

# SEPTEMBER 2015 GROUNDWATER SAMPLING DATA SUMMARY REPORT

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

Prepared for:



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

February 2016

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Norfolk, Virginia 23511

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Contract Number: N62470-11-D-8013  
CTO WE15

February 2016

A handwritten signature in cursive script that reads "Brian Caldwell".

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**Brian Caldwell**  
Contract Task Order Manager

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- Appendix C Analytical Data Validation - ARCADIS
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## List of Acronyms and Abbreviations

DOT	Department of Transportation
IDW	Investigation Derived Waste
Katahdin	Katahdin Analytical Services
NWIRP	Naval Weapons Industrial Reserve Plant
ONCT	Onsite Containment System
OU	Operable Unit
POTW	Publicly Owned Treatment Works
QA	Quality Assurance
QC	Quality Control
SAP	Sampling and Analysis Plan
UFP	Uniform Federal Policy
VOC	Volatile Organic Compounds

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## 1.0 PROJECT BACKGROUND

Resolution Consultants has prepared this Groundwater Sampling Data Summary Report for the Naval Facilities Engineering Command, Mid-Atlantic under contract task order WE15 Contract N62470-11-D-8013. The report describes quarterly sampling activities in September 2015, which is part of the Navy's ongoing Environmental Restoration Program 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).

This data summary report provides information on quarterly sampling of 34 monitoring wells by Resolution Consultants (for the Navy) and ARCADIS (for Northrop Grumman). The purpose of this sampling is to provide information on the extent and magnitude of volatile organic compounds (VOCs) located in a narrow area immediately south of the Onsite Containment System (ONCT) in the western offsite plume, which could represent contamination that has bypassed the ONCT. The locations of monitoring wells sampled as part of this effort are shown in Figure 2 and listed in Table 1.

Per an agreement between the Navy and Northrop Grumman (letter May 6, 2015), Northrop Grumman takes over the quarterly sampling of selected monitoring wells after the initial sampling (post well installation) by Resolution Consultants. A total of 12 monitoring wells were transitioned to Northrop Grumman for the September 2015 quarterly sampling.

Documentation of these activities is included in the appendices of this report. Groundwater sampling forms, and analytical data validation for wells sampled by Resolution Consultants are included in Appendix A and B, respectively. Appendix C contains analytical data validation for wells sampled by ARCADIS.

Additional Navy-owned wells were sampled by ARCADIS in the third quarter of 2015 as part of ongoing OU2 monitoring programs. These activities are documented in their report, *Results of Third Quarter 2015 Groundwater Monitoring* (ARCADIS, 2015), included in Appendix D to this report.

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## 2.0 FIELD PROGRAM

Field tasks were conducted in September of 2015 in accordance with the Uniform Federal Policy (UFP) Sampling and Analysis Plan (SAP) Addendum: *Groundwater Sampling Using Low Stress (Low Flow) Purging and Sampling Protocol* (Resolution Consultants, 2013). The field investigation included purging and sampling of monitoring wells in the quarterly groundwater sampling network.

The September 2015 quarterly sampling round consisted of a total of 34 wells (Table 1). Of these, 22 groundwater wells were sampled by Resolution Consultants and 12 were sampled by ARCADIS, Northrop Grumman's consultant. ARCADIS sampled the following wells after initial sampling by Resolution Consultants in June 2015 (Resolution Consultants, 2015): BPOW5-5, BPOW5-6, BPOW6-5, and BPOW6-6. These wells were transitioned to ARCADIS for the September sampling event.

### 2.1 Sampling

Resolution Consultants purged monitoring wells using a bladder pump with the intake placed at the approximate midpoint of the screened interval. The following field water quality parameters were continuously measured during purging: water temperature, pH, conductivity, oxidation-reduction potential, dissolved oxygen and turbidity. Groundwater analytical samples were collected when field water quality parameters stabilized. Samples were analyzed for VOCs via Method 8260C and 1,4-dioxane via Method 8270C by Katahdin Analytical Services (Katahdin). All purge water was managed as investigation derived waste (IDW). Quality assurance (QA) and quality control (QC) samples were collected during the sampling effort.

Analytical results and stabilized field parameters for wells sampled by Resolution Consultants are summarized in Table 2 and Table 3, respectively. Groundwater sample forms and data validation packages for wells sampled by Resolution Consultants are included in Appendix A and B, respectively.

Results and data validation packages for ARCADIS-sampled wells are provided in Table 4, Table 5 and Appendix C of this report.

### 2.2 Investigation Derived Waste

Resolution Consultants utilized dedicated and disposable sampling equipment when possible to avoid the potential for cross-contamination of samples. The sampling equipment included dedicated disposable polyethylene tubing, disposable gloves, and laboratory supplied sample bottles. Hand

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held equipment was decontaminated using a liquinox and water wash, a potable water rinse followed by a distilled water rinse. Purge water was collected in 5-gallon pails or 55-gallon drums.

Resolution Consultants transported purge water from point of generation to the designated staging area at NWIRP in Department of Transportation (DOT) approved 5-gallon pails. Purge water was then containerized in a frac tank and stored at NWIRP Bethpage for characterization and ultimate disposal to the Nassau County Publicly Owned Treatment Works (POTW) in accordance with the facility's existing discharge permit. A representative water sample was collected from each of the frac tanks and submitted to Katahdin for analysis of VOCs via Method SW 624, pH via Method SW 9040B, PCBs via Method 8082 and Total Metals via Method SW 846. All analytical criteria were met for disposal of water. No solid waste was generated during sampling.



### **3.0 SUMMARY**

Well construction information for all wells sampled by Resolution Consultants and ARCADIS is summarized in Table 1.

Analytical results and stabilized field water quality parameters for wells sampled by Resolution Consultants are summarized in Tables 2 and 3, respectively. Groundwater sample forms and data validation packages for wells sampled by Resolution Consultants are included in Appendix A and B, respectively.

Analytical results for wells sampled by ARCADIS are summarized in Tables 4 and 5. Data validation packages for wells sampled by ARCADIS are included in Appendix C.

Documentation of other ongoing OU2 groundwater sampling activities by ARCADIS in the third quarter 2015 can be found in their report, *Results of Third Quarter 2015 Groundwater Monitoring* (ARCADIS, 2015). This ARCADIS report is included as Appendix D to this report.

#### 4.0 REFERENCES

ARCADIS, 2015. *Results of Third Quarter 2015 Groundwater Monitoring, Operable Unit 2, Northrop Grumman Systems Corporation and Naval Weapons Industrial Reserve Plant (NWIRP) Sites, Bethpage, New York*. November.

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

Resolution Consultants, 2015. *June 2015 Groundwater Sampling Data Summary Report, Bethpage, NY*. October.

**Tables**

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

Well	Total Depth (ft bgs)	Top of Screen (ft bgs)	Bottom of Screen (ft bgs)	Mid-screen (ft bgs)	Sump Length (ft)	VPB affiliation	Sampled by
RE103D1	645	625	640	630	5	VPB137	Resolution
RE103D2	673	653	673	663	0	VPB137	Resolution
RE103D3	735	715	730	720	5	VPB137	Resolution
RE104D1	375	350	370	360	5	VPB138	Resolution
RE104D2	735	710	730	720	5	VPB138	Resolution
RE104D3	785	760	780	770	5	VPB138	Resolution
RE105D1	555	530	550	540	5	VPB139	Resolution
RE105D2	755	730	750	740	5	VPB139	Resolution
RE108D1	555	530	550	540	5	VPB142	Resolution
RE108D2	655	630	650	640	5	VPB142	Resolution
RE120D1	655	630	650	640	5	VPB154	Resolution
RE120D2	713	690	710	700	3	VPB154	Resolution
RE120D3	765	740	760	750	5	VPB154	Resolution
RE122D1	545	520	540	530	5	VPB156	Resolution
RE122D2	615	590	610	600	5	VPB156	Resolution
RE122D3	740	715	735	725	5	VPB156	Resolution
RE123D1	505	480	500	490	5	VPB157	Resolution
RE123D2	660	635	655	645	5	VPB157	Resolution
RE123D3	840	815	835	825	5	VPB157	Resolution
TT101D	350	325	345	335	5	VPB129	Resolution
TT101D1	595	570	590	580	5	VPB129	Resolution
TT101D2	765	740	760	750	5	VPB129	Resolution
BPOW5-1	515	480	510	495	5	VPB132	ARCADIS
BPOW5-2	585	540	580	560	5	VPB132	ARCADIS
BPOW5-3	665	620	660	640	5	VPB132	ARCADIS
BPOW5-4	575	545	570	557.5	5	VPB151	ARCADIS
BPOW5-5	545	515	540	527.5	5	VPB152	ARCADIS
BPOW5-6	615	585	610	597.5	5	VPB152	ARCADIS
BPOW6-1	580	550	575	562.5	5	VPB145	ARCADIS
BPOW6-2	785	755	780	767.5	5	VPB145	ARCADIS
BPOW6-3	780	750	775	762.5	5	VPB146	ARCADIS
BPOW6-4	575	545	570	557.5	5	VPB146	ARCADIS
BPOW6-5	555	525	550	537.5	5	VPB147	ARCADIS
BPOW6-6	800	770	795	782.5	5	VPB147	ARCADIS

ft bgs - feet below ground surface

**TABLE 2. ANALYTICAL DATA SUMMARY FOR  
WELLS SAMPLED BY RESOLUTION CONSULTANTS  
2015 OU2 GROUNDWATER INVESTIGATION**

Location	NYSDEC Groundwater Guidance or Standard Value (Note 1)	RE104D1	RE104D2	RE104D3	RE105D1
Sample Date		9/25/2015	9/25/2015	9/25/2015	9/28/2015
Sample ID		RE104D1-GW-092515	RE104D2-GW-092515	RE104D3-GW-092515	RE105D1-GW-092815
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	<b>4.6</b>	< 0.50 U	< 0.50 U	<b>11</b>
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	<b>0.87 J</b>	< 0.50 U	< 0.50 U	<b>1.6</b>
1,2,4-TRICHLOROENZENE	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	<b>1.2 J</b>	<b>2.0</b>	< 1.0 U	<b>1.6 J</b>
1,2-DICHLOROPROPANE	1	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,3-DICHLOROBENZENE	3	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,4-DICHLOROBENZENE	3	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,4-DIOXANE (Method 8270D_SIM)	NL	<b>12</b>	<b>0.13 J</b>	< 0.17 U	<b>14</b>
2-BUTANONE	50	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ
2-HEXANONE	50	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ
4-METHYL-2-PENTANONE	NL	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ
ACETONE	50	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ
BENZENE	1	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
BROMODICHLOROMETHANE	50	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
BROMOFORM	50	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
BROMOMETHANE	5	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ
CARBON DISULFIDE	60	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
CARBON TETRACHLORIDE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROETHANE	5	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ
CHLOROFORM	7	< 0.50 U	<b>0.53 J</b>	< 0.50 U	<b>0.34 J</b>
CHLOROMETHANE	5	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ
CIS-1,2-DICHLOROETHENE	5	<b>1.2</b>	<b>2.0</b>	< 0.50 U	<b>1.6</b>
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 U
DIBROMOCHLOROMETHANE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
DICHLORODIFLUOROMETHANE	5	<b>0.49 J</b>	< 1.0 UJ	< 1.0 UJ	<b>0.74 J</b>
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 UJ	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
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	<b>2.3 J</b>	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
TOLUENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TRANS-1,2-DICHLOROETHENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TRANS-1,3-DICHLOROPROPENE	0.4	< <b>0.50 U</b>	< <b>0.50 U</b>	< <b>0.50 U</b>	< <b>0.50 U</b>
TRICHLOROETHENE	5	<b>110</b>	<b>4.2</b>	< 0.50 U	<b>94</b>
TRICHLOROFLUOROMETHANE	5	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ
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

**TABLE 2. ANALYTICAL DATA SUMMARY FOR  
WELLS SAMPLED BY RESOLUTION CONSULTANTS  
2015 OJ2 GROUNDWATER INVESTIGATION**

Location	NYSDEC Groundwater Guidance or Standard Value (Note 1)	RE105D2	RE108D1	RE108D2	RE123D1
Sample Date		9/28/2015	9/28/2015	9/28/2015	9/29/2015
Sample ID		RE105D2-GW-092815	RE108D1-GW-092815	RE108D2-GW-092815	RE123D1-GW-092915
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	<b>30</b>	<b>1.2</b>	<b>8.3</b>	< 0.50 U
1,1,2-TRICHLOROETHANE	1	<b>1.1</b>	< 0.50 U	<b>1.5</b>	< 0.50 U
1,1-DICHLOROETHANE	5	<b>1.7</b>	< 0.50 U	<b>5.3</b>	< 0.50 U
1,1-DICHLOROETHENE	5	<b>6.8</b>	< 0.50 U	<b>7.6</b>	<b>0.42 J</b>
1,2,4-TRICHLOROENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DIBROMO-3-CHLOROPROPANE	0.04	<b>&lt; 0.75 U</b>	<b>&lt; 0.75 U</b>	<b>&lt; 0.75 U</b>	<b>&lt; 0.75 U</b>
1,2-DIBROMOETHANE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DICHLOROENZENE	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	<b>3.6</b>	<b>0.39 J</b>	<b>8.6</b>	<b>0.50 J</b>
1,2-DICHLOROPROPANE	1	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,3-DICHLOROENZENE	3	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,4-DICHLOROENZENE	3	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,4-DIOXANE (Method 8270D_SIM)	NL	<b>8.7</b>	<b>7.9</b>	<b>7.5</b>	<b>6.6</b>
2-BUTANONE	50	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ
2-HEXANONE	50	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ
4-METHYL-2-PENTANONE	NL	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ
ACETONE	50	< 2.5 UJ	<b>6.9 J</b>	< 2.5 UJ	<b>5.4 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 UJ	< 1.0 UJ	< 1.0 UJ	< 1.0 U
CARBON DISULFIDE	60	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
CARBON TETRACHLORIDE	5	<b>3.7</b>	< 0.50 U	<b>1.6</b>	< 0.50 U
CHLOROENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROETHANE	5	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ	< 1.0 U
CHLOROFORM	7	<b>2.2</b>	< 0.50 U	<b>3.7</b>	< 0.50 U
CHLOROMETHANE	5	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ	< 1.0 U
CIS-1,2-DICHLOROETHENE	5	<b>3.6</b>	<b>0.39 J</b>	<b>8.6</b>	<b>0.50 J</b>
CIS-1,3-DICHLOROPROPENE	0.4	<b>&lt; 0.50 U</b>	<b>&lt; 0.50 U</b>	<b>&lt; 0.50 U</b>	<b>&lt; 0.50 U</b>
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	<b>0.40 J</b>	< 1.0 UJ	< 1.0 UJ	< 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 UJ	< 0.50 UJ	< 0.50 UJ	< 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 UJ
STYRENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TETRACHLOROETHENE	5	<b>1.4 J</b>	<b>1.3 J</b>	<b>2.5 J</b>	<b>3.8 J</b>
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	<b>&lt; 0.50 U</b>	<b>&lt; 0.50 U</b>	<b>&lt; 0.50 U</b>	<b>&lt; 0.50 U</b>
TRICHLOROETHENE	5	<b>1900</b>	<b>98</b>	<b>3400</b>	<b>12</b>
TRICHLOROFUOROMETHANE	5	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ	< 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

**TABLE 2. ANALYTICAL DATA SUMMARY FOR  
WELLS SAMPLED BY RESOLUTION CONSULTANTS  
2015 UO2 GROUNDWATER INVESTIGATION**

Location	NYSDEC Groundwater Guidance or Standard Value (Note 1)	RE123D2	RE123D3	RE120D1	RE120D2
Sample Date		9/29/2015	9/29/2015	9/29/2015	9/29/2015
Sample ID		RE123D2-GW-092915	RE123D3-GW-092915	RE120D1-GW-092915	RE120D2-GW-092915
Sample type code		N	N	N	N
<b>VOC 8260C (ug/L)</b>					
1,1,1-TRICHLOROETHANE	5	< 0.50 U	< 0.50 U	<b>1.4</b>	< 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	<b>36</b>	<b>21</b>
1,1,2-TRICHLOROETHANE	1	< 0.50 U	< 0.50 U	<b>1.6</b>	<b>0.45 J</b>
1,1-DICHLOROETHANE	5	< 0.50 U	< 0.50 U	<b>2.9</b>	< 0.50 U
1,1-DICHLOROETHENE	5	< 0.50 U	< 0.50 U	<b>20</b>	<b>4.9</b>
1,2,4-TRICHLOROENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DIBROMO-3-CHLOROPROPANE	0.04	<b>&lt; 0.75 U</b>	<b>&lt; 0.75 U</b>	<b>&lt; 0.75 UJ</b>	<b>&lt; 0.75 UJ</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	<b>2.7</b>	<b>2.7</b>
1,2-DICHLOROPROPANE	1	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,3-DICHLOROBENZENE	3	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,4-DICHLOROBENZENE	3	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,4-DIOXANE (Method 8270D_SIM)	NL	<b>0.93</b>	< 0.17 U	<b>25</b>	<b>16</b>
2-BUTANONE	50	< 2.5 UJ	< 2.5 UJ	< 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	< 2.5 UJ	< 2.5 UJ	< 2.5 U	< 2.5 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 UJ	< 0.50 UJ	< 0.50 U	< 0.50 U
CARBON TETRACHLORIDE	5	< 0.50 U	< 0.50 U	<b>0.77 J</b>	<b>0.72 J</b>
CHLOROBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
CHLOROFORM	7	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROMETHANE	5	< 1.0 UJ	< 1.0 UJ	< 1.0 U	< 1.0 U
CIS-1,2-DICHLOROETHENE	5	< 0.50 U	< 0.50 U	<b>2.7</b>	<b>2.7</b>
CIS-1,3-DICHLOROPROPENE	0.4	<b>&lt; 0.50 U</b>	<b>&lt; 0.50 U</b>	<b>&lt; 0.50 U</b>	<b>&lt; 0.50 U</b>
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 UJ	<b>0.31 J</b>
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 UJ	< 0.50 UJ	< 0.50 U	< 0.50 U
STYRENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TETRACHLOROETHENE	5	<b>3.7 J</b>	< 0.50 UJ	<b>1.9</b>	<b>3.1</b>
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	<b>&lt; 0.50 U</b>	<b>&lt; 0.50 U</b>	<b>&lt; 0.50 U</b>	<b>&lt; 0.50 U</b>
TRICHLOROETHENE	5	<b>1.4</b>	< 0.50 U	<b>1300 J</b>	<b>760 J</b>
TRICHLOROFLUOROMETHANE	5	< 1.0 U	< 1.0 U	<b>0.44 J</b>	< 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

**TABLE 2. ANALYTICAL DATA SUMMARY FOR  
WELLS SAMPLED BY RESOLUTION CONSULTANTS  
2015 OU2 GROUNDWATER INVESTIGATION**

Location	NYSDEC Groundwater Guidance or Standard Value (Note 1)	RE120D3	TT101D	TT101D	TT101D1
Sample Date		9/29/2015	9/29/2015	9/29/2015	9/29/2015
Sample ID		RE120D3-GW-092915	TT101D-GW-092915	DUPLICATE-GW-092915	TT101D1-GW-092915
Sample type code		N	N	FD	N
<b>VOC 8260C (ug/L)</b>					
1,1,1-TRICHLOROETHANE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 UJ
1,1,2,2-TETRACHLOROETHANE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 UJ
1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	5	<b>2.9</b>	<b>14</b>	<b>15</b>	<b>13 J</b>
1,1,2-TRICHLOROETHANE	1	< 0.50 U	< 0.50 U	< 0.50 U	<b>0.49 J</b>
1,1-DICHLOROETHANE	5	< 0.50 U	<b>0.43 J</b>	<b>0.61 J</b>	< 0.50 UJ
1,1-DICHLOROETHENE	5	< 0.50 U	<b>3.1</b>	<b>3.3</b>	<b>4.0 J</b>
1,2,4-TRICHLOROENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 UJ
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 UJ</b>
1,2-DIBROMOETHANE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 UJ
1,2-DICHLOROBENZENE	3	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 UJ
1,2-DICHLOROETHANE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 UJ
1,2-DICHLOROETHENE, TOTAL	5	< 1.0 U	<b>1.9 J</b>	<b>1.8 J</b>	<b>0.92 J</b>
1,2-DICHLOROPROPANE	1	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 UJ
1,3-DICHLOROBENZENE	3	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 UJ
1,4-DICHLOROBENZENE	3	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 UJ
1,4-DIOXANE (Method 8270D_SIM)	NL	<b>1.2</b>	<b>13</b>	<b>14</b>	<b>11 J</b>
2-BUTANONE	50	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 UJ
2-HEXANONE	50	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 UJ
4-METHYL-2-PENTANONE	NL	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 UJ
ACETONE	50	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 UJ
BENZENE	1	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 UJ
BROMODICHLOROMETHANE	50	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 UJ
BROMOFORM	50	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 UJ
BROMOMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 UJ
CARBON DISULFIDE	60	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 UJ
CARBON TETRACHLORIDE	5	< 0.50 U	< 0.50 U	< 0.50 U	<b>1.6 J</b>
CHLOROBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 UJ
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 UJ
CHLOROMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 UJ
CIS-1,2-DICHLOROETHENE	5	< 0.50 U	<b>1.9</b>	<b>1.8</b>	<b>0.92 J</b>
CIS-1,3-DICHLOROPROPENE	0.4	< <b>0.50 U</b>	< <b>0.50 U</b>	< <b>0.50 U</b>	< <b>0.50 UJ</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 UJ
DICHLORODIFLUOROMETHANE	5	< 1.0 UJ	<b>2.0 J</b>	<b>2.3 J</b>	<b>1.9 J</b>
ETHYLBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 UJ
ISOPROPYLBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 UJ
M- AND P-XYLENE	NL	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 UJ
METHYL ACETATE	NL	< 0.75 U	< 0.75 U	< 0.75 U	< 0.75 UJ
METHYL CYCLOHEXANE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 UJ
METHYL TERT-BUTYL ETHER	10	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 UJ
METHYLENE CHLORIDE	5	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 UJ
O-XYLENE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 UJ
STYRENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 UJ
TETRACHLOROETHENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 UJ
TOLUENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 UJ
TRANS-1,2-DICHLOROETHENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 UJ
TRANS-1,3-DICHLOROPROPENE	0.4	< <b>0.50 U</b>	< <b>0.50 U</b>	< <b>0.50 U</b>	< <b>0.50 UJ</b>
TRICHLOROETHENE	5	<b>120</b>	<b>67</b>	<b>67 J</b>	<b>170 J</b>
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 UJ



