1,1-Dichloroethene</b>	J	0.61	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
<b>Acetone</b>	J	2.7	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
<b>1,1-Dichloroethane</b>	J	2.0	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
<b>1,1,1-Trichloroethane</b>	J	2.2	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH3956-3  
**Client ID:** 152-060314-150-152  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH3956  
**Lab File ID:** D8826.D

**Sample Date:** 03-JUN-14  
**Received Date:** 06-JUN-14  
**Extract Date:** 09-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144345

**Analysis Date:** 09-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 21-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U <b>0.5</b>	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene	*	72.7	%					
Toluene-d8		91.2	%					
1,2-Dichloroethane-d4	*	125.	%					
Dibromofluoromethane		102.	%					

*Q 6/11/14*

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH3956-4RA  
**Client ID:** VPB152-GWD-060314  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH3956  
**Lab File ID:** D8899.D

**Sample Date:** 03-JUN-14  
**Received Date:** 06-JUN-14  
**Extract Date:** 12-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144586

**Analysis Date:** 12-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 21-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
<b>1,1-Dichloroethene</b>	J	0.54	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
<b>Acetone</b>	J	2.8	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
<b>1,1-Dichloroethane</b>		1.9	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
<b>1,1,1-Trichloroethane</b>		1.9	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH3956-4RA  
**Client ID:** VPB152-GWD-060314  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH3956  
**Lab File ID:** D8899.D

**Sample Date:** 03-JUN-14  
**Received Date:** 06-JUN-14  
**Extract Date:** 12-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144586

**Analysis Date:** 12-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 21-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		99.2	%					
Toluene-d8		110.	%					
1,2-Dichloroethane-d4		117.	%					
Dibromofluoromethane		108.	%					

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH3956-5RA  
**Client ID:** 152-060414-200-202  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH3956  
**Lab File ID:** D8900.D

**Sample Date:** 04-JUN-14  
**Received Date:** 06-JUN-14  
**Extract Date:** 12-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144586

**Analysis Date:** 12-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 21-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	UM	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	UMM	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	UM	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
<b>Acetone</b>		8.2	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	UM	0.50	ug/L	1	1	1.0	0.25	0.50
<b>Methyl tert-butyl Ether</b>	M	6.0	ug/L	1	1	1.0	0.36	0.50
<b>1,1-Dichloroethane</b>	JMM	0.93	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	UMM	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	UMM	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	UMM	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	UMM	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	UMM	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	UMM	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	UMM	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	UMM	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	UMM	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	UMM	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	UMM	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	UMM	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	UMM	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	UM	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	UMM	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	UM	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	UMM	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	UMM	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	UMM	0.50	ug/L	1	1	1.0	0.21	0.50





## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH3956-5RA  
**Client ID:** 152-060414-200-202  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH3956  
**Lab File ID:** D8900.D

**Sample Date:** 04-JUN-14  
**Received Date:** 06-JUN-14  
**Extract Date:** 12-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144586

**Analysis Date:** 12-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 21-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	UMM <i>UJ</i>	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	UMM	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	UMM	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	UMM	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	UMM	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	UMM	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	UMM	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	UMM	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	UMM	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	UM	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	UM	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	UMM	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	UMM	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	UMM	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	UMM	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		99.8	%					
Toluene-d8		112.	%					
1,2-Dichloroethane-d4		114.	%					
Dibromofluoromethane		107.	%					

*Paluh*

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH3956-6  
**Client ID:** 152-060414-220-222  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH3956  
**Lab File ID:** D8829.D

**Sample Date:** 04-JUN-14  
**Received Date:** 06-JUN-14  
**Extract Date:** 09-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144345

**Analysis Date:** 09-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 21-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U <i>UJ</i>	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	<del>U</del> <i>UJ</i>	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U <i>UJ</i>	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	<del>U</del> <i>UJ</i>	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
Acetone	<del>U</del> <i>UJ</i>	2.5	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
<b>Methyl tert-butyl Ether</b>	<i>J</i>	0.71	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	<del>U</del> <i>UJ</i>	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50



## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH3956-6  
**Client ID:** 152-060414-220-222  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH3956  
**Lab File ID:** D8829.D

**Sample Date:** 04-JUN-14  
**Received Date:** 06-JUN-14  
**Extract Date:** 09-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144345

**Analysis Date:** 09-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 21-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	<del>U</del> UJ	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		75.7	%					
Toluene-d8		96.0	%					
1,2-Dichloroethane-d4	*	134.	%					
Dibromofluoromethane		107.	%					

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## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH3956-7RA  
**Client ID:** 152-060514-240-242  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH3956  
**Lab File ID:** D8938.D

**Sample Date:** 05-JUN-14  
**Received Date:** 06-JUN-14  
**Extract Date:** 14-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144718

**Analysis Date:** 14-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 21-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
<b>Acetone</b>	J	7.8	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	J UJ	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	J UJ	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50



## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH3956-7RA  
**Client ID:** 152-060514-240-242  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH3956  
**Lab File ID:** D8938.D

**Sample Date:** 05-JUN-14  
**Received Date:** 06-JUN-14  
**Extract Date:** 14-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144718

**Analysis Date:** 14-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 21-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	<del>U</del> UJ	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		102.	%					
Toluene-d8		113.	%					
1,2-Dichloroethane-d4		110.	%					
Dibromofluoromethane		104.	%					

*Revised*

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH3956-8RA  
**Client ID:** 152-060514-260-262  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH3956  
**Lab File ID:** D8936.D

**Sample Date:** 05-JUN-14  
**Received Date:** 06-JUN-14  
**Extract Date:** 14-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144718

**Analysis Date:** 14-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 21-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
<b>Acetone</b>	J	12	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
<b>Methyl tert-butyl Ether</b>		6.1	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	U 45	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U 45	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50

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## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH3956-8RA  
**Client ID:** 152-060514-260-262  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH3956  
**Lab File ID:** D8936.D

**Sample Date:** 05-JUN-14  
**Received Date:** 06-JUN-14  
**Extract Date:** 14-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144718

**Analysis Date:** 14-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 21-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	<del>U</del> UJ	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		105.	%					
Toluene-d8		117.	%					
1,2-Dichloroethane-d4		116.	%					
Dibromofluoromethane		112.	%					

*R. Katahdin*



## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH3956-9RA  
**Client ID:** 152-060514-280-282  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH3956  
**Lab File ID:** D8939.D

**Sample Date:** 05-JUN-14  
**Received Date:** 06-JUN-14  
**Extract Date:** 14-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144718

**Analysis Date:** 14-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 21-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
Acetone	<del>U</del> J	4.0	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	<del>U</del> UJ	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	<del>U</del> UJ	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50





## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH3956-9RA  
**Client ID:** 152-060514-280-282  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH3956  
**Lab File ID:** D8939.D

**Sample Date:** 05-JUN-14  
**Received Date:** 06-JUN-14  
**Extract Date:** 14-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144718

**Analysis Date:** 14-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 21-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	<del>U</del> UJ	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		100.	%					
Toluene-d8		112.	%					
1,2-Dichloroethane-d4		112.	%					
Dibromofluoromethane		108.	%					

*R. J. / 6/14*

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH3956-10RA  
**Client ID:** 152-060514-300-302  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH3956  
**Lab File ID:** D8934.D

**Sample Date:** 05-JUN-14  
**Received Date:** 06-JUN-14  
**Extract Date:** 14-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144718

**Analysis Date:** 14-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 21-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
<b>Acetone</b>	<del>U</del> J	3.2	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
<b>Methyl tert-butyl Ether</b>		6.3	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	<del>U</del> U J	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	<del>U</del> U J	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50

Rulu/14

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH3956-10RA  
**Client ID:** 152-060514-300-302  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH3956  
**Lab File ID:** D8934.D

**Sample Date:** 05-JUN-14  
**Received Date:** 06-JUN-14  
**Extract Date:** 14-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144718

**Analysis Date:** 14-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 21-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	<del>U</del> UJ	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		100.	%					
Toluene-d8		110.	%					
1,2-Dichloroethane-d4		106.	%					
Dibromofluoromethane		102.	%					

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## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH3956-11RA  
**Client ID:** VPB152-TB-060514  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH3956  
**Lab File ID:** D8897.D

**Sample Date:** 05-JUN-14  
**Received Date:** 06-JUN-14  
**Extract Date:** 12-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144586

**Analysis Date:** 12-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 21-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Dichlorodifluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
Chloromethane	U	1.0	ug/L	1	2	2.0	0.36	1.0
Vinyl Chloride	U	1.0	ug/L	1	2	2.0	0.25	1.0
Bromomethane	U	1.0	ug/L	1	2	2.0	0.49	1.0
Chloroethane	U	1.0	ug/L	1	2	2.0	0.55	1.0
Trichlorofluoromethane	U	1.0	ug/L	1	2	2.0	0.24	1.0
1,1-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.35	0.50
Carbon Disulfide	U	0.50	ug/L	1	1	1.0	0.25	0.50
Freon-113	U	0.50	ug/L	1	1	1.0	0.31	0.50
Methylene Chloride	U	2.5	ug/L	1	5	5.0	1.1	2.5
Acetone	U	2.5	ug/L	1	5	5.0	2.2	2.5
trans-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.25	0.50
Methyl tert-butyl Ether	U	0.50	ug/L	1	1	1.0	0.36	0.50
1,1-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.21	0.50
cis-1,2-Dichloroethene	U	0.50	ug/L	1	1	1.0	0.21	0.50
Chloroform	U	0.50	ug/L	1	1	1.0	0.32	0.50
1,1,1-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
2-Butanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
Cyclohexane	U	0.50	ug/L	1	1	1.0	0.31	0.50
Carbon Tetrachloride	U	0.50	ug/L	1	1	1.0	0.22	0.50
Benzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,2-Dichloroethane	U	0.50	ug/L	1	1	1.0	0.20	0.50
Trichloroethene	U	0.50	ug/L	1	1	1.0	0.28	0.50
1,2-Dichloropropane	U	0.50	ug/L	1	1	1.0	0.25	0.50
Bromodichloromethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
cis-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.19	0.50
Toluene	U	0.50	ug/L	1	1	1.0	0.27	0.50
4-Methyl-2-Pentanone	U	2.5	ug/L	1	5	5.0	1.3	2.5
trans-1,3-Dichloropropene	U	0.50	ug/L	1	1	1.0	0.20	0.50
1,1,2-Trichloroethane	U	0.50	ug/L	1	1	1.0	0.33	0.50
Tetrachloroethene	U	0.50	ug/L	1	1	1.0	0.40	0.50
Dibromochloromethane	U	0.50	ug/L	1	1	1.0	0.30	0.50
2-Hexanone	U	2.5	ug/L	1	5	5.0	1.7	2.5
Chlorobenzene	U	0.50	ug/L	1	1	1.0	0.22	0.50
Ethylbenzene	U	0.50	ug/L	1	1	1.0	0.21	0.50

## Report of Analytical Results

**Client:** ENSAFE  
**Lab ID:** SH3956-11RA  
**Client ID:** VPB152-TB-060514  
**Project:** Navy Clean WE15-03-06 NW  
**SDG:** SH3956  
**Lab File ID:** D8897.D

**Sample Date:** 05-JUN-14  
**Received Date:** 06-JUN-14  
**Extract Date:** 12-JUN-14  
**Extracted By:** DJP  
**Extraction Method:** SW846 5030  
**Lab Prep Batch:** WG144586

**Analysis Date:** 12-JUN-14  
**Analyst:** DJP  
**Analysis Method:** SW846 8260B  
**Matrix:** AQ  
**% Solids:** NA  
**Report Date:** 21-JUN-14

Compound	Qualifier	Result	Units	Dilution	LOQ	ADJ LOQ	ADJ MDL	ADJ LOD
Xylenes (total)	U	1.5	ug/L	1	3	3.0	0.25	1.5
Styrene	U	0.50	ug/L	1	1	1.0	0.23	0.50
Bromoform	U	0.50	ug/L	1	1	1.0	0.23	0.50
Isopropylbenzene	U	0.50	ug/L	1	1	1.0	0.23	0.50
1,1,2,2-Tetrachloroethane	U	0.50	ug/L	1	1	1.0	0.38	0.50
1,3-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.26	0.50
1,4-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.24	0.50
1,2-Dichlorobenzene	U	0.50	ug/L	1	1	1.0	0.15	0.50
1,2,4-Trichlorobenzene	U	0.50	ug/L	1	1	1.0	0.37	0.50
Methyl Acetate	U	0.75	ug/L	1	1	1.0	0.53	0.75
Methylcyclohexane	U	0.50	ug/L	1	1	1.0	0.30	0.50
o-Xylene	U	0.50	ug/L	1	1	1.0	0.25	0.50
M+P-Xylenes	U	1.0	ug/L	1	2	2.0	0.59	1.0
1,2-Dichloroethylene (Total)	U	1.0	ug/L	1	2	2.0	0.21	1.0
1,2-Dibromoethane	U	0.50	ug/L	1	1	1.0	0.22	0.50
1,2-Dibromo-3-Chloropropane	U	0.75	ug/L	1	1	1.0	0.50	0.75
P-Bromofluorobenzene		106.	%					
Toluene-d8		119.	%					
1,2-Dichloroethane-d4	*	124.	%					
Dibromofluoromethane	*	115.	%					