**TABLE 2. ANALYTICAL DATA SUMMARY FOR  
WELLS SAMPLED BY RESOLUTION CONSULTANTS  
2015 UO2 GROUNDWATER INVESTIGATION**

Location	NYSDEC Groundwater Guidance or Standard Value (Note 1)	TT101D2	RE122D1	RE122D2	RE122D3
Sample Date		9/29/2015	9/30/2015	9/30/2015	9/30/2015
Sample ID		TT101D2-GW-092915	RE122D1-GW-093015	RE122D2-GW-093015	RE122D3-GW-093015
Sample type code		N	N	N	N
<b>VOC 8260C (ug/L)</b>					
1,1,1-TRICHLOROETHANE	5	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
1,1,2,2-TETRACHLOROETHANE	5	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	5	<b>19 J</b>	<b>4.4 J</b>	<b>20 J</b>	< 0.50 UJ
1,1,2-TRICHLOROETHANE	1	<b>0.60 J</b>	< 0.50 UJ	<b>2.4 J</b>	< 0.50 UJ
1,1-DICHLOROETHANE	5	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
1,1-DICHLOROETHENE	5	<b>4.7 J</b>	<b>0.45 J</b>	<b>8.9 J</b>	< 0.50 UJ
1,2,4-TRICHLOROBENZENE	5	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
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 UJ</b>
1,2-DIBROMOETHANE	NL	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
1,2-DICHLOROBENZENE	3	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
1,2-DICHLOROETHANE	5	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
1,2-DICHLOROETHENE, TOTAL	5	<b>1.0 J</b>	<b>1.0 J</b>	<b>4.9 J</b>	< 1.0 UJ
1,2-DICHLOROPROPANE	1	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
1,3-DICHLOROBENZENE	3	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
1,4-DICHLOROBENZENE	3	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
1,4-DIOXANE (Method 8270D_SIM)	NL	<b>3.8</b>	<b>13</b>	<b>17</b>	< 0.17 U
2-BUTANONE	50	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ
2-HEXANONE	50	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ
4-METHYL-2-PENTANONE	NL	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ
ACETONE	50	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ
BENZENE	1	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
BROMODICHLOROMETHANE	50	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
BROMOFORM	50	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
BROMOMETHANE	5	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ
CARBON DISULFIDE	60	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
CARBON TETRACHLORIDE	5	<b>1.4 J</b>	< 0.50 UJ	<b>2.0 J</b>	< 0.50 UJ
CHLOROENZENE	5	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
CHLOROETHANE	5	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ
CHLOROFORM	7	< 0.50 UJ	< 0.50 UJ	<b>1.6 J</b>	< 0.50 UJ
CHLOROMETHANE	5	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ
CIS-1,2-DICHLOROETHENE	5	<b>1.0 J</b>	<b>1.0 J</b>	<b>4.9 J</b>	< 0.50 UJ
CIS-1,3-DICHLOROPROPENE	0.4	< <b>0.50 UJ</b>	< <b>0.50 UJ</b>	< <b>0.50 UJ</b>	< <b>0.50 UJ</b>
CYCLOHEXANE	NL	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
DIBROMOCHLOROMETHANE	5	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
DICHLORODIFLUOROMETHANE	5	< 1.0 UJ	< 1.0 UJ	<b>0.36 J</b>	< 1.0 UJ
ETHYLBENZENE	5	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
ISOPROPYLBENZENE	5	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
M- AND P-XYLENE	NL	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ
METHYL ACETATE	NL	< 0.75 UJ	< 0.75 UJ	< 0.75 UJ	< 0.75 UJ
METHYL CYCLOHEXANE	NL	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
METHYL TERT-BUTYL ETHER	10	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
METHYLENE CHLORIDE	5	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ
O-XYLENE	NL	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
STYRENE	5	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
TETRACHLOROETHENE	5	<b>0.84 J</b>	<b>1.0 J</b>	<b>2.6 J</b>	< 0.50 UJ
TOLUENE	5	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
TRANS-1,2-DICHLOROETHENE	5	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
TRANS-1,3-DICHLOROPROPENE	0.4	< <b>0.50 UJ</b>	< <b>0.50 UJ</b>	< <b>0.50 UJ</b>	< <b>0.50 UJ</b>
TRICHLOROETHENE	5	<b>640 J</b>	<b>600 J</b>	<b>5200 J</b>	<b>10 J</b>
TRICHLOROFLUOROMETHANE	5	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ
VINYL CHLORIDE	2	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ
XYLENES, TOTAL	5	< 1.5 UJ	< 1.5 UJ	< 1.5 UJ	< 1.5 UJ

**TABLE 2. ANALYTICAL DATA SUMMARY FOR  
WELLS SAMPLED BY RESOLUTION CONSULTANTS  
2015 OU2 GROUNDWATER INVESTIGATION**

Location	NYSDEC Groundwater Guidance or Standard Value (Note 1)	RE103D1	RE103D2	RE103D3
Sample Date		9/30/2015	9/30/2015	9/30/2015
Sample ID		RE103D1-GW- 093015	RE103D2-GW- 093015	RE103D3-GW- 093015
Sample type code		N	N	N
<b>VOC 8260C (ug/L)</b>				
1,1,1-TRICHLOROETHANE	5	< 0.50 U	< 0.50 U	< 0.50 U
1,1,2,2-TETRACHLOROETHANE	5	< 0.50 U	< 0.50 U	< 0.50 U
1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	5	<b>12</b>	<b>4.9</b>	<b>2.3 J</b>
1,1,2-TRICHLOROETHANE	1	<b>0.52 J</b>	<b>0.54 J</b>	< 0.50 U
1,1-DICHLOROETHANE	5	<b>1.0</b>	<b>0.72 J</b>	<b>0.50 J</b>
1,1-DICHLOROETHENE	5	<b>8.4</b>	<b>1.3</b>	<b>0.64 J</b>
1,2,4-TRICHLOROBENZENE	5	< 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>
1,2-DIBROMOETHANE	NL	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DICHLOROBENZENE	3	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DICHLOROETHANE	5	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DICHLOROETHENE, TOTAL	5	<b>3.6</b>	<b>1.5 J</b>	<b>0.96 J</b>
1,2-DICHLOROPROPANE	1	< 0.50 U	< 0.50 U	< 0.50 U
1,3-DICHLOROBENZENE	3	< 0.50 U	< 0.50 U	< 0.50 U
1,4-DICHLOROBENZENE	3	< 0.50 U	< 0.50 U	< 0.50 U
1,4-DIOXANE (Method 8270D_SIM)	NL	<b>24</b>	<b>2.1</b>	<b>1.0</b>
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	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ
BENZENE	1	< 0.50 U	< 0.50 U	< 0.50 U
BROMODICHLOROMETHANE	50	< 0.50 U	< 0.50 U	< 0.50 U
BROMOFORM	50	< 0.50 U	< 0.50 U	< 0.50 U
BROMOMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U
CARBON DISULFIDE	60	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
CARBON TETRACHLORIDE	5	<b>0.31 J</b>	<b>0.46 J</b>	< 0.50 U
CHLOROBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U
CHLOROFORM	7	<b>0.61 J</b>	<b>1.0</b>	<b>0.74 J</b>
CHLOROMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U
CIS-1,2-DICHLOROETHENE	5	<b>3.6</b>	<b>1.5</b>	<b>0.96 J</b>
CIS-1,3-DICHLOROPROPENE	0.4	< <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
DIBROMOCHLOROMETHANE	5	< 0.50 U	< 0.50 U	< 0.50 U
DICHLORODIFLUOROMETHANE	5	<b>0.26 J</b>	< 1.0 U	< 1.0 U
ETHYLBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U
ISOPROPYLBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U
M- AND P-XYLENE	NL	< 1.0 U	< 1.0 U	< 1.0 U
METHYL ACETATE	NL	< 0.75 U	< 0.75 U	< 0.75 U
METHYL CYCLOHEXANE	NL	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
METHYL TERT-BUTYL ETHER	10	< 0.50 U	< 0.50 U	< 0.50 U
METHYLENE CHLORIDE	5	< 2.5 U	< 2.5 U	< 2.5 U
O-XYLENE	NL	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
STYRENE	5	< 0.50 U	< 0.50 U	< 0.50 U
TETRACHLOROETHENE	5	<b>6.7</b>	<b>4.0</b>	<b>3.6 J</b>
TOLUENE	5	< 0.50 U	< 0.50 U	< 0.50 U
TRANS-1,2-DICHLOROETHENE	5	< 0.50 U	< 0.50 U	< 0.50 U
TRANS-1,3-DICHLOROPROPENE	0.4	< <b>0.50 U</b>	< <b>0.50 U</b>	< <b>0.50 U</b>
TRICHLOROETHENE	5	<b>860</b>	<b>830</b>	<b>470</b>
TRICHLOROFLUOROMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U
VINYL CHLORIDE	2	< 1.0 U	< 1.0 U	< 1.0 U
XYLENES, TOTAL	5	< 1.5 U	< 1.5 U	< 1.5 U

**TABLE 2. ANALYTICAL DATA SUMMARY FOR  
WELLS SAMPLED BY RESOLUTION CONSULTANTS  
2015 OU2 GROUNDWATER INVESTIGATION**

**Notes:**

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

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

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

Yellow highlighted values exceed Groundwater Standards or guidance value

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

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

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

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

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

**TABLE 3**  
**STABILIZED FIELD PARAMETERS FOR WELLS SAMPLED**  
**BY RESOLUTION CONSULTANTS**  
**2015 OU2 GROUNDWATER INVESTIGATION**

Well	Date	Temperature (°C)	pH	Specific Conductance (µS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Depth to water (ft bgs)	Flow rate (ml/min)
TT101D	9/29/2015	15.38	4.50	0.077	0.17	270.4	0.00	33.81	900
TT101D1	9/29/2015	14.83	4.21	0.078	0.60	366.0	0.00	35.38	750
TT101D2	9/29/2015	16.54	4.23	0.037	6.15	447.3	0.00	36.12	900
RE103D1	9/30/2015	15.52	4.41	0.082	4.42	344.6	0.01	40.26	700
RE103D2	9/30/2015	16.80	4.26	0.030	7.17	392.9	0.00	39.62	750
RE103D3	9/30/2015	16.00	4.42	0.026	4.15	312.1	0.00	40.62	500
RE104D1	9/25/2015	14.52	4.21	0.067	4.58	384.5	0.48	36.18	450
RE104D2	9/25/2015	14.73	5.16	0.019	6.65	279.7	1.39	40.56	500
RE104D3	9/25/2015	15.44	3.26	17.00	6.00	450.1	8.7	41.50	500
RE105D1	9/28/2015	14.86	4.48	0.099	2.07	364.4	0.59	38.68	500
RE105D2	9/28/2015	15.68	4.74	0.055	4.71	266.7	0.19	41.31	500
RE108D1	9/28/2015	15.32	4.41	0.086	6.76	374.8	0.32	41.01	750
RE108D2	9/28/2015	15.90	4.98	0.069	4.76	256.5	0.34	41.43	475
RE120D1	9/29/2015	16.22	4.19	0.096	1.41	354.4	0.15	37.55	500
RE120D2	9/29/2015	16.94	4.26	0.063	5.04	383.3	1.21	37.48	500
RE120D3	9/29/2015	16.40	4.58	0.025	3.28	271.7	10.6	37.28	600
RE122D1	9/30/2015	18.27	5.62	0.092	2.62	244.6	1.37	45.60	500
RE122D2	9/30/2015	18.01	4.40	0.087	4.47	364.2	1.64	42.38	500
RE122D3	9/30/2015	17.40	3.65	0.020	2.93	379.1	14.67	43.48	600
RE123D1	9/29/2015	19.82	5.01	0.097	7.64	261.9	5.26	46.62	500
RE123D2	9/29/2015	21.41	4.34	0.025	6.52	324.1	13.2	49.02	500
RE123D3	9/29/2015	19.54	5.12	0.041	0.46	-119.0	9.87	49.31	500

°C - degrees Celsius  
 µS/cm - Microsiemens per Centimeter  
 mg/L - milligrams per liter  
 mV - Millivolts  
 NTU - Nephelometric Turbidity Unit  
 ft bgs - feet below ground surface  
 ml/min - milliliters per minute

Table 4. Concentrations of Volatile Organic Compounds and 1,4-Dioxane in Monitoring Wells BPOW5-1, BPOW5-2, BPOW5-3, BPOW5-4, BPOW5-5, and BPOW5-6, Third Quarter 2015, Operable Unit 2 (Groundwater), Bethpage, New York.

CONSTITUENT (Units in µg/L)	Well: Sample ID: Date:	BPOW5-1 BPOW5-1 8/31/2015	BPOW5-2 BPOW5-2 9/1/2015	BPOW5-3 BPOW5-3 9/2/2015	BPOW5-4 BPOW5-4 9/14/2015	BPOW5-5 BPOW5-5 9/15/2015	BPOW5-6 BPOW5-6 9/16/2015
<b>Volatile Organic Compounds (VOCs)<sup>(1)</sup></b>							
1,1,1-Trichloroethane		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
1,1,2,2-Tetrachloroethane		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
1,1,2-trichloro-1,2,2-trifluoroethane		< 1.0	< 1.0	< 1.0	< 1.0	< 1.0	< 1.0
1,1,2-Trichloroethane		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
1,1-Dichloroethane		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50 J	< 0.50
1,1-Dichloroethene		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
1,2-Dichloroethane		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
1,2-Dichloropropane		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
2-Butanone (MEK)		< 5.0	< 5.0	< 5.0	< 5.0	< 5.0	< 5.0
2-Hexanone		< 2.0	< 2.0	< 2.0	< 2.0	< 2.0	< 2.0
4-methyl-2-pentanone (MIBK)		< 2.0	< 2.0	< 2.0	< 2.0	< 2.0	< 2.0
Acetone		< 5.0	< 5.0	< 5.0	< 5.0 B	< 5.0 B	< 5.0 B
Benzene		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Bromodichloromethane		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Bromoform		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Bromomethane		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Carbon Disulfide		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Carbon tetrachloride		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Chlorobenzene		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Chloroethane		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Chloroform		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Chloromethane		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
cis-1,2-dichloroethene		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
cis-1,3-dichloropropene		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50 J	< 0.50
Dibromochloromethane		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Ethylbenzene		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50 J	< 0.50
Methylene Chloride		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Styrene		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Tetrachloroethane		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Toluene		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
trans-1,2-dichloroethene		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
trans-1,3-dichloropropene		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Trichloroethylene		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Vinyl Chloride		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Xylene-o		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Xylenes - m,p		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
<b>Total VOCs<sup>(2)</sup></b>		0	0	0	0	0	0
<b>1,4-Dioxane<sup>(3)</sup></b>		< 0.21	< 0.21	<b>0.39</b>	<b>0.28</b>	<b>0.39</b>	< 0.23

**Notes and Abbreviations:**

- (1) Samples were analyzed for the TCL VOCs using USEPA Method 524.2.
- (2) Total VOCs are rounded to two significant figures.
- (3) Samples were analyzed for 1,4-Dioxane using USEPA Method 8270D SIM.

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

- Bold** Constituent detected
- TCL** Target Compound List
- VOC** Volatile Organic Compound
- USEPA** United States Environmental Protection Agency
- SIM** Selected Ion Monitoring
- µg/L** Micrograms per liter
- J** Constituent value is estimated
- B** Compound detected in associated blank sample
- <0.50** Compound not detected above its laboratory detection limit

Table 5. Concentrations of Volatile Organic Compounds and 1,4-Dioxane in Monitoring Wells BPOW6-1, BPOW6-2, BPOW6-3, BPOW6-4, BPOW6-5 and BPOW6-6, Third Quarter 2015, Operable Unit 2 (Groundwater), Bethpage, New York.

CONSTITUENT (Units in µg/L)	Well: Sample ID: Date:	BPOW6-1 BPOW6-1 9/3/2015	BPOW6-2 BPOW6-2 9/4/2015	BPOW6-3 BPOW6-3 9/8/2015	BPOW6-4 BPOW6-4 9/9/2015	BPOW6-4 BPOW6-4R <sup>(1)</sup> 9/9/2015	BPOW6-5 BPOW6-5 9/10/2015	BPOW6-6 BPOW6-6 9/11/2015
<b>Volatile Organic Compounds (VOCs)<sup>(2)</sup></b>								
1,1,1-Trichloroethane		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
1,1,2,2-Tetrachloroethane		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
1,1,2-trichloro-1,2,2-trifluoroethane		< 1.0	< 1.0	< 1.0	< 1.0	< 1.0	< 1.0	< 1.0
1,1,2-Trichloroethane		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
1,1-Dichloroethane		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
1,1-Dichloroethene		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
1,2-Dichloroethane		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
1,2-Dichloropropane		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
2-Butanone (MEK)		< 5.0	< 5.0	< 5.0	< 5.0	< 5.0	< 5.0	< 5.0
2-Hexanone		< 2.0	< 2.0	< 2.0	< 2.0	< 2.0	< 2.0	< 2.0
4-methyl-2-pentanone (MIK)		< 2.0	< 2.0	< 2.0	< 2.0	< 2.0	< 2.0	< 2.0
Acetone		< 5.0	< 5.0	< 5.0	< 5.0	< 5.0	< 5.0	< 5.0
Benzene		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Bromodichloromethane		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Bromoform		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Bromomethane		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Carbon Disulfide		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	<b>0.24 J</b>	< 0.50
Carbon tetrachloride		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Chlorobenzene		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Chloroethane		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Chloroform		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Chloromethane		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
cis-1,2-dichloroethene		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
cis-1,3-dichloropropene		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Dibromochloromethane		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Ethylbenzene		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Methylene Chloride		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Styrene		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Tetrachloroethene		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Toluene		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
trans-1,2-dichloroethene		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
trans-1,3-dichloropropene		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Trichloroethylene		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Vinyl Chloride		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Xylene-o		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
Xylenes - m,p		< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50	< 0.50
<b>Total VOCs<sup>(3)</sup></b>		<b>0</b>	<b>0</b>	<b>0</b>	<b>0</b>	<b>0</b>	<b>0.24</b>	<b>0</b>
<hr/>								
<b>1,4-Dioxane<sup>(4)</sup></b>		< 0.21	< 0.29	< 0.22	< 0.22	< 0.23	< 0.22	< 0.22

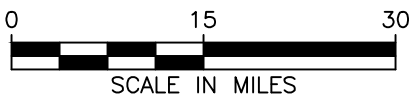
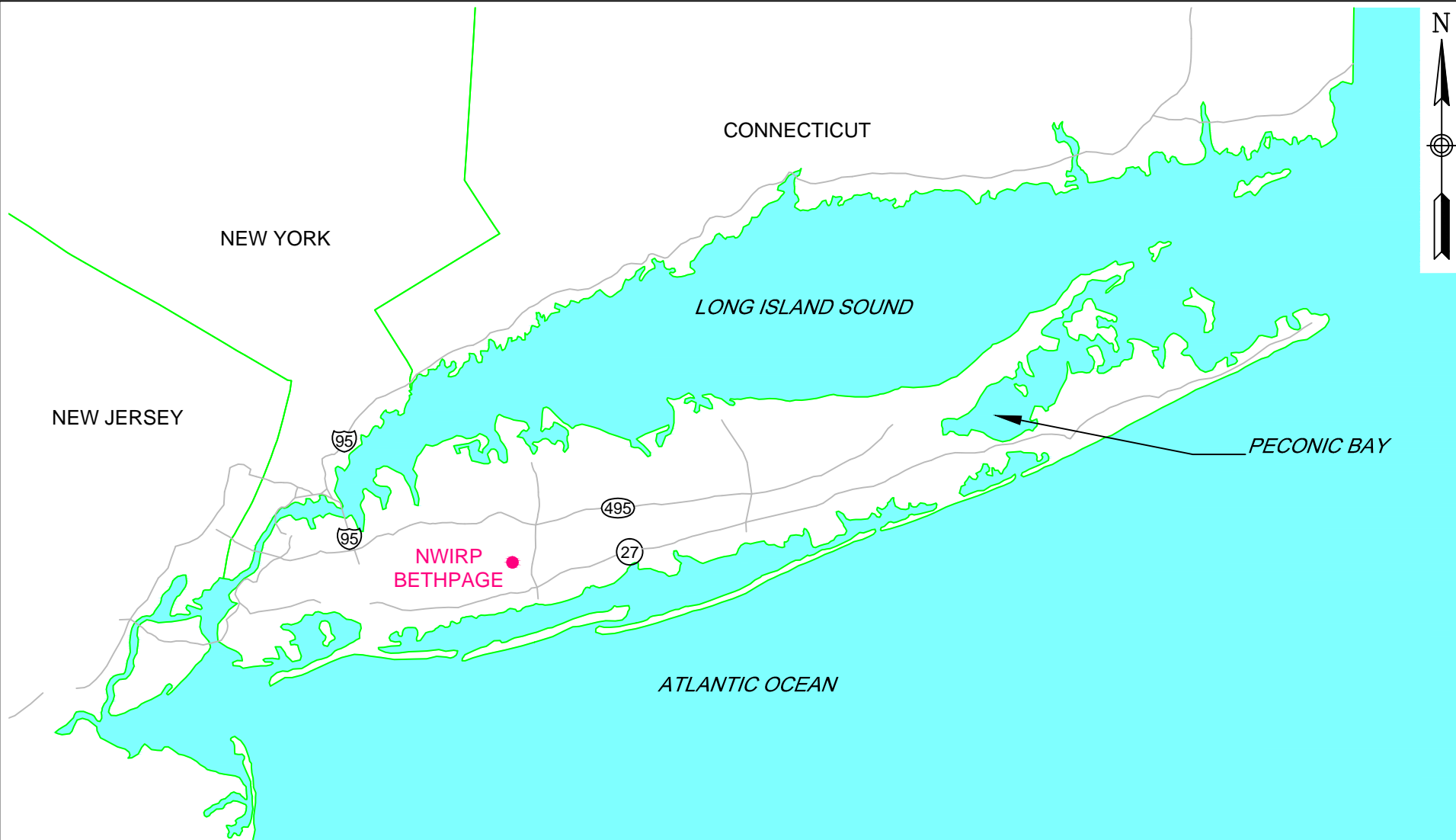
**Notes and Abbreviations:**

- <sup>(1)</sup> BPOW 6-4R is a blind duplicate sample.
- <sup>(2)</sup> Samples were analyzed for the TCL VOCs using USEPA Method 524.2.
- <sup>(3)</sup> Total VOCs are rounded to two significant figures.
- <sup>(4)</sup> Samples were analyzed for 1,4-Dioxane using USEPA Method 8270D SIM.

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

- Bold** Constituent detected
- TCL** Target Compound List
- VOC** Volatile Organic Compound
- USEPA** United States Environmental Protection Agency
- SIM** Selected Ion Monitoring
- µg/L** Micrograms per liter
- B** Compound detected in associated blank sample
- <0.50** Compound not detected above its laboratory detection limit

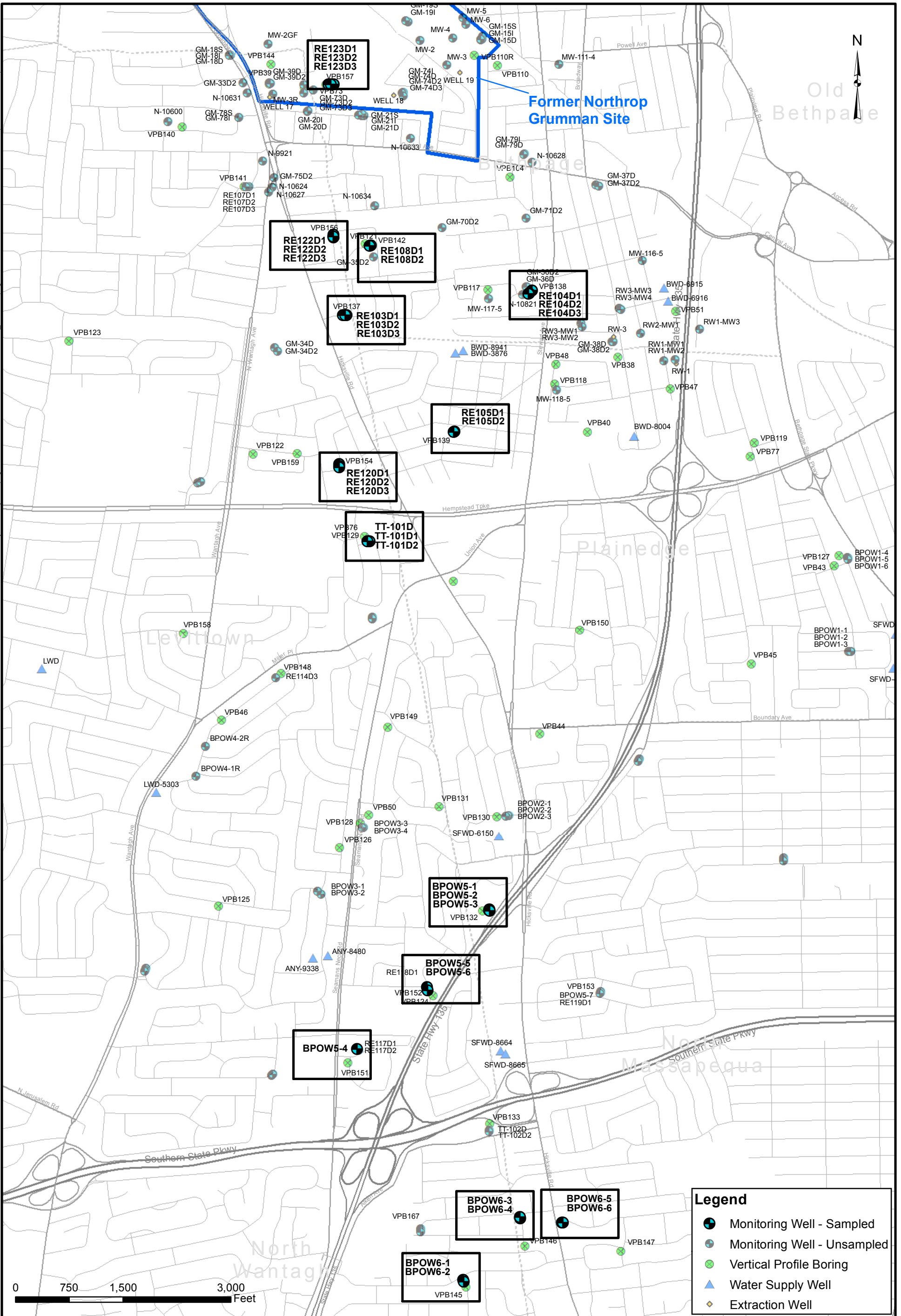
## **Figures**



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





**Legend**

- Monitoring Well - Sampled
- Monitoring Well - Unsampled
- Vertical Profile Boring
- Water Supply Well
- Extraction Well



**MONITORING WELL SAMPLE LOCATIONS**  
**SEPTEMBER 2015 GROUNDWATER SAMPLING**  
**NAVAL WEAPONS INDUSTRIAL RESERVE PLANT**  
**BETHPAGE, NEW YORK**

CONTRACT NUMBER N62470-11-D8013	CTO NUMBER WE 15
APPROVED BY EV	DATE 1/8/2016
APPROVED BY	DATE
FIGURE NO. <b>2</b>	REV 0

## **Appendices**

**Appendix A**

**Groundwater Sampling Forms – Resolution Consultants**



RESOLUTION  
CONSULTANTS

Well ID: RE10401

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 9/25/15 Time: Start 1315 am/pm  
 Project No: 60266526 Finish 1545 am/pm  
 Site Location: Hilltop  
 Weather Conds: Sunny 80° Collector(s): Paul Konezh

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

a. Total Well Length 375 c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material  
 4-inch PVC  
 b. Water Table Depth 36.77 d. Calculated System Volume (see back) 49.6L / 13.1 gal

### 2. WELL PURGE DATA

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

c. Field Testing Equipment used:

Make	Model	Serial Number
YSE	556	RF422118
Hanna	98203	64519

Time (24hr)	Volume Removed (Liters)	Temp. (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
13:51		14.70	3.69	0.068	9.85	390.5		450		OK
13:56		14.49	3.56	0.068	5.99	392.6		450		clear NO
14:01		14:39	3.59	0.070	4.68	396.4				
14:06		14.29	3.65	0.069	5.60	399.2	1.64	450		
14:12		14.34	3.75	0.067	5.25	398.2			36.22	
14:22	5g	14.23	3.80	0.067	4.96	399.8				
14:30		14:30	3.84	0.067	4.88	400.8			36.19	

d. Acceptance criteria pass/fail Yes No N/A (continued on back)

Has required volume been removed	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Has required turbidity been reached	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Have parameters stabilized	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>

If no or N/A - Explain below.