**Section 5**

**VPB 152 Analytical Data Table**

Location	VPB152	VPB152	VPB152	VPB152	
Sample Date	NYSDEC Groundwater Guidance or Standard Value (Note 1)	6/3/2014	6/3/2014	6/3/2014	6/3/2014
Sample ID	VPB152-GW-060314- 58-60	VPB152-GW-060314- 98-100	VPB152-GW-060314- 150-152	VPB152-GWD-060314	
Sample Interval	58 - 60 ft	98 - 100 ft	150 - 152 ft	150 - 152 ft	
Sample type code	N	N	N	FD	
VOC 8260C (ug/L)					
1,1,1-TRICHLOROETHANE	5	< 0.50 UJ	< 0.50 U	<b>2.2 J</b>	<b>1.9</b>
1,1,2,2-TETRACHLOROETHANE	5	< 0.50 UJ	< 0.50 U	< 0.50 UJ	< 0.50 U
1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	5	< 0.50 UJ	< 0.50 U	< 0.50 UJ	< 0.50 U
1,1,2-TRICHLOROETHANE	1	< 0.50 UJ	< 0.50 U	< 0.50 UJ	< 0.50 U
1,1-DICHLOROETHANE	5	< 0.50 UJ	<b>1.8</b>	<b>2.0 J</b>	<b>1.9</b>
1,1-DICHLOROETHENE	5	< 0.50 UJ	< 0.50 U	<b>0.61 J</b>	<b>0.54 J</b>
1,2,4-TRICHLOROBENZENE	5	< 0.50 UJ	< 0.50 U	< 0.50 UJ	< 0.50 U
1,2-DIBROMO-3-CHLOROPROPANE	0.04	< <b>0.75 UJ</b>	< <b>0.75 UJ</b>	< <b>0.75 UJ</b>	< <b>0.75 U</b>
1,2-DIBROMOETHANE	NL	< 0.50 UJ	< 0.50 U	< 0.50 UJ	< 0.50 U
1,2-DICHLOROBENZENE	3	< 0.50 UJ	< 0.50 U	< 0.50 UJ	< 0.50 U
1,2-DICHLOROETHANE	5	< 0.50 UJ	< 0.50 U	< 0.50 UJ	< 0.50 U
1,2-DICHLOROETHENE, TOTAL	5	< 1.0 UJ	< 1.0 U	< 1.0 UJ	< 1.0 U
1,2-DICHLOROPROPANE	1	< 0.50 UJ	< 0.50 U	< 0.50 UJ	< 0.50 U
1,3-DICHLOROBENZENE	3	< 0.50 UJ	< 0.50 U	< 0.50 UJ	< 0.50 U
1,4-DICHLOROBENZENE	3	< 0.50 UJ	< 0.50 U	< 0.50 UJ	< 0.50 U
2-BUTANONE	50	< 2.5 UJ	< 2.5 U	< 2.5 UJ	< 2.5 U
2-HEXANONE	50	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ	< 2.5 U
4-METHYL-2-PENTANONE	NL	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ	< 2.5 U
ACETONE	50	<b>16 J</b>	<b>4.4 J</b>	<b>2.7 J</b>	<b>2.8 J</b>
BENZENE	1	< 0.50 UJ	< 0.50 U	< 0.50 UJ	< 0.50 U
BROMODICHLOROMETHANE	50	< 0.50 UJ	< 0.50 U	< 0.50 UJ	< 0.50 U
BROMOFORM	50	< 0.50 UJ	< 0.50 U	< 0.50 UJ	< 0.50 U
BROMOMETHANE	5	< 1.0 UJ	< 1.0 U	< 1.0 UJ	< 1.0 U
CARBON DISULFIDE	60	< 0.50 UJ	< 0.50 U	< 0.50 UJ	< 0.50 U
CARBON TETRACHLORIDE	5	< 0.50 UJ	< 0.50 U	< 0.50 UJ	< 0.50 U
CHLOROBENZENE	5	< 0.50 UJ	< 0.50 U	< 0.50 UJ	< 0.50 U
CHLOROETHANE	5	< 1.0 UJ	< 1.0 U	< 1.0 UJ	< 1.0 U
CHLOROFORM	7	< 0.50 UJ	< 0.50 U	< 0.50 UJ	< 0.50 U
CHLOROMETHANE	5	< 1.0 UJ	< 1.0 U	< 1.0 UJ	< 1.0 U
CIS-1,2-DICHLOROETHENE	5	< 0.50 UJ	< 0.50 U	< 0.50 UJ	< 0.50 U
CIS-1,3-DICHLOROPROPENE	0.4	< <b>0.50 UJ</b>	< <b>0.50 U</b>	< <b>0.50 UJ</b>	< <b>0.50 U</b>
CYCLOHEXANE	NL	< 0.50 UJ	< 0.50 U	< 0.50 UJ	< 0.50 U
DIBROMOCHLOROMETHANE	5	<b>0.34 J</b>	< 0.50 U	< 0.50 UJ	< 0.50 U
DICHLORODIFLUOROMETHANE	5	< 1.0 UJ	< 1.0 U	< 1.0 UJ	< 1.0 U
ETHYLBENZENE	5	< 0.50 UJ	< 0.50 U	< 0.50 UJ	< 0.50 U
ISOPROPYLBENZENE	5	< 0.50 UJ	< 0.50 U	< 0.50 UJ	< 0.50 U
M- AND P-XYLENE	NL	< 1.0 UJ	< 1.0 U	< 1.0 UJ	< 1.0 U
METHYL ACETATE	NL	< 0.75 UJ	< 0.75 U	< 0.75 UJ	< 0.75 U
METHYL CYCLOHEXANE	NL	< 0.50 UJ	< 0.50 U	< 0.50 UJ	< 0.50 U
METHYL TERT-BUTYL ETHER	10	< 0.50 UJ	< 0.50 U	< 0.50 UJ	< 0.50 U
METHYLENE CHLORIDE	5	< 2.5 UJ	< 2.5 U	< 2.5 UJ	< 2.5 U
O-XYLENE	NL	< 0.50 UJ	< 0.50 U	< 0.50 UJ	< 0.50 U
STYRENE	5	< 0.50 UJ	< 0.50 U	< 0.50 UJ	< 0.50 U
TETRACHLOROETHENE	5	< 0.50 UJ	< 0.50 U	< 0.50 UJ	< 0.50 U
TOLUENE	5	< 0.50 UJ	< 0.50 U	< 0.50 UJ	< 0.50 U
TRANS-1,2-DICHLOROETHENE	5	< 0.50 UJ	< 0.50 U	< 0.50 UJ	< 0.50 U
TRANS-1,3-DICHLOROPROPENE	0.4	< <b>0.50 UJ</b>	< <b>0.50 U</b>	< <b>0.50 UJ</b>	< <b>0.50 U</b>
TRICHLOROETHENE	5	< 0.50 UJ	< 0.50 U	< 0.50 UJ	< 0.50 U
TRICHLOROFLUOROMETHANE	5	< 1.0 UJ	< 1.0 U	< 1.0 UJ	< 1.0 U
VINYL CHLORIDE	2	< 1.0 UJ	< 1.0 U	< 1.0 UJ	< 1.0 U
XYLENES, TOTAL	5	< 1.5 UJ	< 1.5 U	< 1.5 UJ	< 1.5 U

Location		VPB152	VPB152	VPB152	VPB152
Sample Date	NYSDEC	6/4/2014	6/4/2014	6/5/2014	6/5/2014
Sample ID	Groundwater Guidance or Standard Value (Note 1)	VPB152-GW-060414- 200-202	VPB152-GW-060414- 220-222	VPB152-GW-060514- 240-242	VPB152-GW-060514- 260-262
Sample Interval		200 - 202 ft	220 - 222 ft	240 - 242 ft	260 - 262 ft
Sample type code		N	N	N	N
<b>VOC 8260C (ug/L)</b>					
1,1,1-TRICHLOROETHANE	5	< 0.50 UJ	< 0.50 U	< 0.50 U	< 0.50 U
1,1,2,2-TETRACHLOROETHANE	5	< 0.50 UJ	< 0.50 U	< 0.50 U	< 0.50 U
1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	5	< 0.50 UJ	< 0.50 U	< 0.50 U	< 0.50 U
1,1,2-TRICHLOROETHANE	1	< 0.50 UJ	< 0.50 U	< 0.50 U	< 0.50 U
1,1-DICHLOROETHANE	5	<b>0.93 J</b>	< 0.50 U	< 0.50 U	< 0.50 U
1,1-DICHLOROETHENE	5	< 0.50 UJ	< 0.50 U	< 0.50 U	< 0.50 U
1,2,4-TRICHLOROBENZENE	5	< 0.50 UJ	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DIBROMO-3-CHLOROPROPANE	0.04	< <b>0.75 U</b>	< <b>0.75 UJ</b>	< <b>0.75 UJ</b>	< <b>0.75 UJ</b>
1,2-DIBROMOETHANE	NL	< 0.50 UJ	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DICHLOROBENZENE	3	< 0.50 UJ	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DICHLOROETHANE	5	< 0.50 UJ	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DICHLOROETHENE, TOTAL	5	< 1.0 UJ	< 1.0 U	< 1.0 U	< 1.0 U
1,2-DICHLOROPROPANE	1	< 0.50 UJ	< 0.50 U	< 0.50 U	< 0.50 U
1,3-DICHLOROBENZENE	3	< 0.50 UJ	< 0.50 U	< 0.50 U	< 0.50 U
1,4-DICHLOROBENZENE	3	< 0.50 UJ	< 0.50 U	< 0.50 U	< 0.50 U
2-BUTANONE	50	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 U
2-HEXANONE	50	< 2.5 UJ	< 2.5 U	< 2.5 UJ	< 2.5 UJ
4-METHYL-2-PENTANONE	NL	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ
ACETONE	50	<b>8.2 J</b>	< 2.5 UJ	<b>7.8 J</b>	<b>12 J</b>
BENZENE	1	< 0.50 UJ	< 0.50 U	< 0.50 U	< 0.50 U
BROMODICHLOROMETHANE	50	< 0.50 UJ	< 0.50 U	< 0.50 U	< 0.50 U
BROMOFORM	50	< 0.50 UJ	< 0.50 U	< 0.50 U	< 0.50 U
BROMOMETHANE	5	< 1.0 U	< 1.0 UJ	< 1.0 U	< 1.0 U
CARBON DISULFIDE	60	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CARBON TETRACHLORIDE	5	< 0.50 UJ	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROBENZENE	5	< 0.50 UJ	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROETHANE	5	< 1.0 U	< 1.0 UJ	< 1.0 U	< 1.0 U
CHLOROFORM	7	< 0.50 UJ	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROMETHANE	5	< 1.0 U	< 1.0 UJ	< 1.0 U	< 1.0 U
CIS-1,2-DICHLOROETHENE	5	< 0.50 UJ	< 0.50 U	< 0.50 U	< 0.50 U
CIS-1,3-DICHLOROPROPENE	0.4	< <b>0.50 UJ</b>	< <b>0.50 U</b>	< <b>0.50 U</b>	< <b>0.50 U</b>
CYCLOHEXANE	NL	< 0.50 UJ	< 0.50 U	< 0.50 U	< 0.50 U
DIBROMOCHLOROMETHANE	5	< 0.50 UJ	< 0.50 U	< 0.50 U	< 0.50 U
DICHLORODIFLUOROMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
ETHYLBENZENE	5	< 0.50 UJ	< 0.50 U	< 0.50 U	< 0.50 U
ISOPROPYLBENZENE	5	< 0.50 UJ	< 0.50 U	< 0.50 U	< 0.50 U
M- AND P-XYLENE	NL	< 1.0 UJ	< 1.0 U	< 1.0 U	< 1.0 U
METHYL ACETATE	NL	< 0.75 UJ	< 0.75 U	< 0.75 U	< 0.75 U
METHYL CYCLOHEXANE	NL	< 0.50 UJ	< 0.50 U	< 0.50 U	< 0.50 U
METHYL TERT-BUTYL ETHER	10	<b>6.0 J</b>	<b>0.71 J</b>	< 0.50 U	<b>6.1</b>
METHYLENE CHLORIDE	5	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 U
O-XYLENE	NL	< 0.50 UJ	< 0.50 U	< 0.50 U	< 0.50 U
STYRENE	5	< 0.50 UJ	< 0.50 U	< 0.50 U	< 0.50 U
TETRACHLOROETHENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TOLUENE	5	< 0.50 UJ	< 0.50 U	< 0.50 U	< 0.50 U
TRANS-1,2-DICHLOROETHENE	5	< 0.50 UJ	< 0.50 U	< 0.50 U	< 0.50 U
TRANS-1,3-DICHLOROPROPENE	0.4	< <b>0.50 UJ</b>	< <b>0.50 U</b>	< <b>0.50 U</b>	< <b>0.50 U</b>
TRICHLOROETHENE	5	< 0.50 UJ	< 0.50 U	< 0.50 U	< 0.50 U
TRICHLOROFLUOROMETHANE	5	< 1.0 UJ	< 1.0 UJ	< 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 UJ	< 1.5 U	< 1.5 U	< 1.5 U



Location	VPB152	VPB152	VPB152	VPB152	
Sample Date	NYSDEC Groundwater Guidance or Standard Value (Note 1)	6/5/2014	6/5/2014	6/6/2014	6/6/2014
Sample ID		VPB152-GW-060514- 280-282	VPB152-GW-060514- 300-302	VPB152-GW-060614- 320-322	VPB152-GW-060614- 345-347
Sample Interval		280 - 282 ft	300 - 302 ft	320 - 322 ft	345 - 347 ft
Sample type code		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 UJ	< 0.75 UJ	< 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
2-BUTANONE	50	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 U
2-HEXANONE	50	< 2.5 UJ	< 2.5 UJ	< 2.5 U	< 2.5 U
4-METHYL-2-PENTANONE	NL	< 2.5 UJ	< 2.5 UJ	< 2.5 U	< 2.5 U
ACETONE	50	4.0 J	3.2 J	5.3 J	4.4 J
BENZENE	1	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
BROMODICHLOROMETHANE	50	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
BROMOFORM	50	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
BROMOMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
CARBON DISULFIDE	60	< 0.50 U	< 0.50 U	0.26 J	0.42 J
CARBON TETRACHLORIDE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
CHLOROFORM	7	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
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 U	< 0.50 U	< 0.50 U	< 0.50 U
DIBROMOCHLOROMETHANE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
DICHLORODIFLUOROMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
ETHYLBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
ISOPROPYLBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
M- AND P-XYLENE	NL	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
METHYL ACETATE	NL	< 0.75 U	< 0.75 U	< 0.75 U	< 0.75 U
METHYL CYCLOHEXANE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
METHYL TERT-BUTYL ETHER	10	< 0.50 U	6.3	11	< 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 U	< 0.50 U	< 0.50 U	< 0.50 U
TOLUENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TRANS-1,2-DICHLOROETHENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TRANS-1,3-DICHLOROPROPENE	0.4	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TRICHLOROETHENE	5	< 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		VPB152	VPB152	VPB152	VPB152
Sample Date	NYSDEC	6/6/2014	6/9/2014	6/9/2014	6/9/2014
Sample ID	Groundwater Guidance or Standard Value (Note 1)	VPB152-GW-060914- 360-362	VPB152-GW-060914- 380-382	VPB152-GW-060914- 400-402	VPB152-GW-060914- 420-422
Sample Interval		360 - 362 ft	380 - 382 ft	400 - 402 ft	420 - 422 ft
Sample type code		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
2-BUTANONE	50	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 U
2-HEXANONE	50	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 U
4-METHYL-2-PENTANONE	NL	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 U
ACETONE	50	9.9 J	6.9 J	< 2.5 U	3.1 J
BENZENE	1	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
BROMODICHLOROMETHANE	50	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
BROMOFORM	50	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
BROMOMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
CARBON DISULFIDE	60	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CARBON TETRACHLORIDE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
CHLOROFORM	7	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
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 U	< 0.50 U	< 0.50 U	< 0.50 U
DIBROMOCHLOROMETHANE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
DICHLORODIFLUOROMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
ETHYLBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
ISOPROPYLBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
M- AND P-XYLENE	NL	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
METHYL ACETATE	NL	< 0.75 U	< 0.75 U	< 0.75 U	< 0.75 U
METHYL CYCLOHEXANE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
METHYL TERT-BUTYL ETHER	10	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
METHYLENE CHLORIDE	5	< 2.5 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 U	< 0.50 U	< 0.50 U	< 0.50 U
TOLUENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TRANS-1,2-DICHLOROETHENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TRANS-1,3-DICHLOROPROPENE	0.4	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TRICHLOROETHENE	5	< 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	VPB152	VPB152	VPB152	VPB152	
Sample Date	NYSDEC Groundwater Guidance or Standard Value (Note 1)	6/10/2014	6/10/2014	6/10/2014	6/10/2014
Sample ID		VPB152-GW-061014- 440-442	VPB152-GW-061014- 460-462	VPB152-GW-061014- 480-482	VPB152-GW-061014- 500-502
Sample Interval		440 - 442 ft	460 - 462 ft	480 - 482 ft	500 - 502 ft
Sample type code		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
2-BUTANONE	50	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 U
2-HEXANONE	50	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 U
4-METHYL-2-PENTANONE	NL	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 U
ACETONE	50	< 5.0 U	< 2.5 U	< 5.0 U	< 5.0 U
BENZENE	1	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
BROMODICHLOROMETHANE	50	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
BROMOFORM	50	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
BROMOMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
CARBON DISULFIDE	60	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CARBON TETRACHLORIDE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
CHLOROFORM	7	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
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 U	< 0.50 U	< 0.50 U	< 0.50 U
DIBROMOCHLOROMETHANE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
DICHLORODIFLUOROMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
ETHYLBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
ISOPROPYLBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
M- AND P-XYLENE	NL	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
METHYL ACETATE	NL	< 0.75 U	< 0.75 U	< 0.75 U	< 0.75 U
METHYL CYCLOHEXANE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
METHYL TERT-BUTYL ETHER	10	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
METHYLENE CHLORIDE	5	< 2.5 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 U	< 0.50 U	< 0.50 U	< 0.50 U
TOLUENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TRANS-1,2-DICHLOROETHENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TRANS-1,3-DICHLOROPROPENE	0.4	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TRICHLOROETHENE	5	< 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	VPB152	VPB152	VPB152	VPB152	
Sample Date	NYSDEC Groundwater Guidance or Standard Value (Note 1)	6/11/2014	6/11/2014	6/11/2014	6/12/2014
Sample ID		VPB152-GW-061114- 520-522	VPB152-GW-061114- 540-542	VPB152-GW-061114- 560-562	VPB152-GW-061214- 580-582
Sample Interval		520 - 522 ft	540 - 542 ft	560 - 562 ft	580 - 582 ft
Sample type code		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
2-BUTANONE	50	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 U
2-HEXANONE	50	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 U
4-METHYL-2-PENTANONE	NL	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 U
ACETONE	50	< 5.0 U	< 2.5 U	< 5.0 U	< 5.0 U
BENZENE	1	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
BROMODICHLOROMETHANE	50	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
BROMOFORM	50	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
BROMOMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
CARBON DISULFIDE	60	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CARBON TETRACHLORIDE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
CHLOROFORM	7	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
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 U	< 0.50 U	< 0.50 U	< 0.50 U
DIBROMOCHLOROMETHANE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
DICHLORODIFLUOROMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
ETHYLBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
ISOPROPYLBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
M- AND P-XYLENE	NL	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
METHYL ACETATE	NL	< 0.75 U	< 0.75 U	< 0.75 U	< 0.75 U
METHYL CYCLOHEXANE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
METHYL TERT-BUTYL ETHER	10	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
METHYLENE CHLORIDE	5	< 2.5 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 U	< 0.50 U	< 0.50 U	< 0.50 U
TOLUENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TRANS-1,2-DICHLOROETHENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TRANS-1,3-DICHLOROPROPENE	0.4	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TRICHLOROETHENE	5	< 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	VPB152	VPB152	VPB152	VPB152
Sample Date	6/12/2014	6/13/2014	6/13/2014	6/13/2014
Sample ID	NYSDEC Groundwater Guidance or Standard Value (Note 1)	VPB152-GW-061214- 605-607	VPB152-GW-061314- 620-622	VPB152-GW-061314- 640-642
Sample Interval	605 - 607 ft	620 - 622 ft	640 - 642 ft	640 - 642 ft
Sample type code	N	N	N	FD
<b>VOC 8260C (ug/L)</b>				
1,1,1-TRICHLOROETHANE	5	< 0.50 U	< 0.50 UJ	< 0.50 U
1,1,2,2-TETRACHLOROETHANE	5	< 0.50 U	< 0.50 UJ	< 0.50 U
1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	5	< 0.50 U	< 0.50 UJ	< 0.50 U
1,1,2-TRICHLOROETHANE	1	< 0.50 U	< 0.50 UJ	< 0.50 U
1,1-DICHLOROETHANE	5	< 0.50 U	< 0.50 UJ	< 0.50 U
1,1-DICHLOROETHENE	5	< 0.50 U	< 0.50 UJ	< 0.50 U
1,2,4-TRICHLOROBENZENE	5	< 0.50 U	< 0.50 UJ	< 0.50 U
1,2-DIBROMO-3-CHLOROPROPANE	0.04	< <b>0.75 U</b>	< <b>0.75 UJ</b>	< <b>0.75 U</b>
1,2-DIBROMOETHANE	NL	< 0.50 U	< 0.50 UJ	< 0.50 U
1,2-DICHLOROBENZENE	3	< 0.50 U	< 0.50 UJ	< 0.50 U
1,2-DICHLOROETHANE	5	< 0.50 U	< 0.50 UJ	< 0.50 U
1,2-DICHLOROETHENE, TOTAL	5	< 1.0 U	< 1.0 UJ	< 1.0 U
1,2-DICHLOROPROPANE	1	< 0.50 U	< 0.50 UJ	< 0.50 U
1,3-DICHLOROBENZENE	3	< 0.50 U	< 0.50 UJ	< 0.50 U
1,4-DICHLOROBENZENE	3	< 0.50 U	< 0.50 UJ	< 0.50 U
2-BUTANONE	50	< 2.5 U	< 2.5 UJ	< 2.5 UJ
2-HEXANONE	50	< 2.5 U	< 2.5 UJ	< 2.5 U
4-METHYL-2-PENTANONE	NL	< 2.5 U	< 2.5 UJ	< 2.5 UJ
ACETONE	50	< 2.5 U	<b>14 J</b>	<b>7.0 J</b>
BENZENE	1	< 0.50 U	< 0.50 UJ	< 0.50 U
BROMODICHLOROMETHANE	50	< 0.50 U	< 0.50 UJ	< 0.50 U
BROMOFORM	50	< 0.50 U	< 0.50 UJ	< 0.50 U
BROMOMETHANE	5	< 1.0 U	< 1.0 UJ	< 1.0 U
CARBON DISULFIDE	60	< 0.50 U	<b>0.40 J</b>	< 1.0 U
CARBON TETRACHLORIDE	5	< 0.50 U	< 0.50 UJ	< 0.50 U
CHLOROBENZENE	5	< 0.50 U	< 0.50 UJ	< 0.50 U
CHLOROETHANE	5	< 1.0 U	< 1.0 UJ	< 1.0 UJ
CHLOROFORM	7	< 0.50 U	< 0.50 UJ	< 0.50 U
CHLOROMETHANE	5	< 1.0 U	< 1.0 UJ	< 1.0 U
CIS-1,2-DICHLOROETHENE	5	< 0.50 U	< 0.50 UJ	< 0.50 U
CIS-1,3-DICHLOROPROPENE	0.4	< <b>0.50 U</b>	< <b>0.50 UJ</b>	< <b>0.50 U</b>
CYCLOHEXANE	NL	< 0.50 U	< 0.50 UJ	< 0.50 U
DIBROMOCHLOROMETHANE	5	< 0.50 U	< 0.50 UJ	< 0.50 U
DICHLORODIFLUOROMETHANE	5	< 1.0 U	< 1.0 UJ	< 1.0 UJ
ETHYLBENZENE	5	< 0.50 U	< 0.50 UJ	< 0.50 U
ISOPROPYLBENZENE	5	< 0.50 U	< 0.50 UJ	< 0.50 U
M- AND P-XYLENE	NL	< 1.0 U	< 1.0 UJ	< 1.0 U
METHYL ACETATE	NL	< 0.75 U	< 0.75 UJ	< 0.75 U
METHYL CYCLOHEXANE	NL	< 0.50 U	< 0.50 UJ	< 0.50 U
METHYL TERT-BUTYL ETHER	10	< 0.50 U	< 0.50 UJ	< 0.50 U
METHYLENE CHLORIDE	5	< 2.5 U	< 2.5 UJ	< 2.5 U
O-XYLENE	NL	< 0.50 U	< 0.50 UJ	< 0.50 U
STYRENE	5	< 0.50 U	< 0.50 UJ	< 0.50 U
TETRACHLOROETHENE	5	< 0.50 U	< 0.50 UJ	< 0.50 U
TOLUENE	5	< 0.50 U	<b>3.7 J</b>	<b>0.72 J</b>
TRANS-1,2-DICHLOROETHENE	5	< 0.50 U	< 0.50 UJ	< 0.50 U
TRANS-1,3-DICHLOROPROPENE	0.4	< <b>0.50 U</b>	< <b>0.50 UJ</b>	< <b>0.50 U</b>
TRICHLOROETHENE	5	< 0.50 U	< 0.50 UJ	< 0.50 U
TRICHLOROFLUOROMETHANE	5	< 1.0 U	< 1.0 UJ	< 1.0 U
VINYL CHLORIDE	2	< 1.0 U	< 1.0 UJ	< 1.0 U
XYLENES, TOTAL	5	< 1.5 U	< 1.5 UJ	< 1.5 U