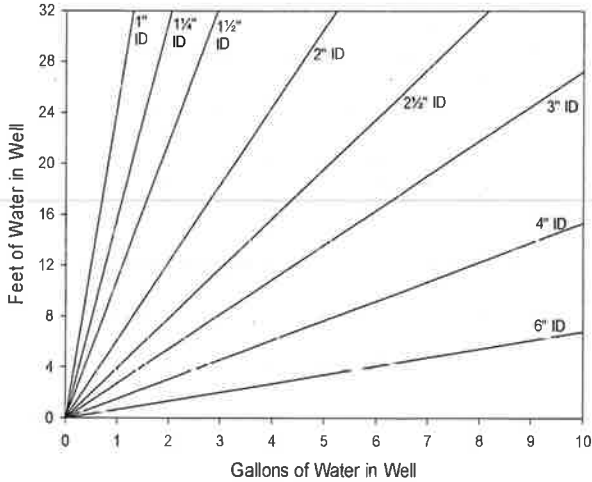
### 3. SAMPLE COLLECTION: Method: Geotech bladder pump with drop tube assembly

Sample ID	Container Type	No. of Containers	Preservation	Analysis Req.	Time
RE10401-GW-092515	40-mL vial	3	HCl	VOCs	1530
	1-L amber	2	none	1,4-Dioxane	

Comments: loggers Robium for NYS DEC

Signature: [Signature] Date: 9/25/15

# Purge Volume Calculation



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

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

Well ID:

(continued from front)

Time (24 hr)	Volume Removed (Liters)	Temp (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
14:36	6.5 g	14.37	3.96	0.067	4.84	395.6	0.95	450	36.20	Clear NO
14:43		14.60	4.14	0.067	4.78	387.9				
14:48		14.35	4.00	0.067	4.78	396.4				
14:53	7 g	14.46	3.96	0.067	4.71	399.7	0.46		36.22	
14:58		14.48	4.22	0.067	4.72	387.5				
15:03		14.48	3.95	0.067	4.70	401			36.18	
15:08	10.5 g	14.24	3.97	0.066	4.71	400.9	0.43			
15:17		14.85	4.19	0.068	4.58	387.5				
15:22	13 g	14.52	4.21	0.067	4.58	384.5	0.48			



RESOLUTION  
CONSULTANTS

Well ID: RE 10402

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 9 / 15 / 15 Time: Start 1315 am/pm  
 Project No: 60266526 Finish: 1545 am/pm  
 Site Location: Hilltop  
 Weather Conds: Sunny 80° Collector(s): Paul Karcch

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

a. Total Well Length 735 c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material  
4-inch PVC  
 b. Water Table Depth 40.78 d. Calculated System Volume (see back) \_\_\_\_\_

### 2. WELL PURGE DATA

a. Purge Method: Geotech bladder pump with drop tube assembly

b. Acceptance Criteria defined (see workplan)

- Temperature ± 3% - D.O. ± 10% (values >0.5 mg/L) Turbidity ± 10%
- pH ± 0.1 unit - ORP ± 10mV
- Sp. Cond. ± 3% - Drawdown < 0.3' Remove a minimum 1 screen volume

c. Field Testing Equipment used:	Make	Model	Serial Number
	<u>YSI</u>	<u>556</u>	<u>RFW 22120</u>
	<u>Hanna</u>	<u>HI 98702</u>	<u>645784</u>

Time (24hr)	Volume Removed (Liters)	Temp. (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
<u>1340</u>										
<u>1350</u>		<u>15.05</u>	<u>5.06</u>	<u>0.022</u>	<u>7.52</u>	<u>265.2</u>		<u>475</u>		
<u>1355</u>		<u>15.06</u>	<u>5.06</u>	<u>0.021</u>	<u>6.47</u>	<u>267.7</u>		<u>500</u>		
<u>1400</u>		<u>14.95</u>	<u>5.07</u>	<u>0.021</u>	<u>6.02</u>	<u>269.4</u>			<u>40.74</u>	
<u>1405</u>		<u>14.94</u>	<u>5.07</u>	<u>0.022</u>	<u>5.93</u>	<u>272.1</u>				
<u>1410</u>		<u>14.90</u>	<u>5.07</u>	<u>0.021</u>	<u>5.98</u>	<u>273.3</u>	<u>3.33</u>			
<u>1415</u>	<u>5 gal</u>	<u>14.89</u>	<u>5.04</u>	<u>0.020</u>	<u>6.24</u>	<u>274.9</u>	<u>3.34</u>			

d. Acceptance criteria pass/fail

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

If no or N/A - Explain below.

(continued on back)

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

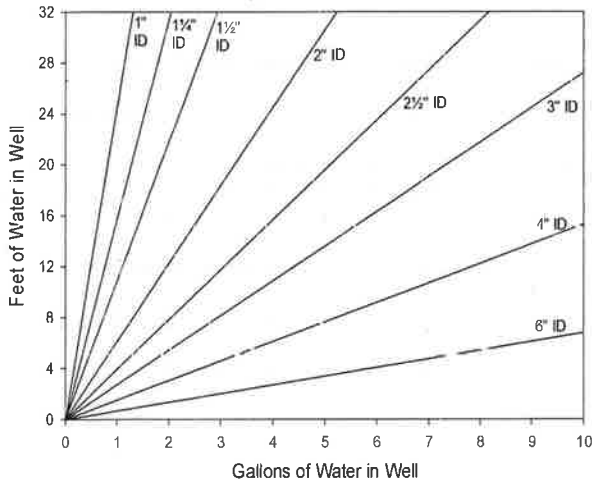
Sample ID	Container Type	No. of Containers	Preservation	Analysis Req.	Time
<u>RE10402-GW-092515</u>	<u>40-mL vial</u>	<u>3</u>	<u>HCl</u>	<u>VOCs</u>	<u>1520</u>
<u>RE10402-GW-092515</u>	<u>1-L amber</u>	<u>2</u>	<u>none</u>	<u>1,4-Dioxane</u>	<u>1520</u>

Radium for DEC (3-3L plastic)

Comments \_\_\_\_\_

Signature: Paul Karcch Date: 9/25/15

Purge Volume Calculation



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

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

Well ID: *RE10402 1340*

(continued from front)

Time (24 hr)	Volume Removed (Liters)	Temp (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
1420		14.84	5.11	0.019	6.42	276.3	3.47	500	40.62	
1425		14.83	5.12	0.019	6.47	276.8				
1430		14.81	5.14	0.019	6.52	277.4	3.11			
1435		14.84	5.13	0.019	6.55	277.7			40.63	
1440		14.82	5.15	0.019	6.56	277.7				
1445		14.83	5.15	0.019	6.62	278.3	2.73			
1450	<i>1092</i>	14.79	5.14	0.018	6.73	278.9		500		
1455		14.82	5.14	0.019	6.68	279.2			40.58	
1500		14.82	5.15	0.018	6.70	279.3				
1505		14.77	5.16	0.019	6.71	279.2				
1510		14.75	5.17	0.018	6.68	279.4	1.39			
1515		14.73	5.16	0.019	6.65	279.7		500	40.56	
1520										<i>Sample</i>



RESOLUTION  
CONSULTANTS

Well ID: RE104D3

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 9/25/15 Time: Start 1300 am/pm  
 Project No: 60266526 Finish 1545 am/pm  
 Site Location: \_\_\_\_\_  
 Weather Conds: \_\_\_\_\_ Collector(s): Leslie Baechler

**1. WATER LEVEL DATA: (measured from Top of Casing)**  
 a. Total Well Length 785 c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material  
 b. Water Table Depth 41.87 d. Calculated System Volume (see back) \_\_\_\_\_  
 4-inch PVC

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

Make	Model	Serial Number
<u>VST</u>	<u>556</u>	<u>24698</u>
<u>Hausza</u>	<u>98703</u>	<u>64518</u>

Time (24hr)	Volume Removed (Liters)	Temp. (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
1335	6	15.95	4.32	20	5.66	351.2	4.26	500	41.75	Clean/none
1340		15.88	4.17	19	4.65	365.1		500		"
1350	10	15.83	4.00	18	5.20	386.6		500		"
1355		15.79	3.99	18	5.46	389.4		500		"
1400	16	15.75	3.98	17	5.72	393.1	4.75	560	41.64	"
1405		15.61	3.92	17	5.80	401.7		500		"
1410	20	15.59	3.86	17	5.85	407.5		500		"

d. Acceptance criteria pass/fail

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

If no or N/A - Explain below.

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

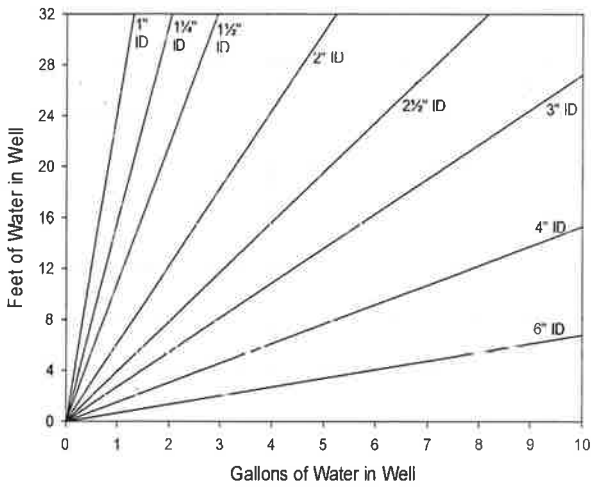
Sample ID	Container Type	No. of Containers	Preservation	Analysis Req.	Time
<u>RE104D3-GW-092515</u>	40-mL vial	3	HCl	VOCs	<u>1455</u>
	1-L amber	2	none	1,4-Dioxane	

Comments: Radium for NYSDEC

Signature: Leslie Baechler Date: 9/25/15



Purge Volume Calculation



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

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

Well ID:

(continued from front)

Time (24 hr)	Volume Removed (Liters)	Temp (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor	
1415	1	15.59	3.81	17	5.85	411.8	5.21	500		clean	none
1420	26	15.47	3.67	17	5.90	421.3		500	41.50	"	"
1425		15.54	3.61	17	5.96	425.8		500		"	"
1430	30	15.50	3.64	17	5.97	424.8		500		"	"
1435		15.58	3.50	17	5.93	426.0	8.89	500	41.50	"	"
1440	36	15.56	3.43	17	5.94	442.3	8.33	500		"	"
1445		15.43	3.39	17	5.99	450.6	8.52	500		"	"
1450	40	15.44	3.26	17	6.00	450.1	8.71	500	41.50	"	"
1455	SAMPLE TIME										



RESOLUTION  
CONSULTANTS

Well ID: RE105'01

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 9/28/15 Time: Start 9:00 am/pm  
 Project No: 60266526 Finish 11:20 am/pm  
 Site Location: RE105 @ Roosevelt + Franklin  
 Weather Conds: 80°F Overcast Collector(s): JC

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

a. Total Well Length 535 c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material  
4-inch PVC  
 b. Water Table Depth 39.82 d. Calculated System Volume (see back) 13.1 gal purge

### 2. WELL PURGE DATA

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

c. Field Testing Equipment used:	Make	Model	Serial Number
<u>YSI</u>		<u>556 MP5</u>	<u>2218</u>
<u>Hanna</u>		<u>HL 98703</u>	<u>U 64518x</u>

Time (24hr)	Volume Removed (Liters)	Temp. (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
920	-	15.15	3.69	0.103	8.26	394.4	1.56	500	38.34	clear/none
925	-	14.69	3.77	0.103	5.15	401.6	0.74	500	38.92	"
930	-	14.67	3.81	0.101	3.72	398.0	0.91	500	39.10	"
935	-	14.66	4.00	0.099	2.38	397.0	0.72	500	39.10	"
940	-	14.61	4.02	0.098	2.21	396.1	0.69	500	39.11	"
945	-	14.57	4.05	0.098	1.93	395.2	0.71	500	39.12	"
950	-	14.65	4.11	0.099	2.28	391.8	0.64	500	39.14	"

d. Acceptance criteria pass/fail Yes No N/A (continued on back)

Has required volume been removed	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>
Has required turbidity been reached	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Have parameters stabilized	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>

If no or N/A - Explain below.

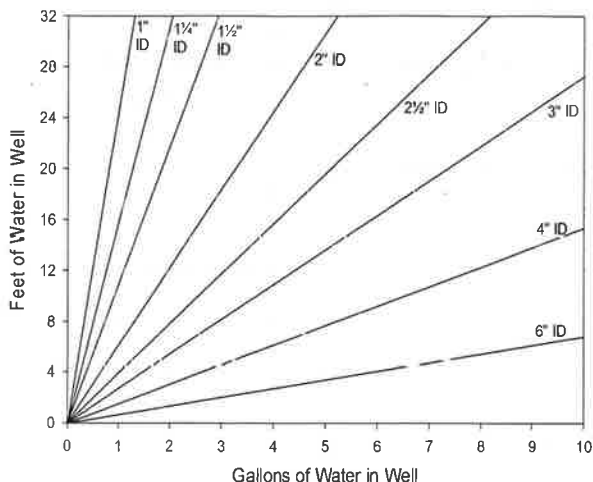
### 3. SAMPLE COLLECTION: Method: Geotech bladder pump with drop tube assembly

Sample ID	Container Type	No. of Containers	Preservation	Analysis Req.	Time
<u>RE105DI-GW-092815</u>	<u>40-mL vial</u>	<u>3</u>	<u>HCl</u>	<u>VOCs</u>	<u>1100</u>
<u>RE105DI-GW-092815</u>	<u>1-L amber</u>	<u>2</u>	<u>none</u>	<u>1,4-Dioxane</u>	<u>1100</u>

Comments Transducer present.

Signature \_\_\_\_\_ Date 9/28/15

Purge Volume Calculation



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

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

Well ID: RE105DI

(continued from front)

Time (24 hr)	Volume Removed (Liters)	Temp (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
9:55	5.9.1	14.67	4.18	0.098	2.21	391.0	0.68	500	39.14	clear / none
10:00	-	14.65	4.20	0.098	2.34	380.1	0.67	500	39.14	"
10:05	-	14.64	4.23	0.098	2.41	389.2	0.71	500	39.15	"
10:10	-	14.82	4.32	0.098	2.62	383.7	0.85	500	39.15	"
10:15	-	14.83	4.32	0.098	2.39	383.4	0.83	500	39.15	"
10:20	-	14.82	4.34	0.099	2.04	380.1	0.84	500	38.87	"
10:25	-	14.62	4.26	0.098	1.87	387.2	0.65	500	38.72	"
10:30	10.9.1	14.68	4.30	0.098	1.79	382.6	0.71	500	38.69	"
10:38	-	14.74	4.37	0.098	2.07	382.5	0.61	500	38.68	"
10:40	-	14.75	4.37	0.098	2.23	378.9	0.59	500	38.68	"
10:45	-	14.76	4.43	0.098	2.05	374.8	0.60	500	28.64	"
10:50	-	14.77	4.43	0.098	2.11	369.1	0.59	500	38.69	"
10:55	13.5	14.86	4.45	0.099	2.07	364.4	0.59	500	38.68	"



RESOLUTION CONSULTANTS

Well ID: RE10502

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 9/28/15 Time: Start 830 am/pm  
 Project No: 60266526 Finish 1130 am/pm  
 Site Location: Lincoln  
 Weather Conds: Sunny 65° Collector(s): Paul Karcch

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

a. Total Well Length 755 c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material  
 4-inch PVC  
 b. Water Table Depth 42.73 d. Calculated System Volume (see back) 41.62 / 13.1 gal

### 2. WELL PURGE DATA

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

c. Field Testing Equipment used:	Make	Model	Serial Number
	YST	556	RF22120
	Hanna	HI 98703	464518X

Time (24hr)	Volume Removed (Liters)	Temp. (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
900										slurb
910		15.42	4.14	0.052	6.19	323.0		475	43.15	
915		15.45	4.42	0.051	4.61	310.4		500	43.13	
920		15.35	4.49	0.052	3.79	305.8				
925		15.51	4.53	0.052	3.50	300.5				
930		15.62	4.54	0.052	4.07	300.5	0.91	500	43.18	
935	Signal	15.75	4.55	0.056	4.53	299.2				

d. Acceptance criteria pass/fail  
 Has required volume been removed  Yes  No  N/A  
 Has required turbidity been reached  Yes  No  N/A  
 Have parameters stabilized  Yes  No  N/A  
 If no or N/A - Explain below.

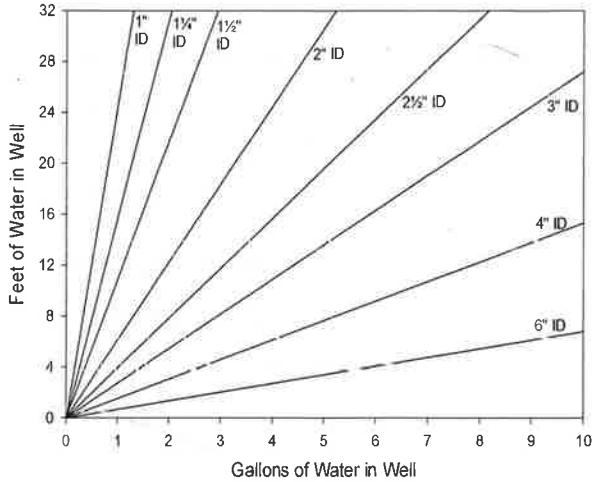
### 3. SAMPLE COLLECTION: Method: Geotech bladder pump with drop tube assembly

Sample ID	Container Type	No. of Containers	Preservation	Analysis Req.	Time
RE10502-GW-092815	40-mL vial	3	HCl	VOCs	1040
RE10502-GW-092815	1-L amber	2	none	1,4-Dioxane	1040

Comments: transducer

Signature: Paul Karcch Date: 9/28/15

Purge Volume Calculation



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

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

Well ID: RE105 DR 7900

(continued from front)

Time (24 hr)	Volume Removed (Liters)	Temp (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
940		15.79	4.61	0.056	4.63	294.1				
945		15.81	4.64	0.057	4.80	292.8	0.26		43.11	
950		15.67	4.67	0.056	4.80	287.3				
955		15.72	4.70	0.056	4.77	280.6				
1000		15.64	4.71	0.056	4.84	277.8			42.15	
1005		15.57	4.71	0.056	4.81	276.5	0.19	520		
1010	13gal	15.63	4.72	0.056	4.72	275.4				
1015		15.75	4.72	0.055	4.74	274.9			41.75	
1020		15.82	4.73	0.056	4.73	272.7				
1025		15.60	4.74	0.055	4.74	270.5			41.48	
1030		15.60	4.73	0.055	4.75	268.1			41.38	
1035	13gal	15.68	4.74	0.055	4.71	266.7			41.31	
1040										Sample



RESOLUTION CONSULTANTS

Well ID: RE10801

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 9/28/15 Time: Start 1310 am/pm  
 Project No: 60266526 Finish 1500 am/pm  
 Site Location: RE108 @ Corona + Ceril  
 Weather Conds: \_\_\_\_\_ Collector(s): JC

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

a. Total Well Length 535 c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material  
 b. Water Table Depth 41.00 d. Calculated System Volume (see back) 13.1 gal purge 4-inch PVC

### 2. WELL PURGE DATA

a. Purge Method: Geotech bladder pump with drop tube assembly

b. Acceptance Criteria defined (see workplan)

- Temperature ± 3% - D.O. ± 10% (values >0.5 mg/L) Turbidity ± 10%
- pH ± 0.1 unit - ORP ± 10mV
- Sp. Cond. ± 3% - Drawdown < 0.3' Remove a minimum 1 screen volume

c. Field Testing Equipment used:	Make	Model	Serial Number
<u>YSI</u>		<u>556 mps</u>	<u>22118</u>
<u>Hanna</u>		<u>HI 98703</u>	<u>U64518X</u>

Time (24hr)	Volume Removed (Liters)	Temp. (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
1320	-	17.04	4.11	0.093	9.86	357.6	5.14	750	41.02	clear from
1325	-	15.62	4.06	0.082	6.03	366.0	8.16	750	41.02	"
1330	-	15.39	4.07	0.087	6.14	368.4	3.89	750	41.02	"
1335	-	15.17	4.06	0.087	6.74	377.9	1.84	750	41.02	"
1340	-	15.15	4.09	0.086	6.86	379.6	0.79	750	41.02	"
1345	<u>5 gal</u>	15.16	4.10	0.086	6.89	380.4	0.53	750	41.02	"
1350	-	15.09	4.14	0.085	6.75	381.2	0.49	750	41.02	"

d. Acceptance criteria pass/fail

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

If no or N/A - Explain below.

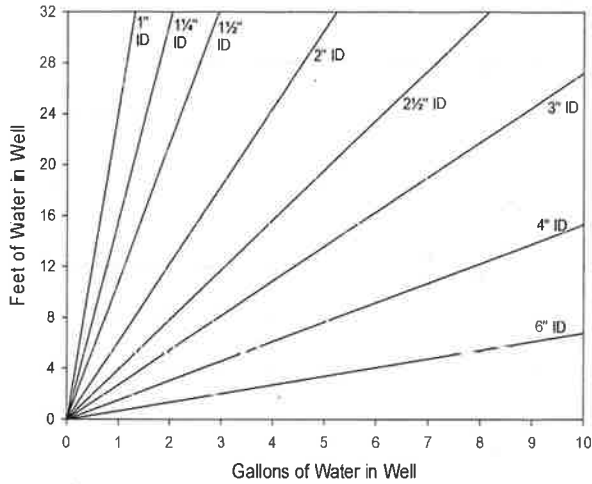
### 3. SAMPLE COLLECTION: Method: Geotech bladder pump with drop tube assembly

Sample ID	Container Type	No. of Containers	Preservation	Analysis Req.	Time
<u>RE10801-GW-092815</u>	<u>40-mL vial</u>	<u>3</u>	<u>HCl</u>	<u>VOCs</u>	<u>1435</u>
<u>RE10801-GW-092815</u>	<u>1-L amber</u>	<u>2</u>	<u>none</u>	<u>1,4-Dioxane</u>	<u>1435</u>
<del>RE10801-GW-092815</del>	<del>1-L plastic</del>		<del>HNO3</del>		

Comments: Liquinox  
Radium sample for NYSDEC

Signature: \_\_\_\_\_ Date: 9/28/15

Purge Volume Calculation



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

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

Well ID: RE108DI

(continued from front)

Time (24 hr)	Volume Removed (Liters)	Temp (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
1355	-	15.15	4.19	0.086	6.76	380.3	0.55	750	41.01	clear / none
1400	-	15.16	4.25	0.086	6.81	380.6	0.63	750	41.00	"
1405	-	15.16	4.26	0.086	6.82	381.4	0.47	750	41.00	"
1410	10.441	15.17	4.26	0.086	6.79	382.6	0.41	750	41.01	"
1415	-	15.18	4.32	0.086	6.74	378.3	0.49	750	40.99	"
1420	-	15.30	4.34	0.086	6.75	377.1	0.55	750	40.99	"
1425	-	15.25	4.31	0.086	6.82	375.9	0.53	750	40.99	"
1430	13.1401	15.22	4.41	0.086	6.76	374.8	0.32	750	41.01	"



**RESOLUTION  
CONSULTANTS**

Well ID: RE10802

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 9/28 / 15 Time: Start 1245 am/pm  
 Project No: 60266526 Finish        am/pm  
 Site Location: Carona & Ceil  
 Weather Conds: 80° partly sunny, recent rain Collector(s):       

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

a. Total Well Length 655 c. Length of Water Column        (a-b) Casing Diameter/Material 4-inch PVC  
 b. Water Table Depth 41.40 d. Calculated System Volume (see back)       

**2. WELL PURGE DATA**

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

c. Field Testing Equipment used:	Make	Model	Serial Number
<u>YSI</u>	<u>YSI</u>	<u>55C</u>	<u>RFW22120</u>
<u>Hydra</u>	<u>Hydra</u>	<u>HI98703</u>	<u>64518</u>

Time (24hr)	Volume Removed (Liters)	Temp. (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
<u>055</u>										<u>2600</u>
<u>1315</u>		<u>16.70</u>	<u>4.54</u>	<u>0.072</u>	<u>5.46</u>	<u>272.4</u>		<u>475</u>	<u>40.65</u>	
<u>1325</u>		<u>16.75</u>	<u>4.95</u>	<u>0.071</u>	<u>5.57</u>	<u>256.6</u>	<u>2.38</u>	<u>475</u>	<u>40.61</u>	
<u>1330</u>		<u>16.55</u>	<u>4.97</u>	<u>0.071</u>	<u>5.77</u>	<u>249.8</u>	<u>2.61</u>		<u>40.57</u>	
<u>1335</u>		<u>16.20</u>	<u>4.91</u>	<u>0.070</u>	<u>5.82</u>	<u>252.6</u>			<u>40.61</u>	
<u>1340</u>	<u>5 gal</u>	<u>16.23</u>	<u>4.94</u>	<u>0.069</u>	<u>5.81</u>	<u>252.0</u>				
<u>1345</u>		<u>16.13</u>	<u>4.95</u>	<u>0.069</u>	<u>5.78</u>	<u>252.3</u>	<u>0.53</u>		<u>40.95</u>	

d. Acceptance criteria pass/fail

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

If no or N/A - Explain below.

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

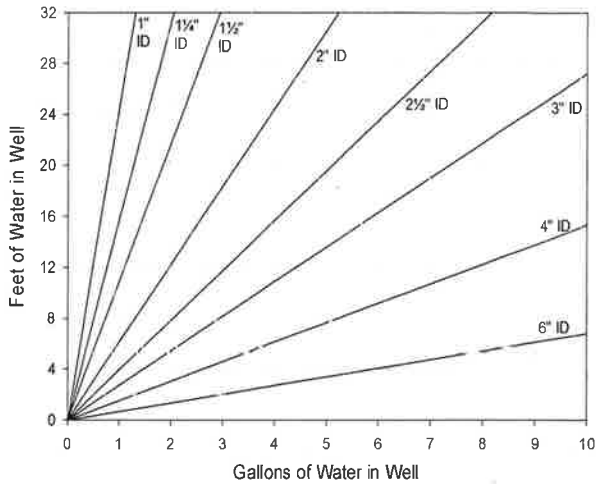
Sample ID	Container Type	No. of Containers	Preservation	Analysis Req.	Time
<u>RE10802-GW-092815</u>	<u>40-mL vial</u>	<u>3</u>	<u>HCl</u>	<u>VOCs</u>	<u>1450</u>
<u>      </u>	<u>1-L amber</u>	<u>2</u>	<u>none</u>	<u>1,4-Dioxane</u>	<u>1450</u>

Comments DEC arrived at 1345 collected radium samples  
transducer - liquinox decou

Signature Paul Hawth Date 9/28/15



Purge Volume Calculation



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

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

Well ID:

RE108DZ

(continued from front)

Time (24 hr)	Volume		pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
	Removed (Liters)	Temp (°C)								
1350		16.14	4.96	0.069	5.71	253.0		475	41.08	
1355		16.18	4.97	0.070	5.88	253.1				
1400		16.20	4.98	0.070	5.61	253.4	0.36		41.11	
1405		16.20	4.97	0.070	5.56	254.9		4.75	41.17	
1410		16.27	4.98	0.069	4.63	255.6	0.29		41.12	
1415	10 gal	16.15	4.98	0.069	4.80	256.2			41.51	
1420		16.04	4.98	0.069	4.73	256.3		475	41.50	
1425		16.07	4.99	0.069	4.73	256.2	0.32		41.43	
1430		16.12	4.98	0.070	4.74	256.2			41.41	
1435		16.04	4.98	0.069	4.75	256.2			41.42	
1440	13 gal	15.90	4.98	0.069	4.76	256.5	0.34		41.43	
1450										sample



Well ID: RE123 D1

RESOLUTION CONSULTANTS

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 9/29/15 Time: Start 9:55 am/pm  
 Project No: 60266526 Finish 12:10 am/pm  
 Site Location: MTA lot  
 Weather Conds: foggy 70° Collector(s): Paul Kureth

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

a. Total Well Length 545 c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material 4-inch PVC  
 b. Water Table Depth 45.43 d. Calculated System Volume (see back) 13.1 gal

### 2. WELL PURGE DATA

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

c. Field Testing Equipment used:	Make	Model	Serial Number
	<u>YSI</u>	<u>SS6 mps</u>	<u>22120</u>
	<u>Hann</u>	<u>HI 90203</u>	<u>U64518X</u>

Time (24hr)	Volume Removed (Liters)	Temp. (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
<u>830</u>										<u>0.1</u>
<u>955</u>									<u>46.62</u>	<u>RESET</u>
<u>1005</u>		<u>20.13</u>	<u>5.97</u>	<u>0.110</u>	<u>10.72</u>	<u>223.8</u>	<u>32.2</u>	<u>500</u>	<u>46.62</u>	
<u>1010</u>		<u>19.91</u>	<u>5.47</u>	<u>0.101</u>	<u>8.67</u>	<u>233.7</u>	<u>—</u>	<u>500</u>	<u>46.62</u>	<u>very cloudy</u>
<u>1015</u>		<u>19.88</u>	<u>5.30</u>	<u>0.097</u>	<u>7.38</u>	<u>233.5</u>	<u>178</u>	<u>500</u>	<u>46.62</u>	<u>cloudy</u>
<u>1020</u>		<u>19.82</u>	<u>5.19</u>	<u>0.096</u>	<u>7.74</u>	<u>242.0</u>	<u>—</u>	<u>500</u>	<u>46.62</u>	<u>"</u>
<u>1025</u>		<u>20.00</u>	<u>5.11</u>	<u>0.096</u>	<u>7.79</u>	<u>251.9</u>	<u>58.6</u>	<u>500</u>	<u>46.62</u>	<u>"</u>

d. Acceptance criteria pass/fail

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

If no or N/A - Explain below.

(continued on back)

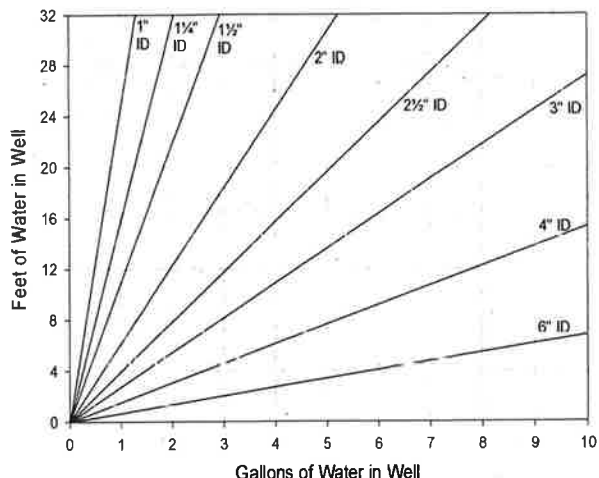
### 3. SAMPLE COLLECTION: Method: Geotech bladder pump with drop tube assembly

Sample ID	Container Type	No. of Containers	Preservation	Analysis Req.	Time
<u>RE123D1-GW-092915</u>	<u>40-mL vial</u>	<u>3</u>	<u>HCl</u>	<u>VOCs</u>	<u>1140</u>
<u>RE123D1-GW-092915</u>	<u>1-L amber</u>	<u>2</u>	<u>none</u>	<u>1,4-Dioxane</u>	<u>1140</u>

Comments Drop tube in sump, reset pump @ 955  
Lequinox

Signature \_\_\_\_\_ Date \_\_\_\_\_

# Purge Volume Calculation



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

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

Well ID: RE123DI

(continued from front)

Time (24 hr)	Volume Removed (Liters)	Temp (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
1030		19.87	5.08	0.096	7.76	256.2	-	500	46.62	clearing
1035	5 gal	19.84	5.07	0.096	7.68	257.7	13.4	500	46.62	"
1040		19.77	5.06	0.096	7.63	259.4	-	500	46.62	"
1045		19.98	5.04	0.097	7.63	262.4	9.11	500	46.62	clear
1050		19.90	5.05	0.096	7.64	261.7	-	500	46.62	"
1055		19.83	5.04	0.097	7.67	261.6	10.08	500	46.62	"
1100		19.77	5.03	0.096	7.60	262.5	-	500	46.62	"
1105	10 gal	19.77	5.03	0.096	7.66	262.6	9.61	500	46.62	"
1110		20.00	5.03	0.097	7.51	262.6	-	500	46.62	"
1115		20.10	5.03	0.097	7.62	262.8	7.24	500	46.62	"
1120		19.84	5.03	0.096	7.64	261.2	-	500	46.62	"
1125		19.94	5.02	0.097	7.65	262.5	5.26	500	46.62	"
1130	13 gal	20.04	5.02	0.097	7.68	261.5	-	500	46.62	"
1135		19.82	5.01	0.097	7.64	261.9	-	500	46.62	"



Well ID: RE12302

RESOLUTION CONSULTANTS

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 9/29/15 Time: Start 8:10 am/pm  
 Project No: 60266526 Finish \_\_\_\_\_ am/pm  
 Site Location: HTA Lab  
 Weather Conds: Foggy, overcast 75° Collector(s): \_\_\_\_\_

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

a. Total Well Length 660 c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material  
 4-inch PVC  
 b. Water Table Depth 49.01 d. Calculated System Volume (see back) 49.62 / 13.190

### 2. WELL PURGE DATA

a. Purge Method: Geotech bladder pump with drop tube assembly

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

c. Field Testing Equipment used:

Make	Model	Serial Number
<u>YSI</u>	<u>556</u>	<u>RRW 24698</u>
<u>HANNA</u>	<u>HI98703</u>	<u>64598X</u>

Time (24hr)	Volume Removed (Liters)	Temp. (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
9:10										
9:50	124/3g	21.02	5.08	0.030	5.83	265.6		500	49.03	cloudy
10:00		21.08	4.73	0.026	6.67	291.0			49.02	
10:05	5gal	21.18	4.49	0.025	7.10	298.7	169			
10:10		21.31	4.51	0.025	6.83	302.1				
10:15		21.33	4.46	0.025	6.38	305.9	20.2		49.01	
10:20		21.38	4.42	0.025	6.16	311.1	278		49.03	

d. Acceptance criteria pass/fail

Has required volume been removed	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Has required turbidity been reached	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Have parameters stabilized	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>

If no or N/A - Explain below.

(continued on back)

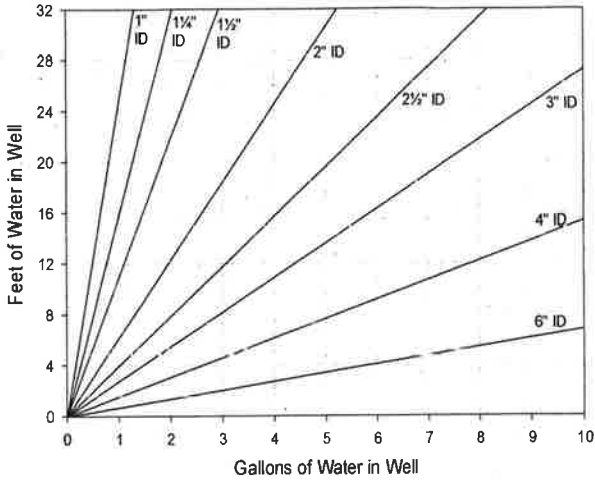
### 3. SAMPLE COLLECTION: Method: Geotech bladder pump with drop tube assembly

Sample ID	Container Type	No. of Containers	Preservation	Analysis Req.	Time
<u>RE12302-GW-092915</u>	<u>40-mL vial</u>	<u>3</u>	<u>HCl</u>	<u>VOCs</u>	<u>1110</u>
<u>RE12302-GW-092915</u>	<u>1-L amber</u>	<u>2</u>	<u>none</u>	<u>1,4-Dioxane</u>	<u>1110</u>

Comments: Site bottom with new tubing, lift 15 ft  
liquinox decor

Signature: Paul Kavitha Date: 9/29/15

**Purge Volume Calculation**



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

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

**Well ID:**

(continued from front)

Time (24 hr)	Volume Removed (Liters)	Temp (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
1025		21.35	4.45	0.025	6.01	311.2	14.5	500	49.02	
1030		21.34	4.44	0.025	6.02	318.6	17.1		49.04	
1035	10 gal	21.81	4.41	0.025	6.05	315.2	12.2		49.03	
1040		21.55	4.36	0.025	6.52	317.8	16.8		49.03	
1045		21.42	4.40	0.025	6.82	317.9	15.0		49.02	
1050		21.41	4.38	0.025	6.61	320.4			49.02	
1055		21.41	4.33	0.025	6.40	324.7	13.2		49.03	
1100		21.41	4.34	0.025	6.52	324.1			49.02	
1110										Sample



Well ID: RE123D3

RESOLUTION CONSULTANTS

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 9/29/15 Time: Start 810 am/pm  
 Project No: 60266526 Finish 1125 am/pm  
 Site Location: MTA Lot  
 Weather Conds: 80°F, muggy Collector(s): JC

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

a. Total Well Length 840 c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material  
 4-inch PVC  
 b. Water Table Depth 49.09 d. Calculated System Volume (see back) 13.1991

### 2. WELL PURGE DATA

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

Make	Model	Serial Number
YSI	556 MFS	22118
Hanna	HI 98703	U64518X

Time (24hr)	Volume Removed (Liters)	Temp. (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
910		19.26	4.56	0.031	7.35	291.7			49.20	ON
950	592	18.97	4.70	0.044	6.44	105.5		500		
955		18.95	4.69	0.044	6.11	48.4	30.3	500	49.28	cloudy / murky
1000		18.98	4.86	0.044	5.77	-25.6			49.29	
1005		19.11	4.94	0.044	2.49	-51.3			49.30	
1010		19.08	5.05	0.044	0.67	-90.1	31.8		49.30	
1015		19.66	5.20	0.044	0.61	-99.7			49.28	

d. Acceptance criteria pass/fail  
 Has required volume been removed  Yes  No  N/A  
 Has required turbidity been reached  Yes  No  N/A  
 Have parameters stabilized  Yes  No  N/A  
 If no or N/A - Explain below.

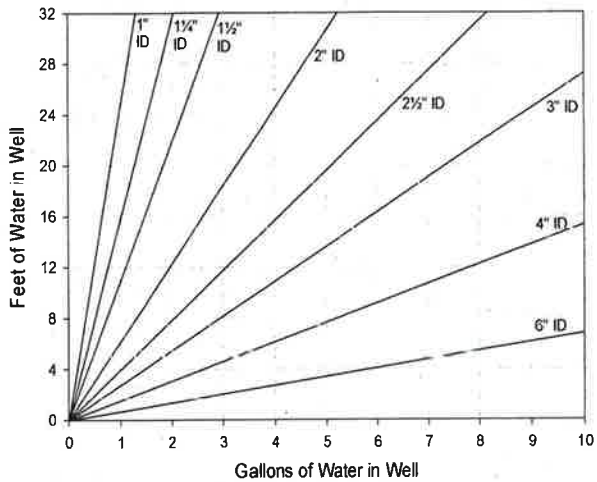
### 3. SAMPLE COLLECTION: Method: Geotech bladder pump with drop tube assembly

Sample ID	Container Type	No. of Containers	Preservation	Analysis Req.	Time
RE123D3-GW-092915	40-mL vial	3	HCl	VOCs	1100
RE123D3-GW-092915	1-L amber	2	none	1,4-Dioxane	1100

Comments: hit bottom with new tubing  
 liquor decon

Signature: Date: 9/29/2015

# Purge Volume Calculation



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

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

## Well ID:

(continued from front)

Time (24 hr)	Volume	Temp (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
	Removed (Liters)									
1020		19.42	5.18	0.043	0.57	-108.0	15.1		49.31	
1025	10 gal	19.31	5.05	0.043	0.54	-107.7				
1030		19.27	4.99	0.042	0.53	-108.1	11.8		49.32	
1035		19.20	5.07	0.042	0.50	-96.7			49.33	
1040		19.67	5.31	0.042	0.49	-117.1	9.87	500	49.32	
1045		19.46	5.11	0.041	0.48	-116.5			49.31	
1050		19.54	5.12	0.041	0.46	-119.0		500	49.31	



Well ID: RE12010

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 9/29/15 Time: Start 1515 am/pm  
 Project No: 60266526 Finish 1545 am/pm  
 Site Location: Shelley  
 Weather Conds: sunny 80° Collector(s): \_\_\_\_\_

**1. WATER LEVEL DATA: (measured from Top of Casing)**  
 a. Total Well Length 655 c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material  
 4-inch PVC

b. Water Table Depth 37.74 d. Calculated System Volume (see back) 20ft screen, 49.62/13.19a

**2. WELL PURGE DATA**

a. Purge Method: Geotech bladder pump with drop tube assembly

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

c. Field Testing Equipment used:

Make	Model	Serial Number
<u>YST</u>	<u>556</u>	<u>22118</u>
<u>Hanna</u>	<u>98703</u>	<u>64516</u>

Time (24hr)	Volume Removed (Liters)	Temp. (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
<u>14:15</u>	<u>-</u>	<u>18.85</u>	<u>4.65</u>	<u>0.105</u>	<u>10.95</u>	<u>294.4</u>	<u>1.85</u>	<u>400</u>	<u>37.69</u>	<u>Clear</u>
<u>14:20</u>	<u>-</u>	<u>17.49</u>	<u>4.63</u>	<u>0.101</u>	<u>3.90</u>	<u>289.6</u>	<u>3.72</u>	<u>375</u>	<u>37.68</u>	<u>Clear</u>
<u>14:25</u>	<u>-</u>	<u>17.31</u>	<u>4.52</u>	<u>0.099</u>	<u>2.20</u>	<u>295.4</u>	<u>-</u>	<u>425</u>	<u>37.68</u>	<u>Clear</u>
<u>14:30</u>	<u>-</u>	<u>17.16</u>	<u>4.36</u>	<u>0.097</u>	<u>1.82</u>	<u>313.5</u>	<u>-</u>	<u>450</u>	<u>37.68</u>	<u>Clear</u>
<u>14:35</u>	<u>-</u>	<u>17.20</u>	<u>4.12</u>	<u>0.098</u>	<u>1.88</u>	<u>326.5</u>	<u>-</u>	<u>400</u>	<u>37.65</u>	<u>Clear</u>
<u>14:40</u>	<u>-</u>	<u>17.05</u>	<u>4.03</u>	<u>0.098</u>	<u>1.88</u>	<u>336.0</u>	<u>-</u>	<u>375</u>	<u>37.65</u>	<u>Clear</u>
<u>14:45</u>	<u>-</u>	<u>17.03</u>	<u>4.06</u>	<u>0.098</u>	<u>1.87</u>	<u>337.4</u>	<u>0.93</u>	<u>-</u>	<u>37.64</u>	<u>Clear</u>

d. Acceptance criteria pass/fail

Has required volume been removed	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Has required turbidity been reached	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Have parameters stabilized	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>

If no or N/A - Explain below.

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

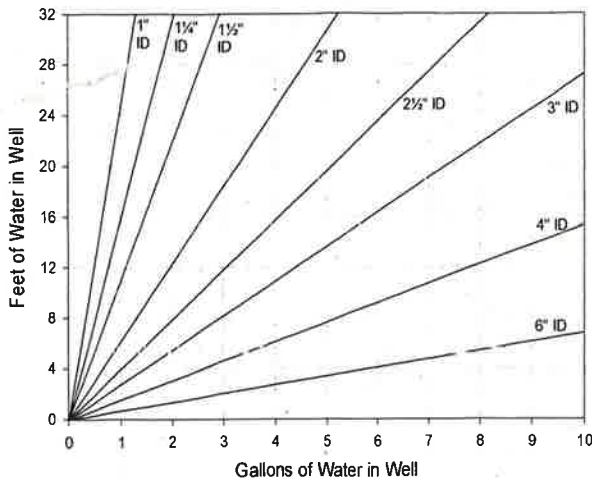
Sample ID	Container Type	No. of Containers	Preservation	Analysis Req.	Time
<u>RE12010-GW-092915</u>	<u>40-mL vial</u>	<u>3</u>	<u>HCl</u>	<u>VOCs</u>	<u>1600</u>
<u>RE12010-GW-092915</u>	<u>1-L amber</u>	<u>2</u>	<u>none</u>	<u>1,4-Dioxane</u>	

Comments Alcanax

Signature \_\_\_\_\_ Date \_\_\_\_\_



Purge Volume Calculation



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

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

Well ID:

(continued from front)

Time (24 hr)	Volume Removed (Liters)	Temp (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
14:50	5 Gal	16.77	4.07	0.098	1.86	338.1	-	450	37.62	Clear
14:55	-	16.92	4.10	0.098	1.80	338.4	-	450	37.64	Clear
15:00	-	16.82	4.12	0.098	1.78	339.1	-	475	37.60	Clear
15:05	-	16.78	4.10	0.098	1.74	339.9	-	475	37.60	Clear
15:10	-	16.66	4.10	0.097	1.72	345.3	0.87	500	37.60	Clear
15:15	-	16.56	4.11	0.097	1.67	347.7	-	-	-	Clear
15:20	-	16.39	4.11	0.097	1.62	349.8	0.65	500	37.60	Clear
15:30	10 Gal	16.38	4.10	0.097	1.57	352.3	0.58	500	37.60	Clear
15:35	-	16.46	4.10	0.097	1.52	346.3	1.2	500	37.59	Clear
15:40	-	16.32	4.11	0.096	1.47	354.8	1.7	475	37.6	Clear
15:45	-	16.30	4.16	0.097	1.46	353.1	0.65	475	37.6	Clear
15:50	-	16.30	4.19	0.096	1.42	353.5	1.2	475	37.59	Clear
15:55	18.5	16.22	4.19	0.096	1.41	354.4	0.15*	500	37.55	Clear



RESOLUTION  
CONSULTANTS

Well ID: RE12002

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 9/29/15 Time: Start 1315 am/pm  
 Project No: 60266526 Finish 1545 am/pm  
 Site Location: Shelley & Hahn  
 Weather Conds: Sunny 80° Collector(s):

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

a. Total Well Length 715 c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material  
 4-inch PVC  
 b. Water Table Depth 37.58 d. Calculated System Volume (see back) 49.64/13.1 gal

### 2. WELL PURGE DATA

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

c. Field Testing Equipment used:

Make	Model	Serial Number
YSI	556	RFW 24698
Hanna	HI 98703	LI 64518 X

Time (24hr)	Volume Removed (Liters)	Temp. (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
1355										NA
1400		18.43	4.94	0.061	5.64	299.4		450	37.56	
1410		18.08	4.65	0.060	4.12	318.8	5.60	500		
1415		17.86	4.57	0.061	4.22	327.4			37.57	
1420		17.70	4.45	0.062	4.60	339.2	3.16		37.58	
1425		17.62	4.34	0.063	4.86	352.4				
1430	5 gal	17.47	4.25	0.063	5.26	364.2		500	37.56	

d. Acceptance criteria pass/fail

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

If no or N/A - Explain below.