Location	VPB152	VPB152	VPB152	VPB152	
Sample Date	NYSDEC Groundwater Guidance or Standard Value (Note 1)	6/16/2014	6/16/2014	6/16/2014	6/17/2014
Sample ID		VPB152-GW-061614- 660-662	VPB152-GW-061614- 680-682	VPB152-GW-061614- 700-702	VPB152-GW-061714- 725-727
Sample Interval		660 - 662 ft	680 - 682 ft	700 - 702 ft	725 - 727 ft
Sample type code		N	N	N	N
<b>VOC 8260C (ug/L)</b>					
1,1,1-TRICHLOROETHANE	5	< 0.50 U	< 0.50 U	< 50 UJ	< 25 UJ
1,1,2,2-TETRACHLOROETHANE	5	< 0.50 U	< 0.50 U	< 50 UJ	< 25 UJ
1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	5	< 0.50 U	< 0.50 U	< 50 UJ	< 25 UJ
1,1,2-TRICHLOROETHANE	1	< 0.50 U	< 0.50 U	< 50 UJ	< 25 UJ
1,1-DICHLOROETHANE	5	< 0.50 U	< 0.50 U	< 50 UJ	< 25 UJ
1,1-DICHLOROETHENE	5	< 0.50 U	< 0.50 U	< 50 UJ	< 25 UJ
1,2,4-TRICHLOROBENZENE	5	< 0.50 U	< 0.50 U	< 50 UJ	< 25 UJ
1,2-DIBROMO-3-CHLOROPROPANE	0.04	< 0.75 U	< 0.75 U	< 75 UJ	< 38 UJ
1,2-DIBROMOETHANE	NL	< 0.50 U	< 0.50 U	< 50 UJ	< 25 UJ
1,2-DICHLOROBENZENE	3	< 0.50 U	< 0.50 U	< 50 UJ	< 25 UJ
1,2-DICHLOROETHANE	5	< 0.50 U	< 0.50 U	< 50 UJ	< 25 UJ
1,2-DICHLOROETHENE, TOTAL	5	< 1.0 U	< 1.0 U	< 100 UJ	< 50 UJ
1,2-DICHLOROPROPANE	1	< 0.50 U	< 0.50 U	< 50 UJ	< 25 UJ
1,3-DICHLOROBENZENE	3	< 0.50 U	< 0.50 U	< 50 UJ	< 25 UJ
1,4-DICHLOROBENZENE	3	< 0.50 U	< 0.50 U	< 50 UJ	< 25 UJ
2-BUTANONE	50	< 2.5 UJ	4.8 J	< 250 UJ	< 120 UJ
2-HEXANONE	50	< 2.5 U	< 2.5 U	< 250 UJ	< 120 UJ
4-METHYL-2-PENTANONE	NL	< 2.5 UJ	< 2.5 UJ	< 250 UJ	< 120 UJ
ACETONE	50	8.4 J	22 J	< 250 UJ	< 120 UJ
BENZENE	1	< 0.50 U	< 0.50 U	< 50 UJ	< 25 UJ
BROMODICHLOROMETHANE	50	< 0.50 U	< 0.50 U	< 50 UJ	< 25 UJ
BROMOFORM	50	< 0.50 U	< 0.50 U	< 50 UJ	< 25 UJ
BROMOMETHANE	5	< 1.0 U	< 1.0 U	< 100 UJ	< 50 UJ
CARBON DISULFIDE	60	< 0.50 U	< 1.0 U	< 50 UJ	< 25 UJ
CARBON TETRACHLORIDE	5	< 0.50 U	< 0.50 U	< 50 UJ	< 25 UJ
CHLOROBENZENE	5	< 0.50 U	< 0.50 U	< 50 UJ	< 25 UJ
CHLOROETHANE	5	< 1.0 UJ	< 1.0 UJ	< 100 UJ	< 50 UJ
CHLOROFORM	7	< 0.50 U	< 0.50 U	< 50 UJ	< 25 UJ
CHLOROMETHANE	5	< 1.0 U	< 1.0 U	< 100 UJ	< 50 UJ
CIS-1,2-DICHLOROETHENE	5	< 0.50 U	< 0.50 U	< 50 UJ	< 25 UJ
CIS-1,3-DICHLOROPROPENE	0.4	< 0.50 U	< 0.50 U	< 50 UJ	< 25 UJ
CYCLOHEXANE	NL	< 0.50 U	< 0.50 U	< 50 UJ	< 25 UJ
DIBROMOCHLOROMETHANE	5	< 0.50 U	< 0.50 U	< 50 UJ	< 25 UJ
DICHLORODIFLUOROMETHANE	5	< 1.0 UJ	< 1.0 UJ	< 100 UJ	< 50 UJ
ETHYLBENZENE	5	< 0.50 U	< 0.50 U	< 50 UJ	< 25 UJ
ISOPROPYLBENZENE	5	< 0.50 U	< 0.50 U	< 50 UJ	< 25 UJ
M- AND P-XYLENE	NL	< 1.0 U	< 1.0 U	< 100 UJ	< 50 UJ
METHYL ACETATE	NL	< 0.75 U	< 0.75 U	< 75 UJ	< 38 UJ
METHYL CYCLOHEXANE	NL	< 0.50 U	< 0.50 U	< 50 UJ	< 25 UJ
METHYL TERT-BUTYL ETHER	10	< 0.50 U	< 0.50 U	< 50 UJ	< 25 UJ
METHYLENE CHLORIDE	5	< 2.5 U	< 2.5 U	< 500 UJ	67 J
O-XYLENE	NL	< 0.50 U	< 0.50 U	< 50 UJ	< 25 UJ
STYRENE	5	< 0.50 U	< 0.50 U	< 50 UJ	< 25 UJ
TETRACHLOROETHENE	5	< 0.50 U	< 0.50 U	< 50 UJ	< 25 UJ
TOLUENE	5	0.52 J	0.69 J	< 50 UJ	< 25 UJ
TRANS-1,2-DICHLOROETHENE	5	< 0.50 U	< 0.50 U	< 50 UJ	< 25 UJ
TRANS-1,3-DICHLOROPROPENE	0.4	< 0.50 U	< 0.50 U	< 50 UJ	< 25 UJ
TRICHLOROETHENE	5	< 0.50 U	< 0.50 U	< 50 UJ	< 25 UJ
TRICHLOROFLUOROMETHANE	5	< 1.0 U	< 1.0 U	< 100 UJ	< 50 UJ
VINYL CHLORIDE	2	< 1.0 U	< 1.0 U	< 100 UJ	< 50 UJ
XYLENES, TOTAL	5	< 1.5 U	< 1.5 U	< 150 UJ	< 75 UJ