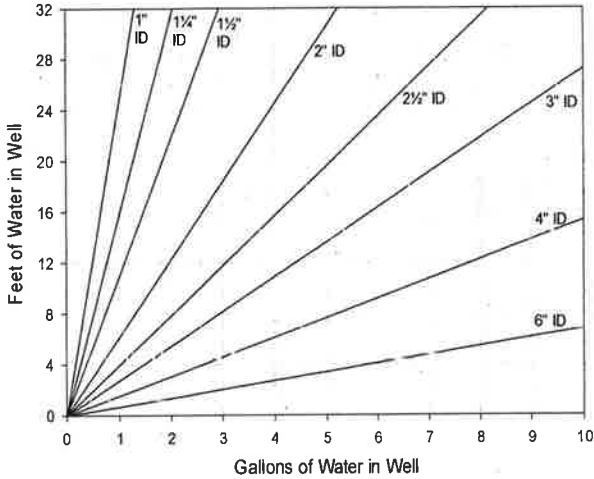
### 3. SAMPLE COLLECTION: Method: Geotech bladder pump with drop tube assembly

Sample ID	Container Type	No. of Containers	Preservation	Analysis Req.	Time
RE12002-GW-092915	40-mL vial	3	HCl	VOCs	1525
RE12002-GW-092915	1-L amber	2	none	1,4-Dioxane	1525

Comments: Alcomox

Signature: Paul Rucott Date: 9/29/15

Purge Volume Calculation



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

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

Well ID: ~~RE-12002~~ RE-12002 @ 1355

(continued from front)

Time (24 hr)	Volume Removed (Liters)	Temp (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
1435		17.43	4.25	0.063	5.31	366.4			37.56	
1440		17.36	4.36	0.063	5.21	363.4	1.39		37.53	
1445		17.31	4.33	0.063	5.17	368.5		500		
1450		17.27	4.31	0.063	5.00	372.0	1.34		37.52	
1455		17.27	4.33	0.063	4.93	373.8			37.52	
1500	10 gal	17.19	4.31	0.063	4.98	377.5	1.20	500	37.51	
1505		17.16	4.27	0.064	5.20	379.6				
1510		17.07	4.34	0.064	5.29	369.6			37.49	
1515		17.03	4.29	0.063	5.22	376.1	1.21		37.50	
1520		16.94	4.26	0.063	5.04	383.3			37.48	
1525										Sample



RESOLUTION CONSULTANTS

Well ID: RE120D3

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 9/29/15 Time: Start 1315 am/pm  
 Project No: 60266526 Finish 1545 am/pm  
 Site Location: Shelley @ Hahn.  
 Weather Conds: Sunny 70° Collector(s): JC

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

a. Total Well Length 265 c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material  
 b. Water Table Depth 38.46 d. Calculated System Volume (see back) 13.1 gal purge  
 4-inch PVC

### 2. WELL PURGE DATA

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

c. Field Testing Equipment used:

Make	Model	Serial Number
<u>YSI</u>	<u>556 MPS</u>	<u>22120</u>
<u>Hanna</u>	<u>HI 98703</u>	<u>U65418X</u>

Time (24hr)	Volume Removed (Liters)	Temp. (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
1400								400		OK
1410	-							600		
1420	-	17.15	4.08	0.021	3.70	289.2	14.0	600	37.54	clear
1440	5 gal	16.75	4.34	0.027	3.58	269.9	17.1	600	37.54	"
1445	-	16.75	4.48	0.026	3.53	267.2	16.9	600	37.54	"
1450	-	16.68	4.53	0.027	3.52	265.9	13.6	600	37.44	"
1455	-	16.65	4.56	0.027	3.52	265.8	12.3	600	37.44	"

d. Acceptance criteria pass/fail

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

If no or N/A - Explain below.

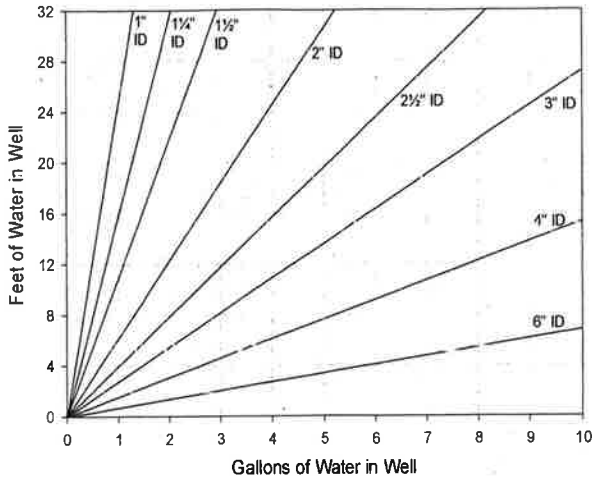
### 3. SAMPLE COLLECTION: Method: Geotech bladder pump with drop tube assembly

Sample ID	Container Type	No. of Containers	Preservation	Analysis Req.	Time
<u>RE120D3-GW-092915</u>	<u>40-mL vial</u>	<u>3</u>	<u>HCl</u>	<u>VOCs</u>	<u>1530</u>
<u>RE120D3-GW-092915</u>	<u>1-L amber</u>	<u>2</u>	<u>none</u>	<u>1,4-Dioxane</u>	<u>1530</u>

Comments Alconox

Signature \_\_\_\_\_ Date 9/29/15

Purge Volume Calculation



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

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

Well ID: RE120D3

(continued from front)

Time (24 hr)	Volume Removed (Liters)	Temp (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
1500	—	16.63	4.58	0.026	3.27	266.5	—	600	37.28	clear
1505	10.991	16.59	4.59	0.026	3.31	266.4	12.3	600	37.28	"
1510	—	16.50	4.60	0.025	3.28	267.9	12.9	600	37.28	"
1515	—	16.50	4.59	0.026	3.31	269.0	12.1	600	37.28	"
1520	—	16.42	4.59	0.025	3.31	270.7	10.0	600	37.28	"
1525	13.1991	16.40	4.58	0.025	3.28	271.7	10.6	600	37.28	"



Well ID: TT1010

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 9/29/15 Time: Start 1645 am/pm  
 Project No: 60266526 Finish 1900 am/pm  
 Site Location: Wentworth  
 Weather Conds: light rain 75° Collector(s): \_\_\_\_\_

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

a. Total Well Length 350 c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material  
4-inch PVC  
 b. Water Table Depth 33.89 d. Calculated System Volume (see back) 13.1 gal

### 2. WELL PURGE DATA

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

c. Field Testing Equipment used:	Make	Model	Serial Number
	<u>YSI</u>	<u>556</u>	<u>24698</u>
	<u>Hanna</u>	<u>98703</u>	<u>64518</u>

Time (24hr)	Volume Removed (Liters)	Temp. (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
1710										
1715								900		
1720		15.56	3.80	0.078	0.24	306	0.00		33.92	
1725		15.47	4.17	0.076	0.15	280.0			33.41	
1730	5.21	15.47	4.37	0.078	0.21	271.5	0.0	910	33.92	
1735		15.45	4.42	0.078	0.21	269.4				
1740		15.42	4.47	0.076	0.19	268.9				

d. Acceptance criteria pass/fail

Has required volume been removed	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Has required turbidity been reached	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Have parameters stabilized	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>

If no or N/A - Explain below.

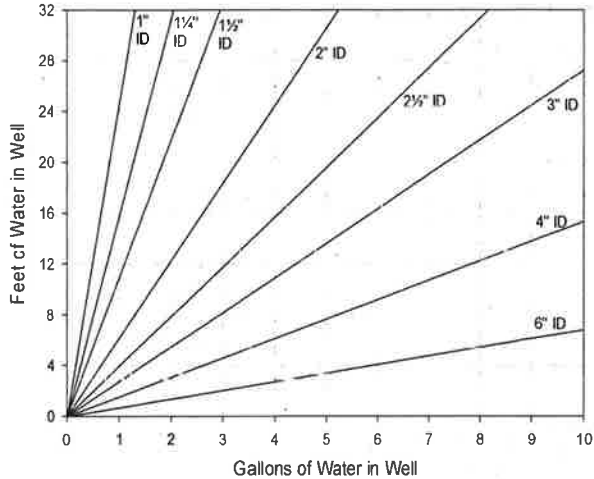
### 3. SAMPLE COLLECTION: Method: Geotech bladder pump with drop tube assembly

Sample ID	Container Type	No. of Containers	Preservation	Analysis Req.	Time
<u>TT1010-GW-092915</u>	40-mL vial	3	HCl	VOCs	1815
<u>Duplicate GW-092915</u>	1-L amber	2	none	1,4-Dioxane	1530

Comments \_\_\_\_\_

Signature Paul Karate Date 9/29/15

Purge Volume Calculation



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

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

Well ID: TT101D

(continued from front)

Time (24 hr)	Volume Removed (Liters)	Temp (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
1745		15.12	4.49	2.076	0.17	268.6			33.89	
1750		15.14	4.50	2.076	0.17	268.1				
1755	1.39	15.34	4.56	2.077	0.18	267.4			33.88	
1800		14.76	4.50	2.077	0.17	272.4	900		33.81	



RESOLUTION CONSULTANTS

Well ID: TT10101

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 9/29/15 Time: Start 1645 am/pm  
 Project No: 60266526 Finish 1900 am/pm  
 Site Location: West mouth  
 Weather Conds: light Rain 75° Collector(s): James Christopher

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

a. Total Well Length 595 c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material  
 b. Water Table Depth 36.41 d. Calculated System Volume (see back) 13.1 gal 4-inch PVC

### 2. WELL PURGE DATA

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

c. Field Testing Equipment used:	Make	Model	Serial Number
<u>YSI</u>	<u>YSI</u>	<u>556 MP3</u>	<u>22118</u>
<u>Hanna</u>	<u>Hanna</u>	<u>HI 98703</u>	<u>064518X</u>

Time (24hr)	Volume Removed (Liters)	Temp. (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
1715		16.41	3.87	0.081	4.41	379.1	-	600	35.40	clear
1720		15.35	3.72	0.079	0.53	383.0	-	600	35.40	"
1730		14.92	3.86	0.078	0.27	377.4	-	650	35.40	"
1745		14.86	4.05	0.078	0.60	371.8	-	750	35.40	"
1750		14.86	4.11	0.078	0.60	369.4	-	750	35.40	"
1755	10 gal	14.86	4.13	0.078	0.61	368.8	-	750	35.40	"
1800		14.83	4.17	0.078	0.61	367.6	0.00	750	35.40	"

d. Acceptance criteria pass/fail

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

If no or N/A - Explain below.

(continued on back)

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

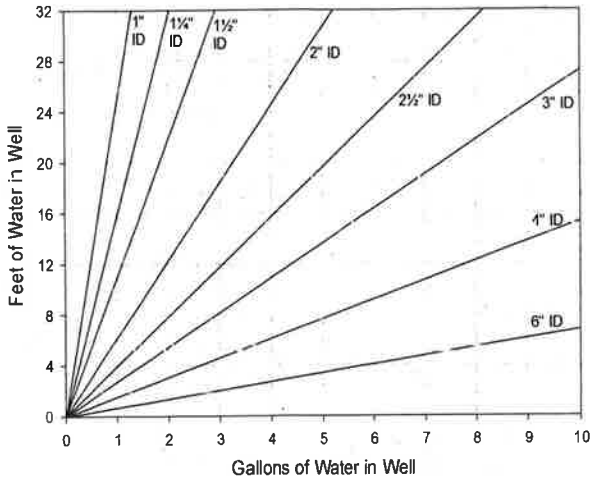
Sample ID	Container Type	No. of Containers	Preservation	Analysis Req.	Time
<u>TT10101-GW-092915</u>	40-mL vial	3	HCl	VOCs	1840
	1-L amber	2	none	1,4-Dioxane	1840

Comments \_\_\_\_\_

Signature \_\_\_\_\_ Date \_\_\_\_\_



### Purge Volume Calculation



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

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

Well ID: TT10:DI

(continued from front)

Time (24 hr)	Volume Removed (Liters)	Temp (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
1805	-	14.80	4.18	0.078	0.61	366.9	-	758	35.38	clear
1810	15.91	14.83	4.21	0.078	0.60	366.0	--	758	35.38	"



RESOLUTION CONSULTANTS

Well ID: TT10102

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 9/29/15 Time: Start 1815 am/pm  
 Project No: 60266526 Finish 1900 am/pm  
 Site Location: near north  
 Weather Conds: light rain 75° Collector(s): \_\_\_\_\_

1. WATER LEVEL DATA: (measured from Top of Casing)  
 a. Total Well Length 765 c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material  
 b. Water Table Depth 3601 d. Calculated System Volume (see back) 13.1 gal 4-inch PVC

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

c. Field Testing Equipment used:

Make	Model	Serial Number
<u>YST</u>	<u>556</u>	<u>22120</u>
<u>Hanna</u>	<u>18703</u>	<u>64518</u>

Time (24hr)	Volume Removed (Liters)	Temp. (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
<u>1705</u>								<u>900</u>		
<u>1710</u>		<u>15.91</u>	<u>4.18</u>	<u>0.037</u>	<u>3.14</u>	<u>356.6</u>				
<u>1715</u>		<u>15.90</u>	<u>4.17</u>	<u>0.037</u>	<u>2.40</u>	<u>358.6</u>	<u>0.0</u>			
<u>1720</u>		<u>15.86</u>	<u>4.17</u>	<u>0.037</u>	<u>3.35</u>	<u>364.0</u>			<u>36.12</u>	
<u>1725</u>		<u>15.83</u>	<u>4.15</u>	<u>0.036</u>	<u>3.86</u>	<u>371.8</u>			<u>36.14</u>	
<u>1730</u>	<u>5 gal</u>	<u>15.83</u>	<u>4.15</u>	<u>0.037</u>	<u>5.39</u>	<u>384.2</u>	<u>0.0</u>	<u>900</u>	<u>36.15</u>	
<u>1735</u>		<u>15.83</u>	<u>4.13</u>	<u>0.037</u>	<u>6.42</u>	<u>391.5</u>				

d. Acceptance criteria pass/fail

Has required volume been removed	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Has required turbidity been reached	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Have parameters stabilized	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>

If no or N/A - Explain below.

(continued on back)

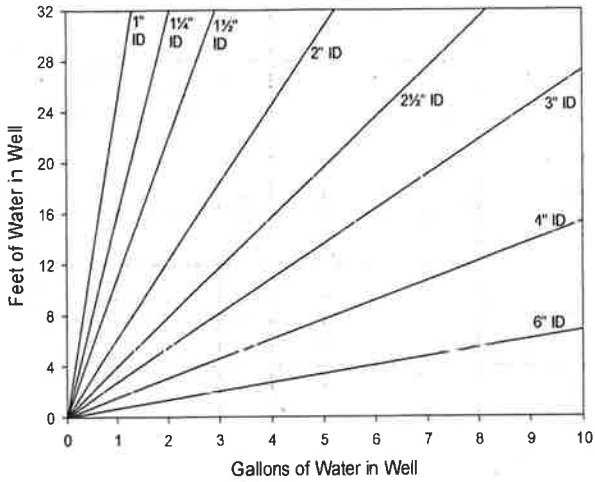
3. SAMPLE COLLECTION: Method: Geotech bladder pump with drop tube assembly

Sample ID	Container Type	No. of Containers	Preservation	Analysis Req.	Time
<u>TT10102-GW-692915</u>	<u>40-mL vial</u>	<u>3</u>	<u>HCl</u>	<u>VOCs</u>	<u>1830</u>
	<u>1-L amber</u>	<u>2</u>	<u>none</u>	<u>1,4-Dioxane</u>	<u>MS, MSD</u>

Comments \_\_\_\_\_

Signature Paul Kweth Date 9/29/15

Purge Volume Calculation



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

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

Well ID: TT 101 D2

(continued from front)

Time (24 hr)	Volume Removed (Liters)	Temp (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
1740		15.76	4.06	1.037	6.61	401.0			36.11	
1745		15.76	4.09	0.037	6.70	415.6				
1750		15.79	4.10	0.037	6.70	410.3				
1755	13 gal	15.75	4.11	0.036	6.67	430.4			35.10	
1800		16.54	4.23	0.037	6.15	447.3		9.00	36.12	
1830										Sample



RESOLUTION CONSULTANTS

Well ID: RE12201

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 9/30/15 Time: Start 8:35 am/pm  
 Project No: 60266526 Finish 10:25 am/pm  
 Site Location: Carbis  
 Weather Conds: rain 75° Collector(s): JC

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

a. Total Well Length 545 c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material  
 b. Water Table Depth 42.95 d. Calculated System Volume (see back) 13.1 gal Purge  
 4-inch PVC

### 2. WELL PURGE DATA

a. Purge Method: Geotech bladder pump with drop tube assembly

b. Acceptance Criteria defined (see workplan)

- Temperature ± 3% - D.O. ± 10% (values >0.5 mg/L) Turbidity ± 10%
- pH ± 0.1 unit - ORP ± 10mV
- Sp. Cond. ± 3% - Drawdown < 0.3' Remove a minimum 1 screen volume

c. Field Testing Equipment used:

Make	Model	Serial Number
<u>YSI</u>	<u>556 MP3</u>	<u>22120</u>
<u>Hanna</u>	<u>HI97803</u>	<u>82</u>
<u>Scientific</u>	<u>Micro TPN</u>	<u>U64518X</u>

Time (24hr)	Volume Removed (Liters)	Temp. (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
<del>8:45</del>								<u>500</u>		
<del>8:45</del>										
<u>8:45</u>		<u>17.93</u>	<u>4.91</u>	<u>0.089</u>	<u>4.19</u>	<u>270.7</u>	<u>1134</u>		<u>42.98</u>	<u>muridity</u>
<u>8:50</u>		<u>16.9</u>	<u>5.06</u>	<u>0.078</u>	<u>3.01</u>	<u>274.5</u>				
<u>8:55</u>		<u>17.26</u>	<u>5.26</u>	<u>0.091</u>	<u>3.77</u>	<u>268.0</u>		<u>500</u>	<u>42.82</u>	
<u>9:00</u>		<u>18.14</u>	<u>5.40</u>	<u>0.091</u>	<u>3.60</u>	<u>259.2</u>	<u>-</u>	<u>500</u>	<u>-</u>	<u>clear</u>
<u>9:05</u>		<u>18.28</u>	<u>5.42</u>	<u>0.091</u>	<u>3.52</u>	<u>257.9</u>	<u>3.06</u>	<u>500</u>	<u>42.82</u>	<u>"</u>

d. Acceptance criteria pass/fail

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

If no or N/A - Explain below.

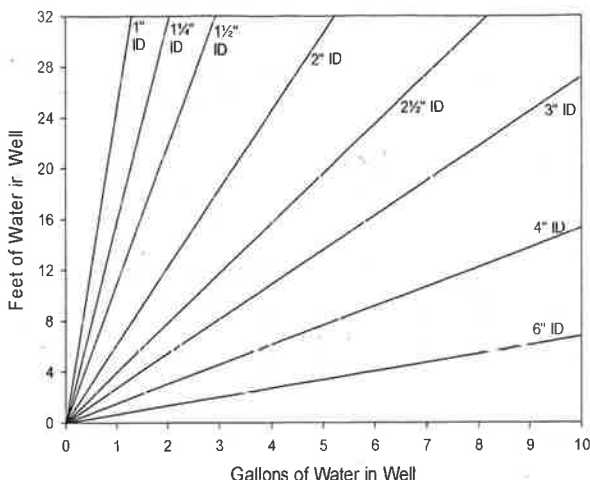
### 3. SAMPLE COLLECTION: Method: Geotech bladder pump with drop tube assembly

Sample ID	Container Type	No. of Containers	Preservation	Analysis Req.	Time
<u>RE12201-GW-092015</u>	<u>40-mL vial</u>	<u>3</u>	<u>HCl</u>	<u>VOCs</u>	<u>1015</u>
<u>RE12201-GW-092015</u>	<u>1-L amber</u>	<u>2</u>	<u>none</u>	<u>1,4-Dioxane</u>	<u>1015</u>

Comments muridity is giving poor turbidity  
luminox

Signature \_\_\_\_\_ Date 9/30/2015

Purge Volume Calculation



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

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

Well ID: REZZDI

(continued from front)

Time (24 hr)	Volume Removed (Liters)	Temp (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
910	5 gal	18.28	5.47	0.092	3.47	255.8	-	500	42.82	clear / none
915	-	18.06	5.48	0.091	3.35	253.8	-	500	42.77	"
920	-	18.14	5.52	0.091	3.25	253.2	4.27	500	42.68	"
925	-	18.13	5.52	0.092	3.18	252.7	-	500	42.68	"
930	-	18.07	5.55	0.091	3.11	251.1	-	500	42.48	"
935	-	18.03	5.55	0.091	3.10	251.8	-	500	42.68	"
940	-	18.07	5.55	0.092	3.01	250.4	-	500	42.68	"
945	10 gal	18.12	5.57	0.092	2.89	250.0	3.21	500	42.58	"
950	-	18.20	5.60	0.092	2.84	247.4	3.01	500	42.68	"
955	-	18.27	5.59	0.092	2.76	247.0	3.41	500	42.61	"
1000	-	18.32	5.61	0.092	2.68	245.7	-	500	42.59	"
1005	13.5 gal	18.21	5.62	0.092	2.62	244.6	1.37	500	45.60	"



RESOLUTION CONSULTANTS

Well ID: RC12202

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 9/30/15 Time: Start 115 am/pm  
 Project No: 60266526 Finish 1100 am/pm  
 Site Location: Curtis  
 Weather Conds: rain 75 Collector(s): \_\_\_\_\_

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

a. Total Well Length 615 c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material  
4-inch PVC  
 b. Water Table Depth 43.18 d. Calculated System Volume (see back) \_\_\_\_\_

### 2. WELL PURGE DATA

a. Purge Method: Geotech bladder pump with drop tube assembly

b. Acceptance Criteria defined (see workplan)

- Temperature	± 3%	- D.O.	± 10% (values >0.5 mg/L)	Turbidity	± 10%
- pH	± 0.1 unit	- ORP	± 10mV		
- Sp. Cond.	± 3%	- Drawdown	< 0.3'	Remove a minimum 1 screen volume	

c. Field Testing Equipment used:

Make	Model	Serial Number
<u>VSI</u>	<u>556</u>	<u>24698</u>
<u>Micro</u>	<u>20000</u>	<u>201205051</u>

Time (24hr)	Volume Removed (Liters)	Temp. (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
<u>8:28</u>										
<u>8:33</u>		<u>18.41</u>	<u>4.72</u>	<u>0.090</u>	<u>2.07</u>	<u>355</u>				
<u>8:47</u>		<u>17.92</u>	<u>4.64</u>	<u>0.086</u>	<u>2.68</u>	<u>359.7</u>			<u>43.12</u>	
<u>9:00</u>		<u>18.46</u>	<u>4.68</u>	<u>0.087</u>	<u>3.63</u>	<u>352.2</u>		<u>500</u>		<u>parameters stopped</u>
<u>9:05</u>		<u>20.92</u>	<u>4.60</u>	<u>0.087</u>	<u>3.73</u>	<u>348.7</u>	<u>0.98</u>			
<u>9:10</u>		<u>18.28</u>	<u>4.61</u>	<u>0.087</u>	<u>4.24</u>	<u>351.4</u>			<u>42.92</u>	
<u>9:15</u>		<u>18.00</u>	<u>4.58</u>	<u>0.087</u>	<u>4.31</u>	<u>352.0</u>				

d. Acceptance criteria pass/fail

	Yes	No	N/A	(continued on back)
Has required volume been removed	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Has required turbidity been reached	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Have parameters stabilized	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	

If no or N/A - Explain below.