Location	VPB152	VPB152	VPB152	VPB152	
Sample Date	NYSDEC Groundwater Guidance or Standard Value (Note 1)	6/17/2014	6/19/2014	6/19/2014	6/20/2014
Sample ID		VPB152-GW-061714- 740-742	VPB152-GW-061914- 780-782	VPB152-GW-061914- 800-802	VPB152-GW-062014- 830-832
Sample Interval		740 - 742 ft	780 - 782 ft	800 - 802 ft	830 - 832 ft
Sample type code		N	N	N	N
<b>VOC 8260C (ug/L)</b>					
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	< <b>0.75 U</b>	< <b>0.75 U</b>	< <b>0.75 U</b>	< <b>0.75 U</b>
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
2-BUTANONE	50	<b>2.1 J</b>	< 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 UJ	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ
ACETONE	50	<b>11 J</b>	<b>7.5 J</b>	<b>11 J</b>	<b>10 J</b>
BENZENE	1	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
BROMODICHLOROMETHANE	50	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
BROMOFORM	50	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
BROMOMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
CARBON DISULFIDE	60	< 0.50 U	< 0.50 U	< 1.0 U	< 0.50 U
CARBON TETRACHLORIDE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 UJ
CHLOROFORM	7	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
CIS-1,2-DICHLOROETHENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CIS-1,3-DICHLOROPROPENE	0.4	< <b>0.50 U</b>	< <b>0.50 U</b>	< <b>0.50 U</b>	< <b>0.50 U</b>
CYCLOHEXANE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 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 U	< 0.50 U	< 0.50 U	< 0.50 U
TOLUENE	5	<b>5.6</b>	< 0.50 U	< 0.50 U	<b>3.5</b>
TRANS-1,2-DICHLOROETHENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TRANS-1,3-DICHLOROPROPENE	0.4	< <b>0.50 U</b>	< <b>0.50 U</b>	< <b>0.50 U</b>	< <b>0.50 U</b>
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 UJ
VINYL CHLORIDE	2	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 UJ
XYLENES, TOTAL	5	< 1.5 U	< 1.5 U	< 1.5 U	< 1.5 U