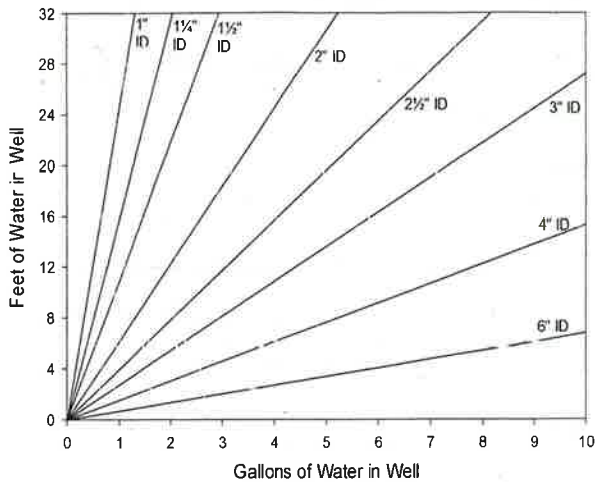
### 3. SAMPLE COLLECTION: Method: Geotech bladder pump with drop tube assembly

Sample ID	Container Type	No. of Containers	Preservation	Analysis Req.	Time
	40-mL vial	3	HCl	VOCs	<u>1020</u>
	1-L amber	2	none	1,4-Dioxane	

Comments Summary from

Signature \_\_\_\_\_ Date \_\_\_\_\_

Purge Volume Calculation



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

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

Well ID: RE12202

(continued from front)

Time (24 hr)	Volume Removed (Liters)	Temp (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
930	5991	18.24	<del>4.35</del> 4.57	0.087	4.30	350.6	5.00	500	42.82	
935		18.09	4.56	0.087	4.23	349.9	5.04			
930		18.08	4.53	0.087	4.30	353.0				
935		18.08	4.50	0.087	4.33	353.8				
940		18.07	4.49	0.087	4.38	354.2		500	42.62	
945		18.19	4.48	0.087	4.35	354.9	1.74			
950	Logal	17.98	4.47	0.087	4.38	355.4			42.58	
955		18.07	4.42	0.087	4.37	357.8		500		
1000		17.95	4.41	0.087	4.46	359.4			42.48	
1005		17.99	4.41	0.087	4.44	360.4	2.02	500	42.42	
1010		18.02	4.40	0.087	4.51	362.9			42.41	
1015		18.01	4.40	0.087	4.47	364.2	1.84		42.38	
1020										sample



Well ID: RE12203

**RESOLUTION CONSULTANTS**

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 9/30/15 Time: Start 8:00 am/pm  
 Project No: 60266526 Finish 11:00 am/pm  
 Site Location: sub 15  
 Weather Conds: rain 75° Collector(s): \_\_\_\_\_

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

a. Total Well Length 740 c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material  
4-inch PVC  
 b. Water Table Depth 44.24 d. Calculated System Volume (see back) \_\_\_\_\_

### 2. WELL PURGE DATA

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

Make	Model	Serial Number
<u>Y51</u>	<u>556 MPS</u>	<u>RFW22118</u>
<u>Micro TP4</u>	<u>26000</u>	<u>201205051</u>

Time (24hr)	Volume Removed (Liters)	Temp. (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
<u>8:20</u>										<u>OK</u>
<u>8:28</u>		<u>14.81</u>	<u>4.50</u>	<u>0.028</u>	<u>6.52</u>	<u>369.0</u>	<u>92.94</u>	<u>400</u>	<u>44.28</u>	
<u>8:47</u>			<u>4.05</u>	<u>0.025</u>	<u>3.10</u>	<u>347.5</u>		<u>300</u>		
<u>8:55</u>			<u>4.22</u>	<u>0.025</u>	<u>2.86</u>	<u>338.4</u>	<u>34.86</u>	<u>3.50</u>	<u>54.18</u>	
<u>9:00</u>		<u>20.22</u>	<u>4.15</u>	<u>0.025</u>	<u>2.81</u>	<u>333.5</u>				
<u>9:06</u>	<u>4g</u>	<u>20.39</u>	<u>4.22</u>	<u>0.023</u>	<u>2.82</u>	<u>338.0</u>			<u>44.18</u>	<u>Clear NO</u>
<u>9:11</u>		<u>20.45</u>	<u>4.26</u>	<u>0.022</u>	<u>2.84</u>	<u>338.4</u>	<u>11.74</u>	<u>275</u>		

d. Acceptance criteria pass/fail

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

If no or N/A - Explain below.

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

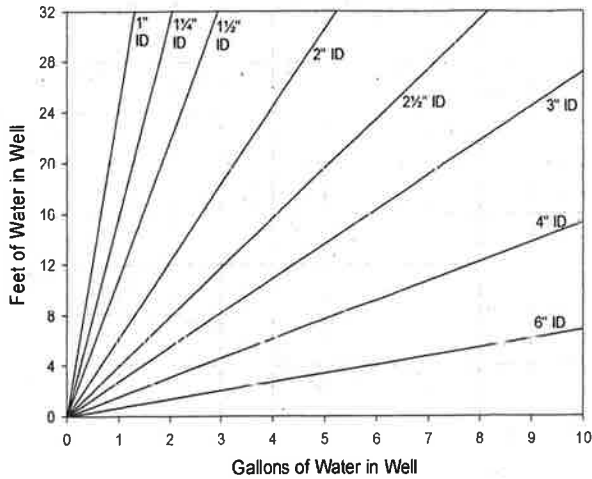
Sample ID	Container Type	No. of Containers	Preservation	Analysis Req.	Time
	<u>40-mL vial</u>	<u>3</u>	<u>HCl</u>	<u>VOCs</u>	<u>1036</u>
	<u>1-L amber</u>	<u>2</u>	<u>none</u>	<u>1,4-Dioxane</u>	

Comments Luminox decom

Signature [Signature] Date 9/30/15



Purge Volume Calculation



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

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

Well ID:

(continued from front)

Time (24 hr)	Volume Removed (Liters)	Temp (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
9:17		20.94	4.35	6.022	3.05	336.4			44.18	
9:22		20.85	4.33	0.022	3.06	337.4	8.15			
				Pull Pump - Replace Bladder						
9:43		18.65	4.13	0.020	5.12	300.5		600	43.28	
9:48	6.5g	17.59	3.18	0.020	3.91	404.4	13.42			
9:56	7.5g	17.76	3.42	0.021	3.20	393.1	12.95	640		
10:01	9.0g	17.55	3.41	0.020	3.06	394.9				
10:06	10g	17.46	3.52	0.020	2.89	386.8	13.17			
10:11		17.48	3.57	0.020	2.84	387.5			43.46	
10:16	11.5g	17.42	3.57	0.020	2.96	387.4	13.51			
10:21		17.57	3.61	0.020	2.96	382.9				
10:26	13	17.41	3.60	0.020	2.96	383.7	15.13	600	43.48	
10:31	14.7	17.40	3.65	0.020	2.93	378.1	14.67			OK
										Sample 10:36



RESOLUTION CONSULTANTS

Well ID: RE10301

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 9/30/15 Time: Start 1200 am/pm  
 Project No: 60266526 Finish 1330 am/pm  
 Site Location: Avoca  
 Weather Conds: overcast, drizzle PS Collector(s):

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

a. Total Well Length 645 c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material 4-inch PVC  
 b. Water Table Depth 40.19 d. Calculated System Volume (see back) 1586 = 9.8 gal / 37.16

### 2. WELL PURGE DATA

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

c. Field Testing Equipment used:

Make	Model	Serial Number
YSI	556	RFW 22118
Micro	20000	20125051

Time (24hr)	Volume Removed (Liters)	Temp. (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
1215										011
1225		15.74	4.29	0.076	1.30	326.0		700		
1230		15.69	4.37	0.076	0.94	323.0	7.59		40.27	
1235		15.47	4.28	0.079	3.11	335.5		700		
1240	5 gal	15.67	4.16	0.080	4.16	347.6	2.23		40.25	
1245		15.53	4.16	0.080	4.20	350.2				
1250	7 gal	15.30	4.12	0.080	4.27	353.7		700	40.24	

d. Acceptance criteria pass/fail Yes No N/A (continued on back)

Has required volume been removed	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Has required turbidity been reached	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Have parameters stabilized	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>

If no or N/A - Explain below.

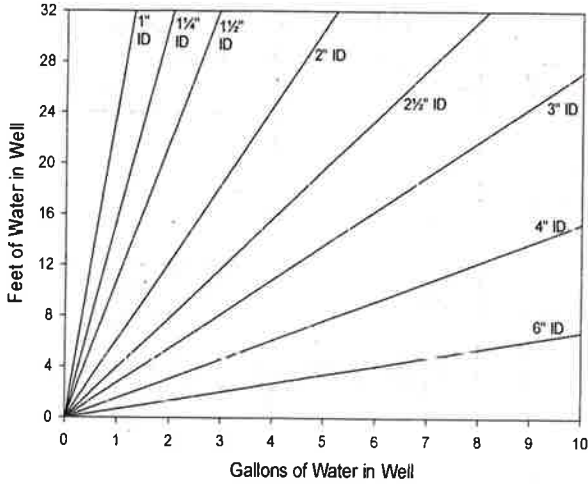
### 3. SAMPLE COLLECTION: Method: Geotech bladder pump with drop tube assembly

Sample ID	Container Type	No. of Containers	Preservation	Analysis Req.	Time
RE10301-610-093015	40-mL vial	3	HCl	VOCs	1315
	1-L amber	2	none	1,4-Dioxane	1315

Comments: Micro 90 dacon transducer

Signature: Paul Kozicki Date: 9/30/15

# Purge Volume Calculation



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

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

Well ID: RE10301

(continued from front)

Time (24 hr)	Volume		Temp (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
	Removed (Liters)										
0955			15.44	4.22	0.078	4.13	350.7	0.20	700	40.25	
1300	8.5		15.39	4.25	0.080	4.35	351.4				
1305			15.58	4.32	0.082	4.33	348.4	0.01		40.26	
1310			15.52	4.41	0.082	4.42	344.6				
1315											Sample



RESOLUTION CONSULTANTS

Well ID: RE10302

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 9/30/15 Time: Start 1200 am/pm  
 Project No: 60266526 Finish 1405 am/pm  
 Site Location: Avoca  
 Weather Conds: 750F overcast - showers Collector(s): JC

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

a. Total Well Length 673 c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material  
 4-inch PVC  
 b. Water Table Depth 39.98 d. Calculated System Volume (see back) 13.5 gal surge

### 2. WELL PURGE DATA

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

c. Field Testing Equipment used: Make Model Serial Number  
YSI 556 MPS 24698  
Scientific Micro TPW 201205051

+2 hrs!  
↓

Time (24hr)	Volume Removed (Liters)	Temp. (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
1030	-	22.84	5.21	0.036	11.62	342.9	-	500	39.61	clear/none
1035	-	17.76	4.86	0.037	12.08	358.2	1.30	500	39.61	"
1040	-	17.10	4.51	0.035	4.85	267.2	-	750	39.60	"
1045	-	17.09	4.47	0.034	5.33	367.9	0.07	750	39.61	"
1050	-	16.94	4.44	0.031	5.97	371.3	-	750	39.61	"
1100	5gal	16.88	4.31	0.030	6.95	381.6	0.00	750	39.62	"
1105	-	16.90	4.29	0.030	7.01	384.5	-	750	39.62	"

d. Acceptance criteria pass/fail Yes No N/A (continued on back)  
 Has required volume been removed     
 Has required turbidity been reached     
 Have parameters stabilized     
 If no or N/A - Explain below.

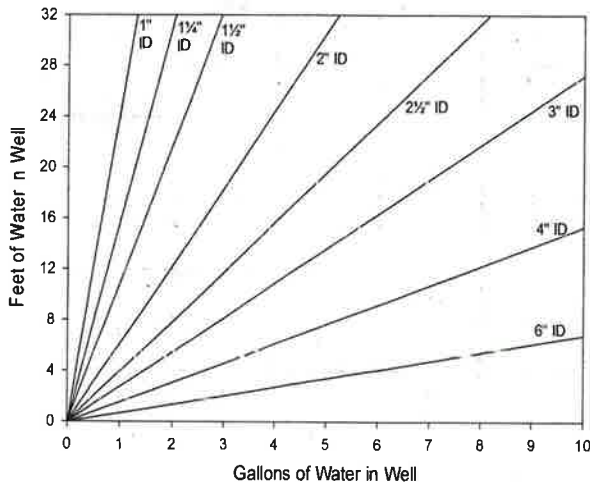
### 3. SAMPLE COLLECTION: Method: Geotech bladder pump with drop tube assembly

Sample ID	Container Type	No. of Containers	Preservation	Analysis Req.	Time
RE10302-GW-093015	40-mL vial	3	HCl	VOCs	1345
RE10302-GW-093015	1-L amber	2	none	1,4-Dioxane	1345

Comments Micro 90 decm

Signature [Signature] Date 9/30/15

# Purge Volume Calculation



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

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

## Well ID:

(continued from front)

12 hrs  
 ↓

Time (24 hr)	Volume Removed (Liters)	Temp (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
1110	—	16.84	4.24	0.030	6.83	388.4	0.00	750	39.62	clear
1115	—	16.84	4.19	0.029	6.78	392.6	0.00	750	39.62	"
1120	10 gal	16.86	4.21	0.029	6.74	390.9	0.00	750	39.62	"
1130	—	16.89	4.25	0.030	6.64	390.2	0.00	750	39.62	"
1135	—	16.79	4.22	0.030	7.12	394.7	0.00	750	39.62	"
1140	13.5 gal	16.80	4.26	0.030	7.17	392.9	0.00	750	39.62	"



RESOLUTION CONSULTANTS

Well ID: RE103D3

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 9/30/15 Time: Start 1200 am/pm  
 Project No: 60266526 Finish 1400 am/pm  
 Site Location: Avoca  
 Weather Conds: Overcast drizzle 75° Collector(s):

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

a. Total Well Length 735 c. Length of Water Column 15 (a-b) Casing Diameter/Material 4-inch PVC  
 b. Water Table Depth 40.82 d. Calculated System Volume (see back) 9.8 g

### 2. WELL PURGE DATA

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

c. Field Testing Equipment used: Make Model Serial Number  
 YSI 556 MPS RFW22120

Time (24hr)	Volume Removed (Liters)	Temp. (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
12:20										
12:25		16.81	4.92	0.027	11.69	281.0				
12:30		16.24	3.89	0.026	8.03	336.0	2.85	450	40.78	clear NO
12:35	1.5g	16.15	4.17	0.026	5.09	320.0				
12:40	2 g	16.19	4.35	0.025	3.96	311.4				
12:45	2.5g	16.09	4.45	0.025	3.61	303.9				
12:50	3 g	16.03	4.37	0.025	3.80	307.2	0.10			

d. Acceptance criteria pass/fail Yes No N/A (continued on back)  
 Has required volume been removed     
 Has required turbidity been reached     
 Have parameters stabilized     
 If no or N/A - Explain below.

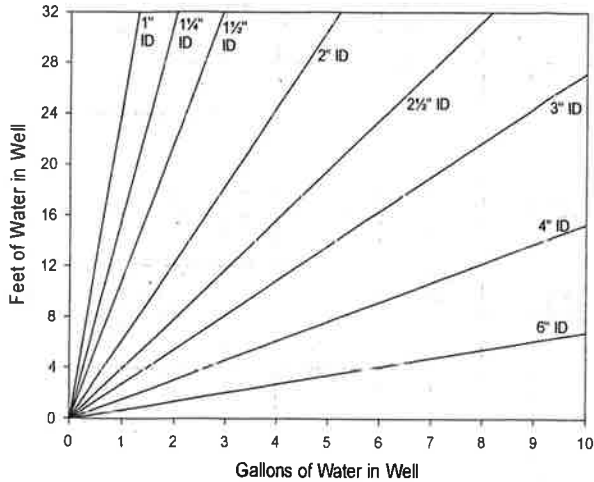
### 3. SAMPLE COLLECTION: Method: Geotech bladder pump with drop tube assembly

Sample ID	Container Type	No. of Containers	Preservation	Analysis Req.	Time
RE103D3-GW-093015	40-mL vial	3	HCl	VOCs	1335
	1-L amber	2	none	1,4-Dioxane	

Comments: Micro 90 decan

Signature: [Signature] Date: 9-30-15

Purge Volume Calculation



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

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

Well ID:

(continued from front)

Time (24 hr)	Volume Removed (Liters)	Temp (°C)	pH	Spec. Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
12:55		16.05	4.36	0.025	3.98	307.9	0.0	500		
13:00	5g	15.98	4.37	0.025	4.01	308.6			40.78	
13:05	5g	16.14	4.39	0.025	4.13	310.4				
13:10	7g	16.12	4.41	0.025	4.10	310.9	0.0			
13:18	8g	16.05	4.41	0.025	4.09	311.6		500	40.62	
13:24		16.03	4.41	0.026	4.13	311.9				
13:30	10g	16.00	4.42	0.026	4.15	312.1				

## **Appendix B**

### **Analytical Data Validation – Resolution Consultants**



**DATA VALIDATION REPORT**

Project:	Regional Groundwater Investigation — NWIRP Bethpage		
Laboratory:	Katahdin Analytical		
Sample Delivery Group:	BETHPAGE-2		
Analyses/Method:	Volatile Organic Compounds by U.S. EPA SW-846 Method 8260C 1,4-Dioxane by U.S. EPA SW-846 Method 8270D via Selective Ion Monitoring (SIM)		
Validation Level:	3		
Project Number:	0888812477.SA.DV		
Prepared by:	Dana Miller/Resolution Consultants	Completed on: 11/30/2015	
Reviewed by:	Tina Clemmey/Resolution Consultants	File Name: BETHPAGE2_8260C_8270D	

**SUMMARY**

This report summarizes data review findings for samples listed below, collected by Resolution Consultants from the Regional Groundwater Investigation — NWIRP Bethpage Site on 25 and 30 September 2015 in accordance with the following Sampling and Analysis Plans:

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

Sample ID	Lab ID	Matrix/Sample Type	Analysis
RE104D1-GW-092515	SI7583-1	Groundwater	8260C/8270D_SIM
RE108D2-GW-092815	SI7583-10	Groundwater	8260C/8270D_SIM
TB01-092815	SI7583-11	Trip Blank	8260C
TB02-092815	SI7583-18RA	Trip Blank	8260C
RE104D2-GW-092515	SI7583-2	Groundwater	8260C/8270D_SIM
RE104D3-GW-092515	SI7583-3	Groundwater	8260C/8270D_SIM
TT309S-GW-092515	SI7583-4	Groundwater	8260C/8270D_SIM
TT309I-GW-092515	SI7583-5	Groundwater	8260C/8270D_SIM
TT309D-GW-092515	SI7583-6	Groundwater	8260C/8270D_SIM
RE105D1-GW-092815	SI7583-7	Groundwater	8260C/8270D_SIM

Sample ID	Lab ID	Matrix/Sample Type	Analysis
RE105D2-GW-092815	SI7583-8	Groundwater	8260C/8270D_SIM
RE108D1-GW-092815	SI7583-9	Groundwater	8260C/8270D_SIM
DUPLICATE-GW-092915	SI7681-10	Duplicate of TT101D-GW-092915	8260C/8270D_SIM
RE122D1-GW-093015	SI7681-11	Groundwater	8260C/8270D_SIM
RE122D2-GW-093015	SI7681-12	Groundwater	8260C/8270D_SIM
RE122D3-GW-093015	SI7681-13	Groundwater	8260C/8270D_SIM
RE103D1-GW-093015	SI7681-14	Groundwater	8260C/8270D_SIM
RE103D2-GW-093015	SI7681-15	Groundwater	8260C/8270D_SIM
RE103D3-GW-093015	SI7681-16	Groundwater	8260C/8270D_SIM
TB01-093015	SI7681-17	Trip Blank	8260C
RE123D1-GW-092915	SI7681-1RA	Groundwater	8260C/8270D_SIM
RE123D2-GW-092915	SI7681-2RA	Groundwater	8260C/8270D_SIM
RE123D3-GW-092915	SI7681-3RA	Groundwater	8260C/8270D_SIM
RE120D1-GW-092915	SI7681-4	Groundwater	8260C/8270D_SIM
RE120D2-GW-092915	SI7681-5	Groundwater	8260C/8270D_SIM
RE120D3-GW-092915	SI7681-6	Groundwater	8260C/8270D_SIM
TT101D-GW-092915	SI7681-7	Groundwater	8260C/8270D_SIM
TT101D1-GW-092915	SI7681-8	Groundwater	8260C/8270D_SIM
TT101D2-GW-092915	SI7681-9	Groundwater	8260C/8270D_SIM

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

## REVIEW ELEMENTS

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

- ✓ Data completeness (chain-of-custody)/sample integrity
- ✓ Holding times and sample preservation

- ✓ Gas chromatography/Mass spectrometer performance checks
- X Initial calibration verification (ICV)/continuing calibration verification (CCV)
- X Laboratory blanks/trip blanks
- X Surrogate spike recoveries
- X Matrix spike and/or matrix spike duplicate results
- ✓ Laboratory control sample/laboratory control sample duplicate results
- ✓ Field duplicates
- ✓ Internal standards
- ✓ Sample results/reporting issues

The symbol (✓) indicates that no validation qualifiers were applied based on this parameter. Acceptable data parameters for which all criteria were met and no qualification was performed and non-conformance or other issues that were noted during validation, but did not result in qualification of data are not discussed further. The symbol (X) indicates that a QC non-conformance resulted in the qualification of data. Any QC non-conformance that resulted in the qualification of data is discussed below.

## **RESULTS**

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

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

- The initial calibration percent relative standard deviation, correlation coefficient/coefficient of determination, and/or response factor method acceptance criteria were met
- The ICV standard percent recovery acceptance criteria were met
- The CCV method percent difference or percent drift and response factor acceptance criteria were met
- The retention time method acceptance criteria were met

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

**ICAL Linearity Non-conformance:**

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

**Notes:**

%RSD = Relative standard deviation  
 J = Estimated  
 UJ = Undetected and estimated

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

**ICV Recovery Non-conformance:**

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

**Notes:**

J = Estimated  
 UJ = Undetected and estimated

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

**CCV Linearity Non-conformance:**

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

**Notes:**

J = Estimated  
 UJ = Undetected and estimated

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

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

Laboratory blanks, equipment blanks, field blanks, and trip blanks were analyzed with samples to assess contamination imparted by sample preparation and/or analysis. All results associated with a particular blank were evaluated to determine whether there was an inherent variability in the data, or if a problem was an isolated occurrence that did not affect the data. Samples were flagged in accordance with *Functional Guidelines* (shown below) where detections were not believed to be site-related.

**Blank Non-conformance Charts:**

<i>For common lab contaminants (methylene chloride, acetone, 2-butanone):</i>			
Blank type	Blank result	Sample result	Action for samples
Method, Storage, Trip, Field, or Equipment	Detects	Not detected	No qualification
	≤ 2x LOQ	< 2x LOQ	Report sample LOQ value with a U
		≥ 2x LOQ and ≤ 4x the LOQ	Report the sample result with a U**
		≥ 4x the LOQ	No qualifications
	> 2x LOQ	< LOD	Report sample LOD value with a U**
		≥ LOD and < 2x LOQ	Report sample LOQ value with a U
		≥ 2x LOQ and < blank contamination	Report the blank result with a U or reject the sample result as unusable R
		≥ 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.**
<b>**Based on Resolution Consultants professional judgment</b>			

<i>For all other compounds:</i>			
Blank type	Blank result	Sample result	Action for samples
Method, Storage, Trip, Field, or Equipment	Detects	Not detected	No qualification
	< 2x LOQ	< 2x LOQ	Report sample LOQ value with a U
		≥ 2x LOQ	Use professional judgment
	> 2x LOQ	< 2x LOQ	Report sample LOQ value with a U
		≥ 2x LOQ and < blank contamination	Report the blank result with a U or reject the sample result as unusable R
		≥ 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.
	= 2x LOQ	< 2x LOQ	Report sample LOQ value with a U
		≥ 2x LOQ	Use professional judgment
	Gross contamination	Detects	Qualify results as unusable R

**Notes:**

LOQ	=	Limit of quantitation
LOD	=	Limit of detection
U	=	Undetected
R	=	Rejected

Lab blank and trip blank non-conformances are summarized in Attachment A in Table's A-4, and A-5.

**Surrogate Spike Recoveries**

Surrogates provide information needed to assess the accuracy of analyses. Known amounts of surrogate compounds, or compounds which are not likely to be found in the actual samples, are added to each organic sample to check for accuracy. If surrogate percent recoveries (%Rs) are close

to the known concentrations, the reported target compound concentrations are assumed to be accurate. Data qualification on the basis of surrogate recovery was as follows:

**Surrogate Recovery Non-conformance Chart:**

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

**Notes:**

- %R = Percent recovery
- J = Estimated
- UJ = Undetected and estimated

Surrogate recovery non-conformance is summarized in Attachment A in Table A-6.

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

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

**MS/MSD Non-conformances Chart:**

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

**Notes:**

- %R = Percent recovery
- RPD = Relative percent difference
- J = Estimated
- UJ = Undetected and estimated

MS/MSD non-conformances are summarized in Attachment A in Table A-7.

### **Qualifications Actions**

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

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

### **ATTACHMENTS**

Attachment A: Non-Conformance Summary Tables

Attachment B: Qualifier Codes and Explanations

Attachment C: Reason Codes and Explanations

Attachment D: Final Results after Data Review

**Attachment A**  
**Non-Conformance Summary Table**



Table A-1 Initial Calibration Non-Conformance					
Method	Analyte	%RSD	Limit	Associated Samples	Qualifier
8260C	Chloromethane	15.21596	<15%	SI7583-1 through SI7583-11, SI7583-18RA, SI7583-10DL, SI7583-8DL	Detects: J Non-detects: UJ
8260C	Bromomethane	15.30744	<15%	SI7583-1 through SI7583-11, SI7583-18RA, SI7583-10DL, SI7583-8DL	Detects: J Non-detects: UJ
8260C	Chloroethane	22.54919	<15%	SI7583-1 through SI7583-11, SI7583-18RA, SI7583-10DL, SI7583-8DL	Detects: J Non-detects: UJ
8260C	Acetone	16.25899	<15%	SI7583-1 through SI7583-11, SI7583-18RA, SI7583-10DL, SI7583-8DL	Detects: J Non-detects: UJ
8260C	Methyl cyclohexane	15.66496	<15%	SI7583-1 through SI7583-11, SI7583-18RA, SI7583-10DL, SI7583-8DL	Detects: J Non-detects: UJ
8260C	2-Hexanone	15.46497	<15%	SI7681-14 through SI7681-17, SI7681-1RA through SI7681-3RA, SI7681-9DL, SI7681-11DL, SI7681-12DL,	Detects: J Non-detects: UJ
8260C	O-Xylene	15.32272	<15%	SI7681-14 through SI7681-17, SI7681-1RA through SI7681-3RA, SI7681-9DL, SI7681-11DL, SI7681-12DL,	Detects: J Non-detects: UJ
8260C	1,2-Dibromo-3-chloropropane	29.69098	<15%	SI7681-4 through SI7681-13, SI7681-14DL through SI7681-16DL	Detects: J Non-detects: UJ

**Notes:**

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

Table A-2 Initial Calibration Verification Non-Conformance						
Method	Analyte	ICV ID	%R	Limit	Associated Samples	Qualifier
8260C	Dichlorodifluoromethane	WG171044-7	67.47	80-120	S17583-1 through S17583-11, S17583-18RA, S17583-10DL, S17583-8DL	Detects: J Non-detects: UJ
8260C	Trichlorofluoromethane	WG171044-7	79.95	80-120	S17583-1 through S17583-11, S17583-18RA, S17583-10DL, S17583-8DL	Detects: J Non-detects: UJ
8260C	Carbon Disulfide	WG171044-7	79.65	80-120	S17583-1 through S17583-11, S17583-18RA, S17583-10DL, S17583-8DL	Detects: J Non-detects: UJ
8260C	Acetone	WG171044-7	138.88	80-120	S17583-1 through S17583-11, S17583-18RA, S17583-10DL, S17583-8DL	Detects: J Non-detects: UJ
8260C	2-Butanone	WG171044-7	159.36	80-120	S17583-1 through S17583-11, S17583-18RA, S17583-10DL, S17583-8DL	Detects: J Non-detects: UJ
8260C	4-methyl-2-pentanone	WG171044-7	146.19	80-120	S17583-1 through S17583-11, S17583-18RA, S17583-10DL, S17583-8DL	Detects: J Non-detects: UJ
8260C	2-Hexanone	WG171044-7	153.44	80-120	S17583-1 through S17583-11, S17583-18RA, S17583-10DL, S17583-8DL	Detects: J Non-detects: UJ
8260C	Carbon Disulfide	WG171658-7	79.48	80-120	S17681-14 through S17681-17, S17681-1RA through S17681-3RA, S17681-9DL, S17681-11DL, S17681-12DL,	Detects: J Non-detects: UJ
8260C	Acetone	WG171658-7	134.34	80-120	S17681-14 through S17681-17, S17681-1RA through S17681-3RA, S17681-9DL, S17681-11DL, S17681-12DL,	Detects: J Non-detects: UJ
8260C	2-Butanone	WG171658-7	143.56	80-120	S17681-14 through S17681-17, S17681-1RA through S17681-3RA, S17681-9DL, S17681-11DL, S17681-12DL,	Detects: J Non-detects: UJ
8260C	4-methyl-2-pentanone	WG171658-7	150.32	80-120	S17681-14 through S17681-17, S17681-1RA through S17681-3RA, S17681-9DL, S17681-11DL, S17681-12DL,	Detects: J Non-detects: UJ
8260C	2-Hexanone	WG171658-7	156.23	80-120	S17681-14 through S17681-17, S17681-1RA through S17681-3RA, S17681-9DL, S17681-11DL, S17681-12DL,	Detects: J Non-detects: UJ
8260C	Dichlorodifluoromethane	WG171660-7	123.49	80-120	S17681-4 through S17681-13, S17681-14DL through S17681-16DL	Detects: J Non-detects: UJ

**Notes:**

ICV ID = Initial calibration verification identification  
ID = Identification  
%R = Percent recovery  
UJ = Non-detect estimated value  
J = Estimated value

Table A-3 Continuing Calibration Verification Non-Conformance					
Lab ID /Calibration ID	Analyte	%D	%D Limit	Associated Samples	Qualifier
WG171352-4 / C4922.D	Tetrachloroethene	34.59924	+/- 20	SI7583-1, SI7583-2, SI7583-3, SI7583-4, SI7583-5, SI7583-6, SI7583-7, SI7583-8, SI7583-9, SI7583-10, and SI7583-11	Detects: J Non-detects: UJ
WG171374-4 / C4942.D	Methyl cyclohexane	21.55731	+/- 20	SI7583-18RA, SI7583-10DL, SI7583-8DL, SI7681-17, SI7681-14, SI7681-15, and SI7681-16	Detects: J Non-detects: UJ
WG171659-4 / C5030.D	Tetrachloroethene	-29.10536	+/- 20	SI7681-1RA, SI7681-2RA, SI7681-3RA, SI7681-4DL, SI7681-5DL, SI7681-8RA, SI7681-9DL, SI7681-11DL, SI7681-12DL	Detects: J Non-detects: UJ
WG171660-4 / P3089.D	Dichlorodifluoromethane	41.50045	+/- 20	SI7681-4, SI7681-5, SI7681-6, SI7681-7, SI7681-8, SI7681-9, SI7681-10, SI7681-11, SI7681-12, SI7681-13	Detects: J Non-detects: UJ
WG171601-4 / P3113.D	Dichlorodifluoromethane	40.79053	+/- 20	SI7681-14DL, SI7681-15DL, SI7681-16DL	Detects: J Non-detects: UJ
WG171601-4 / P3113.D	Chloromethane	22.86899	+/- 20	SI7681-14DL, SI7681-15DL, SI7681-16DL	Detects: J Non-detects: UJ
WG171601-4 / P3113.D	Tetrachloroethene	-30.50148	+/- 20	SI7681-14DL, SI7681-15DL, SI7681-16DL	Detects: J Non-detects: UJ

**Notes:**

ID = Identification  
 %D = Percent difference  
 UJ = Non-detect estimated value  
 J = Detected estimated value

Table A-4 Lab Blank Non-Conformance (Micrograms per liter)					
Blank ID	Analyte	Blank Result	LOQ	Associated Sample	Qualifier
WG171658-9	Carbon Disulfide	0.41	1	TB01-093015	U

**Notes:**

ID = Identification  
 LOQ = Limit of quantitation  
 U = Detected analyte qualified as non-detect due to sample result being less than 2 times the LOQ.

Table A-5 Field Blank Non-Conformance (Micrograms per liter)					
Blank ID	Analyte	Blank Result	LOQ	Associated Sample	Qualifier
TB01-092815	Chloromethane	0.46	2	TT309S-GW-092515	U
TB01-093015	Chloromethane	0.82	2	RE123D2-GW-092915	U
TB01-093015	Chloromethane	0.82	2	RE123D3-GW-092915	U

**Notes:**

ID = Identification  
 LOQ = Limit of quantitation  
 U = Detected analyte qualified as non-detect due to sample result being less than 2 times the LOQ.

Table A-6 Surrogate Non-Conformance					
Method	Surrogate	%R	Limits	Associated Sample	Qualifier
8260C	1,2-Dichloroethane-d4	121	70-120	RE103D3-GW-093015	Detects: J
8260C	1,2-Dichloroethane-d4	122	70-120	RE120D1-GW-092915	Detects: J
8260C	1,2-Dichloroethane-d4	124	70-120	RE120D2-GW-092915	Detects: J
8260C	Dibromofluoromethane	81.3	85-115	RE122D1-GW-093015	Detects: J and Non-detects: UJ
8260C	1,2-Dichloroethane-d4	124	70-120	RE122D1-GW-093015	Detects: J
8260C	Dibromofluoromethane	83.4	85-115	RE122D2-GW-093015	Detects: J and Non-detects: UJ
8260C	1,2-Dichloroethane-d4	122	70-120	RE122D2-GW-093015	Detects: J
8260C	Dibromofluoromethane	84.7	85-115	RE122D3-GW-093015	Detects: J and Non-detects: UJ
8260C	Dibromofluoromethane	80.4	85-115	TT101D1-GW-092915	Detects: J and Non-detects: UJ
8260C	1,2-Dichloroethane-d4	122	70-120	TT101D1-GW-092915	Detects: J
8260C	Dibromofluoromethane	84.5	85-115	TT101D2-GW-092915	Detects: J and Non-detects: UJ
8260C	1,2-Dichloroethane-d4	121	70-120	TT101D2-GW-092915	Detects: J

**Notes:**

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

Table A-7 Matrix Spike/Matrix Spike Duplicate Non-Conformance (Micrograms per liter)							
Spiked Sample	Analyte	Sample Result	Spike Added	MS %R	MSD %R	%R Limits	Qualifier
TT101D2-GW-092915	Trichloroethene	640	50.0	<b>154*</b>	<b>252*</b>	70-125	J

**Notes:**

MS = Matrix spike  
 MSD = Matrix spike duplicate  
 %R = Percent recovery  
**Bold\*** = Percent recovery less than lower control limit  
 J = Detected analyte in associated sample qualified estimated "J" because %R is greater than control limit in associated sample.

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

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

**Attachment C**  
**Reason Codes and Explanations**

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



**Attachment D**  
**Final Results after Data Review**

Sample Delivery Group				BETHPAGE-2		
Lab ID				SI7583-1		
Sample ID				RE104D1-GW-092515		
Sample Date				9/25/2015		
Sample Type				Groundwater		
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	U	
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	U	
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	4.6		
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.5	U	
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	0.5	U	
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	0.87	J	
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	U	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	U	
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	U	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	U	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	1.2	J	
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	U	
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	U	
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	U	
8260C	2-BUTANONE	78-93-3	UG_L	2.5	UJ	c
8260C	2-HEXANONE	591-78-6	UG_L	2.5	UJ	c
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	UJ	c
8260C	ACETONE	67-64-1	UG_L	2.5	UJ	c
8260C	BENZENE	71-43-2	UG_L	0.5	U	
8260C	BROMODICHLOROMETHANE	75-27-4	UG_L	0.5	U	
8260C	BROMOFORM	75-25-2	UG_L	0.5	U	
8260C	BROMOMETHANE	74-83-9	UG_L	1	UJ	c
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	UJ	c
8260C	CARBON TETRACHLORIDE	56-23-5	UG_L	0.5	U	
8260C	CHLOROBENZENE	108-90-7	UG_L	0.5	U	
8260C	CHLOROETHANE	75-00-3	UG_L	1	UJ	c
8260C	CHLOROFORM	67-66-3	UG_L	0.5	U	
8260C	CHLOROMETHANE	74-87-3	UG_L	1	UJ	c
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	1.2		
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG_L	0.5	U	
8260C	CYCLOHEXANE	110-82-7	UG_L	0.5	U	
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG_L	0.5	U	
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	0.49	J	c
8260C	ETHYLBENZENE	100-41-4	UG_L	0.5	U	
8260C	ISOPROPYLBENZENE	98-82-8	UG_L	0.5	U	
8260C	M- AND P-XYLENE	108-38-3/106-42	UG_L	1	U	
8260C	METHYL ACETATE	79-20-9	UG_L	0.75	U	
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	UJ	c
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	U	
8260C	METHYLENE CHLORIDE	75-09-2	UG_L	2.5	U	
8260C	O-XYLENE	95-47-6	UG_L	0.5	U	
8260C	STYRENE	100-42-5	UG_L	0.5	U	
8260C	TETRACHLOROETHENE	127-18-4	UG_L	2.3	J	c
8260C	TOLUENE	108-88-3	UG_L	0.5	U	
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	U	
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	U	
8260C	TRICHLOROETHENE	79-01-6	UG_L	110		
8260C	TRICHLOROFLUOROMETHANE	75-69-4	UG_L	1	UJ	c
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	U	
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	U	
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	12		

Sample Delivery Group				BETHPAGE-2		
Lab ID				SI7583-10		
Sample ID				RE108D2-GW-092815		
Sample Date				9/28/2015		
Sample Type				Groundwater		
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	U	
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	U	
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	8.3		
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	1.5		
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	5.3		
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	7.6		
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	U	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	U	
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	U	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	U	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	8.6		
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	U	
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	U	
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	U	
8260C	2-BUTANONE	78-93-3	UG_L	2.5	UJ	c
8260C	2-HEXANONE	591-78-6	UG_L	2.5	UJ	c
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	UJ	c
8260C	ACETONE	67-64-1	UG_L	2.5	UJ	c
8260C	BENZENE	71-43-2	UG_L	0.5	U	
8260C	BROMODICHLOROMETHANE	75-27-4	UG_L	0.5	U	
8260C	BROMOFORM	75-25-2	UG_L	0.5	U	
8260C	BROMOMETHANE	74-83-9	UG_L	1	UJ	c
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	UJ	c
8260C	CARBON TETRACHLORIDE	56-23-5	UG_L	1.6		
8260C	CHLOROBENZENE	108-90-7	UG_L	0.5	U	
8260C	CHLOROETHANE	75-00-3	UG_L	1	UJ	c
8260C	CHLOROFORM	67-66-3	UG_L	3.7		
8260C	CHLOROMETHANE	74-87-3	UG_L	1	UJ	c
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	8.6		
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG_L	0.5	U	
8260C	CYCLOHEXANE	110-82-7	UG_L	0.5	U	
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG_L	0.5	U	
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	1	UJ	c
8260C	ETHYLBENZENE	100-41-4	UG_L	0.5	U	
8260C	ISOPROPYLBENZENE	98-82-8	UG_L	0.5	U	
8260C	M- AND P-XYLENE	108-38-3/106-42	UG_L	1	U	
8260C	METHYL ACETATE	79-20-9	UG_L	0.75	U	
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	UJ	c
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	U	
8260C	METHYLENE CHLORIDE	75-09-2	UG_L	2.5	U	
8260C	O-XYLENE	95-47-6	UG_L	0.5	U	
8260C	STYRENE	100-42-5	UG_L	0.5	U	
8260C	TETRACHLOROETHENE	127-18-4	UG_L	2.5	J	c
8260C	TOLUENE	108-88-3	UG_L	0.5	U	
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	U	
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	U	
8260C	TRICHLOROETHENE	79-01-6	UG_L	3400		
8260C	TRICHLOROFUOROMETHANE	75-69-4	UG_L	1	UJ	c
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	U	
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	U	
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	7.5		

Sample Delivery Group				BETHPAGE-2		
Lab ID				SI7583-11		
Sample ID				TB01-092815		
Sample Date				9/28/2015		
Sample Type				Trip Blank		
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	U	
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	U	
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	0.5	U	
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.5	U	
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	0.5	U	
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	0.5	U	
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	U	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	U	
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	U	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	U	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	1	U	
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	U	
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	U	
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	U	
8260C	2-BUTANONE	78-93-3	UG_L	2.5	UJ	c
8260C	2-HEXANONE	591-78-6	UG_L	2.5	UJ	c
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	UJ	c
8260C	ACETONE	67-64-1	UG_L	2.5	UJ	c
8260C	BENZENE	71-43-2	UG_L	0.5	U	
8260C	BROMODICHLOROMETHANE	75-27-4	UG_L	0.5	U	
8260C	BROMOFORM	75-25-2	UG_L	0.5	U	
8260C	BROMOMETHANE	74-83-9	UG_L	1	UJ	c
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	UJ	c
8260C	CARBON TETRACHLORIDE	56-23-5	UG_L	0.5	U	
8260C	CHLOROBENZENE	108-90-7	UG_L	0.5	U	
8260C	CHLOROETHANE	75-00-3	UG_L	1	UJ	c
8260C	CHLOROFORM	67-66-3	UG_L	0.5	U	
8260C	CHLOROMETHANE	74-87-3	UG_L	0.46	J	c
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	0.5	U	
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG_L	0.5	U	
8260C	CYCLOHEXANE	110-82-7	UG_L	0.5	U	
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG_L	0.5	U	
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	1	UJ	c
8260C	ETHYLBENZENE	100-41-4	UG_L	0.5	U	
8260C	ISOPROPYLBENZENE	98-82-8	UG_L	0.5	U	
8260C	M- AND P-XYLENE	108-38-3/106-42	UG_L	1	U	
8260C	METHYL ACETATE	79-20-9	UG_L	0.75	U	
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	UJ	c
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	U	
8260C	METHYLENE CHLORIDE	75-09-2	UG_L	1.8	J	
8260C	O-XYLENE	95-47-6	UG_L	0.5	U	
8260C	STYRENE	100-42-5	UG_L	0.5	U	
8260C	TETRACHLOROETHENE	127-18-4	UG_L	0.5	UJ	c
8260C	TOLUENE	108-88-3	UG_L	0.5	U	
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	U	
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	U	
8260C	TRICHLOROETHENE	79-01-6	UG_L	0.5	U	
8260C	TRICHLOROFUOROMETHANE	75-69-4	UG_L	1	UJ	c
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	U	
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	U	
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	NA		

Sample Delivery Group				BETHPAGE-2		
Lab ID				SI7583-18RA		
Sample ID				TB02-092815		
Sample Date				9/28/2015		
Sample Type				Trip Blank		
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	U	
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	U	
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	0.5	U	
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.5	U	
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	0.5	U	
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	0.5	U	
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	U	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	U	
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	U	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	U	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	1	U	
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	U	
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	U	
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	U	
8260C	2-BUTANONE	78-93-3	UG_L	2.5	UJ	c
8260C	2-HEXANONE	591-78-6	UG_L	2.5	UJ	c
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	UJ	c
8260C	ACETONE	67-64-1	UG_L	2.5	UJ	c
8260C	BENZENE	71-43-2	UG_L	0.5	U	
8260C	BROMODICHLOROMETHANE	75-27-4	UG_L	0.5	U	
8260C	BROMOFORM	75-25-2	UG_L	0.5	U	
8260C	BROMOMETHANE	74-83-9	UG_L	1	UJ	c
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	UJ	c
8260C	CARBON TETRACHLORIDE	56-23-5	UG_L	0.5	U	
8260C	CHLOROBENZENE	108-90-7	UG_L	0.5	U	
8260C	CHLOROETHANE	75-00-3	UG_L	1	UJ	c
8260C	CHLOROFORM	67-66-3	UG_L	0.5	U	
8260C	CHLOROMETHANE	74-87-3	UG_L	0.39	J	c
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	0.5	U	
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG_L	0.5	U	
8260C	CYCLOHEXANE	110-82-7	UG_L	0.5	U	
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG_L	0.5	U	
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	1	UJ	c
8260C	ETHYLBENZENE	100-41-4	UG_L	0.5	U	
8260C	ISOPROPYLBENZENE	98-82-8	UG_L	0.5	U	
8260C	M- AND P-XYLENE	108-38-3/106-42	UG_L	1	U	
8260C	METHYL ACETATE	79-20-9	UG_L	0.75	U	
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	UJ	c
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	U	
8260C	METHYLENE CHLORIDE	75-09-2	UG_L	2	J	
8260C	O-XYLENE	95-47-6	UG_L	0.5	U	
8260C	STYRENE	100-42-5	UG_L	0.5	U	
8260C	TETRACHLOROETHENE	127-18-4	UG_L	0.5	U	
8260C	TOLUENE	108-88-3	UG_L	0.5	U	
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	U	
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	U	
8260C	TRICHLOROETHENE	79-01-6	UG_L	0.5	U	
8260C	TRICHLOROFUOROMETHANE	75-69-4	UG_L	1	UJ	c
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	U	
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	U	
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	NA		

Sample Delivery Group				BETHPAGE-2		
Lab ID				SI7583-2		
Sample ID				RE104D2-GW-092515		
Sample Date				9/25/2015		
Sample Type				Groundwater		
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	U	
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	U	
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	0.5	U	
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.5	U	
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	0.5	U	
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	0.5	U	
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	U	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	U	
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	U	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	U	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	2		
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	U	
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	U	
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	U	
8260C	2-BUTANONE	78-93-3	UG_L	2.5	UJ	c
8260C	2-HEXANONE	591-78-6	UG_L	2.5	UJ	c
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	UJ	c
8260C	ACETONE	67-64-1	UG_L	2.5	UJ	c
8260C	BENZENE	71-43-2	UG_L	0.5	U	
8260C	BROMODICHLOROMETHANE	75-27-4	UG_L	0.5	U	
8260C	BROMOFORM	75-25-2	UG_L	0.5	U	
8260C	BROMOMETHANE	74-83-9	UG_L	1	UJ	c
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	UJ	c
8260C	CARBON TETRACHLORIDE	56-23-5	UG_L	0.5	U	
8260C	CHLOROBENZENE	108-90-7	UG_L	0.5	U	
8260C	CHLOROETHANE	75-00-3	UG_L	1	UJ	c
8260C	CHLOROFORM	67-66-3	UG_L	0.53	J	
8260C	CHLOROMETHANE	74-87-3	UG_L	1	UJ	c
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	2		
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG_L	0.5	U	
8260C	CYCLOHEXANE	110-82-7	UG_L	0.5	U	
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG_L	0.5	U	
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	1	UJ	c
8260C	ETHYLBENZENE	100-41-4	UG_L	0.5	U	
8260C	ISOPROPYLBENZENE	98-82-8	UG_L	0.5	U	
8260C	M- AND P-XYLENE	108-38-3/106-42	UG_L	1	U	
8260C	METHYL ACETATE	79-20-9	UG_L	0.75	U	
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	UJ	c
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	U	
8260C	METHYLENE CHLORIDE	75-09-2	UG_L	2.5	U	
8260C	O-XYLENE	95-47-6	UG_L	0.5	U	
8260C	STYRENE	100-42-5	UG_L	0.5	U	
8260C	TETRACHLOROETHENE	127-18-4	UG_L	0.5	UJ	c
8260C	TOLUENE	108-88-3	UG_L	0.5	U	
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	U	
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	U	
8260C	TRICHLOROETHENE	79-01-6	UG_L	4.2		
8260C	TRICHLOROFUOROMETHANE	75-69-4	UG_L	1	UJ	c
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	U	
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	U	
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	0.13	J	