Location	VPB152	VPB152	VPB152	VPB152
Sample Date	6/23/2014	6/24/2014	6/24/2014	6/25/2014
Sample ID	VPB152-GW-062314-850-852	VPB152-GW-062414-870-882	VPB152-GW-062414-880-882	VPB152-GW-062514-910-912
Sample Interval	850 - 852 ft	870 - 872 ft	880 - 882 ft	910 - 912 ft
Sample type code	N	N	N	N
VOC 8260C (ug/L)				
1,1,1-TRICHLOROETHANE	5	< 0.50 U	< 0.50 UJ	< 0.50 UJ
1,1,2,2-TETRACHLOROETHANE	5	< 0.50 U	< 0.50 UJ	< 0.50 UJ
1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	5	< 0.50 U	< 0.50 UJ	< 0.50 UJ
1,1,2-TRICHLOROETHANE	1	< 0.50 U	< 0.50 UJ	< 0.50 UJ
1,1-DICHLOROETHANE	5	< 0.50 U	< 0.50 UJ	< 0.50 UJ
1,1-DICHLOROETHENE	5	< 0.50 U	< 0.50 UJ	< 0.50 UJ
1,2,4-TRICHLOROBENZENE	5	< 0.50 U	< 0.50 UJ	< 0.50 UJ
1,2-DIBROMO-3-CHLOROPROPANE	0.04	< 0.75 U	< 0.75 UJ	< 0.75 UJ
1,2-DIBROMOETHANE	NL	< 0.50 U	< 0.50 UJ	< 0.50 UJ
1,2-DICHLOROBENZENE	3	< 0.50 U	< 0.50 UJ	< 0.50 UJ
1,2-DICHLOROETHANE	5	< 0.50 U	< 0.50 UJ	< 0.50 UJ
1,2-DICHLOROETHENE, TOTAL	5	< 1.0 U	< 1.0 UJ	< 1.0 UJ
1,2-DICHLOROPROPANE	1	< 0.50 U	< 0.50 UJ	< 0.50 UJ
1,3-DICHLOROBENZENE	3	< 0.50 U	< 0.50 UJ	< 0.50 UJ
1,4-DICHLOROBENZENE	3	< 0.50 U	< 0.50 UJ	< 0.50 UJ
2-BUTANONE	50	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ
2-HEXANONE	50	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ
4-METHYL-2-PENTANONE	NL	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ
ACETONE	50	7.7 J	< 5.0 UJ	< 13 UJ
BENZENE	1	< 0.50 U	< 0.50 UJ	< 0.50 UJ
BROMODICHLOROMETHANE	50	< 0.50 U	< 0.50 UJ	< 0.50 UJ
BROMOFORM	50	< 0.50 U	< 0.50 UJ	< 0.50 UJ
BROMOMETHANE	5	< 1.0 U	< 1.0 UJ	< 1.0 UJ
CARBON DISULFIDE	60	< 0.50 U	< 0.50 UJ	< 0.50 UJ
CARBON TETRACHLORIDE	5	< 0.50 U	< 0.50 UJ	< 0.50 UJ
CHLOROBENZENE	5	< 0.50 U	< 0.50 UJ	< 0.50 UJ
CHLOROETHANE	5	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ
CHLOROFORM	7	< 0.50 U	< 0.50 UJ	< 0.50 UJ
CHLOROMETHANE	5	< 1.0 U	< 1.0 UJ	< 1.0 UJ
CIS-1,2-DICHLOROETHENE	5	< 0.50 U	< 0.50 UJ	< 0.50 UJ
CIS-1,3-DICHLOROPROPENE	0.4	< 0.50 U	< 0.50 UJ	< 0.50 UJ
CYCLOHEXANE	NL	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
DIBROMOCHLOROMETHANE	5	< 0.50 U	< 0.50 UJ	< 0.50 UJ
DICHLORODIFLUOROMETHANE	5	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ
ETHYLBENZENE	5	< 0.50 U	< 0.50 UJ	< 0.50 UJ
ISOPROPYLBENZENE	5	< 0.50 U	< 0.50 UJ	< 0.50 UJ
M- AND P-XYLENE	NL	< 1.0 U	< 1.0 UJ	< 1.0 UJ
METHYL ACETATE	NL	< 0.75 U	< 0.75 UJ	< 0.75 UJ
METHYL CYCLOHEXANE	NL	< 0.50 U	< 0.50 UJ	< 0.50 UJ
METHYL TERT-BUTYL ETHER	10	< 0.50 U	< 0.50 UJ	< 0.50 UJ
METHYLENE CHLORIDE	5	< 2.5 U	< 2.5 UJ	< 2.5 UJ
O-XYLENE	NL	< 0.50 U	< 0.50 UJ	< 0.50 UJ
STYRENE	5	< 0.50 U	< 0.50 UJ	< 0.50 UJ
TETRACHLOROETHENE	5	< 0.50 U	< 0.50 UJ	< 0.50 UJ
TOLUENE	5	< 0.50 U	< 0.50 UJ	1.4 J
TRANS-1,2-DICHLOROETHENE	5	< 0.50 U	< 0.50 UJ	< 0.50 UJ
TRANS-1,3-DICHLOROPROPENE	0.4	< 0.50 U	< 0.50 UJ	< 0.50 UJ
TRICHLOROETHENE	5	< 0.50 U	< 0.50 UJ	< 0.50 UJ
TRICHLOROFLUOROMETHANE	5	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ
VINYL CHLORIDE	2	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ
XYLENES, TOTAL	5	< 1.5 U	< 1.5 UJ	< 1.5 UJ



Location		VPB152	VPB152
Sample Date	NYSDEC	6/26/2014	6/27/2014
Sample ID	Groundwater Guidance or Standard Value (Note 1)	VPB152-GW-062614- 930-932	VPB152-GW-062714- 950-952
Sample Interval		930 - 932 ft	950 - 952 ft
Sample type code		N	N
<b>VOC 8260C (ug/L)</b>			
1,1,1-TRICHLOROETHANE	5	< 1.0 UJ	< 1.0 UJ
1,1,2,2-TETRACHLOROETHANE	5	< 1.0 UJ	< 1.0 UJ
1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	5	< 1.0 UJ	< 1.0 UJ
1,1,2-TRICHLOROETHANE	1	< 1.0 UJ	< 1.0 UJ
1,1-DICHLOROETHANE	5	< 1.0 UJ	< 1.0 UJ
1,1-DICHLOROETHENE	5	< 1.0 UJ	< 1.0 UJ
1,2,4-TRICHLOROBENZENE	5	< 1.0 UJ	< 1.0 UJ
1,2-DIBROMO-3-CHLOROPROPANE	0.04	< 1.5 UJ	< 1.5 UJ
1,2-DIBROMOETHANE	NL	< 1.0 UJ	< 1.0 UJ
1,2-DICHLOROBENZENE	3	< 1.0 UJ	< 1.0 UJ
1,2-DICHLOROETHANE	5	< 1.0 UJ	< 1.0 UJ
1,2-DICHLOROETHENE, TOTAL	5	< 2.0 UJ	< 2.0 UJ
1,2-DICHLOROPROPANE	1	< 1.0 UJ	< 1.0 UJ
1,3-DICHLOROBENZENE	3	< 1.0 UJ	< 1.0 UJ
1,4-DICHLOROBENZENE	3	< 1.0 UJ	< 1.0 UJ
2-BUTANONE	50	< 5.0 UJ	< 5.0 UJ
2-HEXANONE	50	< 5.0 UJ	< 5.0 UJ
4-METHYL-2-PENTANONE	NL	< 5.0 UJ	< 5.0 UJ
ACETONE	50	< 32 UJ	46 J
BENZENE	1	< 1.0 UJ	< 1.0 UJ
BROMODICHLOROMETHANE	50	< 1.0 UJ	< 1.0 UJ
BROMOFORM	50	< 1.0 UJ	< 1.0 UJ
BROMOMETHANE	5	< 2.0 UJ	< 2.0 UJ
CARBON DISULFIDE	60	< 1.0 UJ	< 1.0 UJ
CARBON TETRACHLORIDE	5	< 1.0 UJ	< 1.0 UJ
CHLOROBENZENE	5	< 1.0 UJ	< 1.0 UJ
CHLOROETHANE	5	< 2.0 UJ	< 2.0 UJ
CHLOROFORM	7	< 1.0 UJ	< 1.0 UJ
CHLOROMETHANE	5	< 2.0 UJ	< 2.0 UJ
CIS-1,2-DICHLOROETHENE	5	< 1.0 UJ	< 1.0 UJ
CIS-1,3-DICHLOROPROPENE	0.4	< 1.0 UJ	< 1.0 UJ
CYCLOHEXANE	NL	< 1.0 UJ	< 1.0 UJ
DIBROMOCHLOROMETHANE	5	< 1.0 UJ	< 1.0 UJ
DICHLORODIFLUOROMETHANE	5	< 2.0 UJ	< 2.0 UJ
ETHYLBENZENE	5	< 1.0 UJ	< 1.0 UJ
ISOPROPYLBENZENE	5	< 1.0 UJ	< 1.0 UJ
M- AND P-XYLENE	NL	< 2.0 UJ	< 2.0 UJ
METHYL ACETATE	NL	< 1.5 UJ	< 1.5 UJ
METHYL CYCLOHEXANE	NL	< 1.0 UJ	< 1.0 UJ
METHYL TERT-BUTYL ETHER	10	< 1.0 UJ	< 1.0 UJ
METHYLENE CHLORIDE	5	3.2 J	6.6 J
O-XYLENE	NL	< 1.0 UJ	< 1.0 UJ
STYRENE	5	< 1.0 UJ	< 1.0 UJ
TETRACHLOROETHENE	5	< 1.0 UJ	< 1.0 UJ
TOLUENE	5	3.8 J	5.4 J
TRANS-1,2-DICHLOROETHENE	5	< 1.0 UJ	< 1.0 UJ
TRANS-1,3-DICHLOROPROPENE	0.4	< 1.0 UJ	< 1.0 UJ
TRICHLOROETHENE	5	< 1.0 UJ	< 1.0 UJ
TRICHLOROFLUOROMETHANE	5	< 2.0 UJ	< 2.0 UJ
VINYL CHLORIDE	2	< 2.0 UJ	< 2.0 UJ
XYLENES, TOTAL	5	< 3.0 UJ	< 3.0 UJ

**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*** = Detection limit 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.

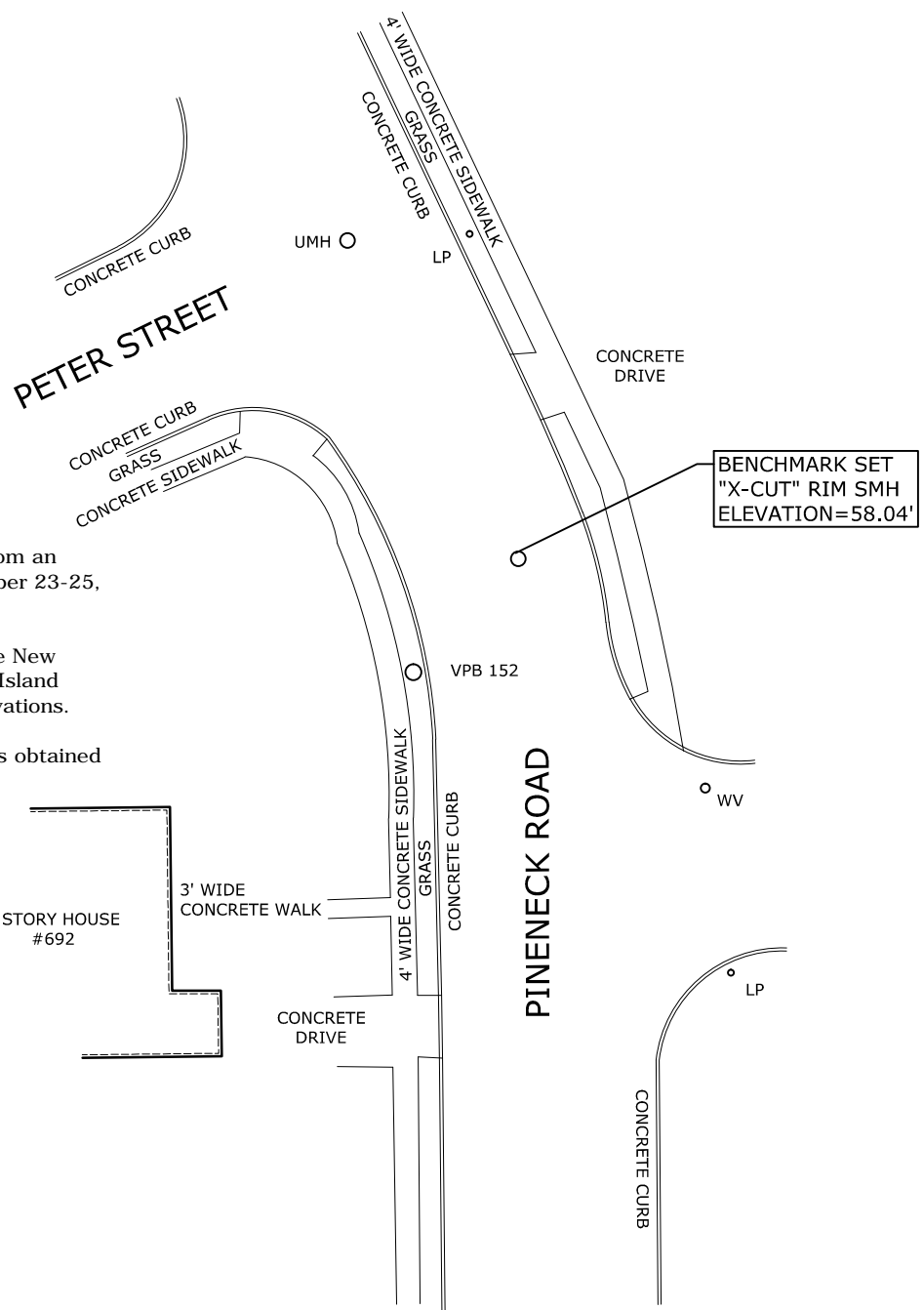
J = The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.

## **Section 6**

### **Survey**

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

Description	Northing	Easting	Latitude	Longitude	Ground	Rim	PVC
VPB 152	197311.98	1126289.12	N40-42-25.60	W73-29-15.67	58.16	NA	NA

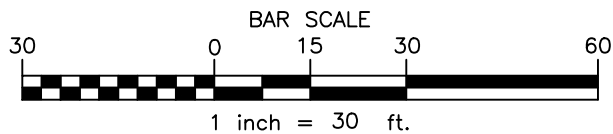
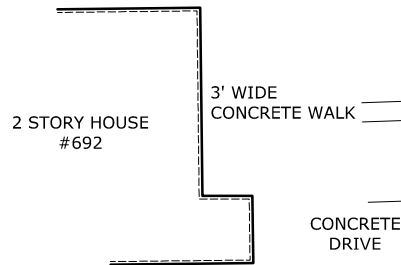


**Map Notes**

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

**Legend**

- LP Light Post
- UMH Unknown Manhole
- VPB Vertical Profile Boring
- WV Water Valve



DWG NO. 14-502

Date	RECORD OF WORK	Appr.	VERTICAL PROFILE BORING 152 SURVEY LOCATION 692 PINENECK ROAD	
			TOWN OF SEAFORD	NASSAU COUNTY, NEW YORK
			<b>C.T. MALE ASSOCIATES</b> 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	
Drafter: LMK		Checker: JFC		
Appr. by: JFC		Proj. No. 14.4121		