Sample Delivery Group				BETHPAGE-2		
Lab ID				SI7583-3		
Sample ID				RE104D3-GW-092515		
Sample Date				9/25/2015		
Sample Type				Groundwater		
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	U	
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	U	
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	0.5	U	
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.5	U	
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	0.5	U	
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	0.5	U	
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	U	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	U	
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	U	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	U	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	1	U	
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	U	
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	U	
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	U	
8260C	2-BUTANONE	78-93-3	UG_L	2.5	UJ	c
8260C	2-HEXANONE	591-78-6	UG_L	2.5	UJ	c
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	UJ	c
8260C	ACETONE	67-64-1	UG_L	2.5	UJ	c
8260C	BENZENE	71-43-2	UG_L	0.5	U	
8260C	BROMODICHLOROMETHANE	75-27-4	UG_L	0.5	U	
8260C	BROMOFORM	75-25-2	UG_L	0.5	U	
8260C	BROMOMETHANE	74-83-9	UG_L	1	UJ	c
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	UJ	c
8260C	CARBON TETRACHLORIDE	56-23-5	UG_L	0.5	U	
8260C	CHLOROBENZENE	108-90-7	UG_L	0.5	U	
8260C	CHLOROETHANE	75-00-3	UG_L	1	UJ	c
8260C	CHLOROFORM	67-66-3	UG_L	0.5	U	
8260C	CHLOROMETHANE	74-87-3	UG_L	1	UJ	c
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	0.5	U	
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG_L	0.5	U	
8260C	CYCLOHEXANE	110-82-7	UG_L	0.5	U	
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG_L	0.5	U	
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	1	UJ	c
8260C	ETHYLBENZENE	100-41-4	UG_L	0.5	U	
8260C	ISOPROPYLBENZENE	98-82-8	UG_L	0.5	U	
8260C	M- AND P-XYLENE	108-38-3/106-42	UG_L	1	U	
8260C	METHYL ACETATE	79-20-9	UG_L	0.75	U	
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	UJ	c
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	U	
8260C	METHYLENE CHLORIDE	75-09-2	UG_L	2.5	U	
8260C	O-XYLENE	95-47-6	UG_L	0.5	U	
8260C	STYRENE	100-42-5	UG_L	0.5	U	
8260C	TETRACHLOROETHENE	127-18-4	UG_L	0.5	UJ	c
8260C	TOLUENE	108-88-3	UG_L	0.5	U	
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	U	
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	U	
8260C	TRICHLOROETHENE	79-01-6	UG_L	0.5	U	
8260C	TRICHLOROFUOROMETHANE	75-69-4	UG_L	1	UJ	c
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	U	
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	U	
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	0.17	U	

Sample Delivery Group				BETHPAGE-2		
Lab ID				SI7583-4		
Sample ID				TT309S-GW-092515		
Sample Date				9/25/2015		
Sample Type				Groundwater		
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	U	
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	U	
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	0.5	U	
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.5	U	
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	0.5	U	
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	0.5	U	
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	U	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	U	
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	U	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	U	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	1	U	
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	U	
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	U	
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	U	
8260C	2-BUTANONE	78-93-3	UG_L	2.5	UJ	c
8260C	2-HEXANONE	591-78-6	UG_L	2.5	UJ	c
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	UJ	c
8260C	ACETONE	67-64-1	UG_L	2.5	UJ	c
8260C	BENZENE	71-43-2	UG_L	0.5	U	
8260C	BROMODICHLOROMETHANE	75-27-4	UG_L	0.5	U	
8260C	BROMOFORM	75-25-2	UG_L	0.5	U	
8260C	BROMOMETHANE	74-83-9	UG_L	1	UJ	c
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	UJ	c
8260C	CARBON TETRACHLORIDE	56-23-5	UG_L	0.5	U	
8260C	CHLOROBENZENE	108-90-7	UG_L	0.5	U	
8260C	CHLOROETHANE	75-00-3	UG_L	1	UJ	c
8260C	CHLOROFORM	67-66-3	UG_L	0.5	U	
8260C	CHLOROMETHANE	74-87-3	UG_L	1	UJ	bt,c
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	0.5	U	
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG_L	0.5	U	
8260C	CYCLOHEXANE	110-82-7	UG_L	0.5	U	
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG_L	0.5	U	
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	1	UJ	c
8260C	ETHYLBENZENE	100-41-4	UG_L	0.5	U	
8260C	ISOPROPYLBENZENE	98-82-8	UG_L	0.5	U	
8260C	M- AND P-XYLENE	108-38-3/106-42	UG_L	1	U	
8260C	METHYL ACETATE	79-20-9	UG_L	0.75	U	
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	UJ	c
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	U	
8260C	METHYLENE CHLORIDE	75-09-2	UG_L	2.5	U	
8260C	O-XYLENE	95-47-6	UG_L	0.5	U	
8260C	STYRENE	100-42-5	UG_L	0.5	U	
8260C	TETRACHLOROETHENE	127-18-4	UG_L	0.5	UJ	c
8260C	TOLUENE	108-88-3	UG_L	0.5	U	
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	U	
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	U	
8260C	TRICHLOROETHENE	79-01-6	UG_L	0.89	J	
8260C	TRICHLOROFUOROMETHANE	75-69-4	UG_L	1	UJ	c
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	U	
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	U	
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	0.35		



Sample Delivery Group				BETHPAGE-2		
Lab ID				SI7583-5		
Sample ID				TT309I-GW-092515		
Sample Date				9/25/2015		
Sample Type				Groundwater		
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	U	
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	U	
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	0.5	U	
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.5	U	
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	0.5	U	
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	0.5	U	
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	U	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	U	
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	U	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	U	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	1	U	
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	U	
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	U	
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	U	
8260C	2-BUTANONE	78-93-3	UG_L	2.5	UJ	c
8260C	2-HEXANONE	591-78-6	UG_L	2.5	UJ	c
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	UJ	c
8260C	ACETONE	67-64-1	UG_L	2.5	UJ	c
8260C	BENZENE	71-43-2	UG_L	0.5	U	
8260C	BROMODICHLOROMETHANE	75-27-4	UG_L	0.5	U	
8260C	BROMOFORM	75-25-2	UG_L	0.5	U	
8260C	BROMOMETHANE	74-83-9	UG_L	1	UJ	c
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	UJ	c
8260C	CARBON TETRACHLORIDE	56-23-5	UG_L	0.5	U	
8260C	CHLOROBENZENE	108-90-7	UG_L	0.5	U	
8260C	CHLOROETHANE	75-00-3	UG_L	1	UJ	c
8260C	CHLOROFORM	67-66-3	UG_L	0.5	U	
8260C	CHLOROMETHANE	74-87-3	UG_L	1	UJ	c
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	0.5	U	
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG_L	0.5	U	
8260C	CYCLOHEXANE	110-82-7	UG_L	0.5	U	
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG_L	0.5	U	
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	1	UJ	c
8260C	ETHYLBENZENE	100-41-4	UG_L	0.5	U	
8260C	ISOPROPYLBENZENE	98-82-8	UG_L	0.5	U	
8260C	M- AND P-XYLENE	108-38-3/106-42	UG_L	1	U	
8260C	METHYL ACETATE	79-20-9	UG_L	0.75	U	
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	UJ	c
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	U	
8260C	METHYLENE CHLORIDE	75-09-2	UG_L	2.5	U	
8260C	O-XYLENE	95-47-6	UG_L	0.5	U	
8260C	STYRENE	100-42-5	UG_L	0.5	U	
8260C	TETRACHLOROETHENE	127-18-4	UG_L	0.5	UJ	c
8260C	TOLUENE	108-88-3	UG_L	0.5	U	
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	U	
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	U	
8260C	TRICHLOROETHENE	79-01-6	UG_L	0.5	U	
8260C	TRICHLOROFUOROMETHANE	75-69-4	UG_L	1	UJ	c
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	U	
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	U	
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	0.61		

Sample Delivery Group				BETHPAGE-2		
Lab ID				SI7583-6		
Sample ID				TT309D-GW-092515		
Sample Date				9/25/2015		
Sample Type				Groundwater		
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	U	
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	U	
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	0.5	U	
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.5	U	
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	0.5	U	
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	0.5	U	
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	U	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	U	
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	U	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	U	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	1	U	
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	U	
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	U	
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	U	
8260C	2-BUTANONE	78-93-3	UG_L	2.5	UJ	c
8260C	2-HEXANONE	591-78-6	UG_L	2.5	UJ	c
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	UJ	c
8260C	ACETONE	67-64-1	UG_L	2.5	UJ	c
8260C	BENZENE	71-43-2	UG_L	0.5	U	
8260C	BROMODICHLOROMETHANE	75-27-4	UG_L	0.5	U	
8260C	BROMOFORM	75-25-2	UG_L	0.5	U	
8260C	BROMOMETHANE	74-83-9	UG_L	1	UJ	c
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	UJ	c
8260C	CARBON TETRACHLORIDE	56-23-5	UG_L	0.5	U	
8260C	CHLOROBENZENE	108-90-7	UG_L	0.5	U	
8260C	CHLOROETHANE	75-00-3	UG_L	1	UJ	c
8260C	CHLOROFORM	67-66-3	UG_L	0.5	U	
8260C	CHLOROMETHANE	74-87-3	UG_L	1	UJ	c
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	0.5	U	
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG_L	0.5	U	
8260C	CYCLOHEXANE	110-82-7	UG_L	0.5	U	
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG_L	0.5	U	
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	1	UJ	c
8260C	ETHYLBENZENE	100-41-4	UG_L	0.5	U	
8260C	ISOPROPYLBENZENE	98-82-8	UG_L	0.5	U	
8260C	M- AND P-XYLENE	108-38-3/106-42	UG_L	1	U	
8260C	METHYL ACETATE	79-20-9	UG_L	0.75	U	
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	UJ	c
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	U	
8260C	METHYLENE CHLORIDE	75-09-2	UG_L	2.5	U	
8260C	O-XYLENE	95-47-6	UG_L	0.5	U	
8260C	STYRENE	100-42-5	UG_L	0.5	U	
8260C	TETRACHLOROETHENE	127-18-4	UG_L	0.97	J	c
8260C	TOLUENE	108-88-3	UG_L	0.5	U	
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	U	
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	U	
8260C	TRICHLOROETHENE	79-01-6	UG_L	0.7	J	
8260C	TRICHLOROFUOROMETHANE	75-69-4	UG_L	1	UJ	c
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	U	
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	U	
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	0.1	J	

Sample Delivery Group				BETHPAGE-2		
Lab ID				SI7583-7		
Sample ID				RE105D1-GW-092815		
Sample Date				9/28/2015		
Sample Type				Groundwater		
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	U	
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	U	
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	11		
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.5	U	
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	0.5	U	
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	1.6		
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	U	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	U	
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	U	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	U	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	1.6	J	
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	U	
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	U	
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	U	
8260C	2-BUTANONE	78-93-3	UG_L	2.5	UJ	c
8260C	2-HEXANONE	591-78-6	UG_L	2.5	UJ	c
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	UJ	c
8260C	ACETONE	67-64-1	UG_L	2.5	UJ	c
8260C	BENZENE	71-43-2	UG_L	0.5	U	
8260C	BROMODICHLOROMETHANE	75-27-4	UG_L	0.5	U	
8260C	BROMOFORM	75-25-2	UG_L	0.5	U	
8260C	BROMOMETHANE	74-83-9	UG_L	1	UJ	c
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	UJ	c
8260C	CARBON TETRACHLORIDE	56-23-5	UG_L	0.5	U	
8260C	CHLOROBENZENE	108-90-7	UG_L	0.5	U	
8260C	CHLOROETHANE	75-00-3	UG_L	1	UJ	c
8260C	CHLOROFORM	67-66-3	UG_L	0.34	J	
8260C	CHLOROMETHANE	74-87-3	UG_L	1	UJ	c
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	1.6		
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG_L	0.5	U	
8260C	CYCLOHEXANE	110-82-7	UG_L	0.5	U	
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG_L	0.5	U	
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	0.74	J	c
8260C	ETHYLBENZENE	100-41-4	UG_L	0.5	U	
8260C	ISOPROPYLBENZENE	98-82-8	UG_L	0.5	U	
8260C	M- AND P-XYLENE	108-38-3/106-42	UG_L	1	U	
8260C	METHYL ACETATE	79-20-9	UG_L	0.75	U	
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	UJ	c
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	U	
8260C	METHYLENE CHLORIDE	75-09-2	UG_L	2.5	U	
8260C	O-XYLENE	95-47-6	UG_L	0.5	U	
8260C	STYRENE	100-42-5	UG_L	0.5	U	
8260C	TETRACHLOROETHENE	127-18-4	UG_L	0.5	UJ	c
8260C	TOLUENE	108-88-3	UG_L	0.5	U	
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	U	
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	U	
8260C	TRICHLOROETHENE	79-01-6	UG_L	94		
8260C	TRICHLOROFUOROMETHANE	75-69-4	UG_L	1	UJ	c
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	U	
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	U	
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	14		

Sample Delivery Group				BETHPAGE-2		
Lab ID				SI7583-8		
Sample ID				RE105D2-GW-092815		
Sample Date				9/28/2015		
Sample Type				Groundwater		
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	U	
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	U	
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	30		
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	1.1		
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	1.7		
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	6.8		
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	U	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	U	
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	U	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	U	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	3.6		
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	U	
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	U	
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	U	
8260C	2-BUTANONE	78-93-3	UG_L	2.5	UJ	c
8260C	2-HEXANONE	591-78-6	UG_L	2.5	UJ	c
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	UJ	c
8260C	ACETONE	67-64-1	UG_L	2.5	UJ	c
8260C	BENZENE	71-43-2	UG_L	0.5	U	
8260C	BROMODICHLOROMETHANE	75-27-4	UG_L	0.5	U	
8260C	BROMOFORM	75-25-2	UG_L	0.5	U	
8260C	BROMOMETHANE	74-83-9	UG_L	1	UJ	c
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	UJ	c
8260C	CARBON TETRACHLORIDE	56-23-5	UG_L	3.7		
8260C	CHLOROBENZENE	108-90-7	UG_L	0.5	U	
8260C	CHLOROETHANE	75-00-3	UG_L	1	UJ	c
8260C	CHLOROFORM	67-66-3	UG_L	2.2		
8260C	CHLOROMETHANE	74-87-3	UG_L	1	UJ	c
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	3.6		
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG_L	0.5	U	
8260C	CYCLOHEXANE	110-82-7	UG_L	0.5	U	
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG_L	0.5	U	
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	0.4	J	c
8260C	ETHYLBENZENE	100-41-4	UG_L	0.5	U	
8260C	ISOPROPYLBENZENE	98-82-8	UG_L	0.5	U	
8260C	M- AND P-XYLENE	108-38-3/106-42	UG_L	1	U	
8260C	METHYL ACETATE	79-20-9	UG_L	0.75	U	
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	UJ	c
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	U	
8260C	METHYLENE CHLORIDE	75-09-2	UG_L	2.5	U	
8260C	O-XYLENE	95-47-6	UG_L	0.5	U	
8260C	STYRENE	100-42-5	UG_L	0.5	U	
8260C	TETRACHLOROETHENE	127-18-4	UG_L	1.4	J	c
8260C	TOLUENE	108-88-3	UG_L	0.5	U	
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	U	
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	U	
8260C	TRICHLOROETHENE	79-01-6	UG_L	1900		
8260C	TRICHLOROFUOROMETHANE	75-69-4	UG_L	1	UJ	c
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	U	
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	U	
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	8.7		

Sample Delivery Group				BETHPAGE-2		
Lab ID				SI7583-9		
Sample ID				RE108D1-GW-092815		
Sample Date				9/28/2015		
Sample Type				Groundwater		
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	U	
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	U	
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	1.2		
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.5	U	
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	0.5	U	
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	0.5	U	
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	U	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	U	
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	U	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	U	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	0.39	J	
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	U	
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	U	
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	U	
8260C	2-BUTANONE	78-93-3	UG_L	2.5	UJ	c
8260C	2-HEXANONE	591-78-6	UG_L	2.5	UJ	c
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	UJ	c
8260C	ACETONE	67-64-1	UG_L	6.9	J	c
8260C	BENZENE	71-43-2	UG_L	0.5	U	
8260C	BROMODICHLOROMETHANE	75-27-4	UG_L	0.5	U	
8260C	BROMOFORM	75-25-2	UG_L	0.5	U	
8260C	BROMOMETHANE	74-83-9	UG_L	1	UJ	c
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	UJ	c
8260C	CARBON TETRACHLORIDE	56-23-5	UG_L	0.5	U	
8260C	CHLOROBENZENE	108-90-7	UG_L	0.5	U	
8260C	CHLOROETHANE	75-00-3	UG_L	1	UJ	c
8260C	CHLOROFORM	67-66-3	UG_L	0.5	U	
8260C	CHLOROMETHANE	74-87-3	UG_L	1	UJ	c
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	0.39	J	
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG_L	0.5	U	
8260C	CYCLOHEXANE	110-82-7	UG_L	0.5	U	
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG_L	0.5	U	
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	1	UJ	c
8260C	ETHYLBENZENE	100-41-4	UG_L	0.5	U	
8260C	ISOPROPYLBENZENE	98-82-8	UG_L	0.5	U	
8260C	M- AND P-XYLENE	108-38-3/106-42	UG_L	1	U	
8260C	METHYL ACETATE	79-20-9	UG_L	0.75	U	
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	UJ	c
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	U	
8260C	METHYLENE CHLORIDE	75-09-2	UG_L	2.5	U	
8260C	O-XYLENE	95-47-6	UG_L	0.5	U	
8260C	STYRENE	100-42-5	UG_L	0.5	U	
8260C	TETRACHLOROETHENE	127-18-4	UG_L	1.3	J	c
8260C	TOLUENE	108-88-3	UG_L	0.5	U	
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	U	
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	U	
8260C	TRICHLOROETHENE	79-01-6	UG_L	98		
8260C	TRICHLOROFUOROMETHANE	75-69-4	UG_L	1	UJ	c
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	U	
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	U	
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	7.9		

Sample Delivery Group				BETHPAGE-2		
Lab ID				SI7681-10		
Sample ID				DUPLICATE-GW-092915		
Sample Date				9/29/2015		
Sample Type				Field Duplicate		
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	U	
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	U	
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	15		
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.5	U	
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	0.61	J	
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	3.3		
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	U	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	UJ	c
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	U	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	U	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	1.8	J	
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	U	
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	U	
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	U	
8260C	2-BUTANONE	78-93-3	UG_L	2.5	U	
8260C	2-HEXANONE	591-78-6	UG_L	2.5	U	
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	U	
8260C	ACETONE	67-64-1	UG_L	2.5	U	
8260C	BENZENE	71-43-2	UG_L	0.5	U	
8260C	BROMODICHLOROMETHANE	75-27-4	UG_L	0.5	U	
8260C	BROMOFORM	75-25-2	UG_L	0.5	U	
8260C	BROMOMETHANE	74-83-9	UG_L	1	U	
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	U	
8260C	CARBON TETRACHLORIDE	56-23-5	UG_L	0.5	U	
8260C	CHLOROBENZENE	108-90-7	UG_L	0.5	U	
8260C	CHLOROETHANE	75-00-3	UG_L	1	U	
8260C	CHLOROFORM	67-66-3	UG_L	0.5	U	
8260C	CHLOROMETHANE	74-87-3	UG_L	1	U	
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	1.8		
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG_L	0.5	U	
8260C	CYCLOHEXANE	110-82-7	UG_L	0.5	U	
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG_L	0.5	U	
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	2.3	J	c
8260C	ETHYLBENZENE	100-41-4	UG_L	0.5	U	
8260C	ISOPROPYLBENZENE	98-82-8	UG_L	0.5	U	
8260C	M- AND P-XYLENE	108-38-3/106-42	UG_L	1	U	
8260C	METHYL ACETATE	79-20-9	UG_L	0.75	U	
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	U	
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	U	
8260C	METHYLENE CHLORIDE	75-09-2	UG_L	2.5	U	
8260C	O-XYLENE	95-47-6	UG_L	0.5	U	
8260C	STYRENE	100-42-5	UG_L	0.5	U	
8260C	TETRACHLOROETHENE	127-18-4	UG_L	0.5	U	
8260C	TOLUENE	108-88-3	UG_L	0.5	U	
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	U	
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	U	
8260C	TRICHLOROETHENE	79-01-6	UG_L	67		
8260C	TRICHLOROFUOROMETHANE	75-69-4	UG_L	1	U	
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	U	
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	U	
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	14		

Sample Delivery Group				BETHPAGE-2		
Lab ID				SI7681-11		
Sample ID				RE122D1-GW-093015		
Sample Date				9/30/2015		
Sample Type				Groundwater		
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	UJ	s
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	UJ	s
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	4.4	J	s
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.5	UJ	s
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	0.5	UJ	s
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	0.45	J	s
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	UJ	s
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	UJ	s,c
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	UJ	s
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	UJ	s
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	UJ	s
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	1	J	s
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	UJ	s
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	UJ	s
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	UJ	s
8260C	2-BUTANONE	78-93-3	UG_L	2.5	UJ	s
8260C	2-HEXANONE	591-78-6	UG_L	2.5	UJ	s
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	UJ	s
8260C	ACETONE	67-64-1	UG_L	2.5	UJ	s
8260C	BENZENE	71-43-2	UG_L	0.5	UJ	s
8260C	BROMODICHLOROMETHANE	75-27-4	UG_L	0.5	UJ	s
8260C	BROMOFORM	75-25-2	UG_L	0.5	UJ	s
8260C	BROMOMETHANE	74-83-9	UG_L	1	UJ	s
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	UJ	s
8260C	CARBON TETRACHLORIDE	56-23-5	UG_L	0.5	UJ	s
8260C	CHLOROBENZENE	108-90-7	UG_L	0.5	UJ	s
8260C	CHLOROETHANE	75-00-3	UG_L	1	UJ	s
8260C	CHLOROFORM	67-66-3	UG_L	0.5	UJ	s
8260C	CHLOROMETHANE	74-87-3	UG_L	1	UJ	s
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	1	J	s
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG_L	0.5	UJ	s
8260C	CYCLOHEXANE	110-82-7	UG_L	0.5	UJ	s
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG_L	0.5	UJ	s
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	1	UJ	s,c
8260C	ETHYLBENZENE	100-41-4	UG_L	0.5	UJ	s
8260C	ISOPROPYLBENZENE	98-82-8	UG_L	0.5	UJ	s
8260C	M- AND P-XYLENE	108-38-3/106-42	UG_L	1	UJ	s
8260C	METHYL ACETATE	79-20-9	UG_L	0.75	UJ	s
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	UJ	s
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	UJ	s
8260C	METHYLENE CHLORIDE	75-09-2	UG_L	2.5	UJ	s
8260C	O-XYLENE	95-47-6	UG_L	0.5	UJ	s
8260C	STYRENE	100-42-5	UG_L	0.5	UJ	s
8260C	TETRACHLOROETHENE	127-18-4	UG_L	1	J	s
8260C	TOLUENE	108-88-3	UG_L	0.5	UJ	s
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	UJ	s
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	UJ	s
8260C	TRICHLOROETHENE	79-01-6	UG_L	600	J	s
8260C	TRICHLOROFUOROMETHANE	75-69-4	UG_L	1	UJ	s
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	UJ	s
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	UJ	s
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	13		

Sample Delivery Group				BETHPAGE-2		
Lab ID				SI7681-12		
Sample ID				RE122D2-GW-093015		
Sample Date				9/30/2015		
Sample Type				Groundwater		
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	UJ	s
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	UJ	s
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	20	J	s
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	2.4	J	s
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	0.5	UJ	s
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	8.9	J	s
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	UJ	s
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	UJ	s,c
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	UJ	s
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	UJ	s
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	UJ	s
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	4.9	J	s
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	UJ	s
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	UJ	s
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	UJ	s
8260C	2-BUTANONE	78-93-3	UG_L	2.5	UJ	s
8260C	2-HEXANONE	591-78-6	UG_L	2.5	UJ	s
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	UJ	s
8260C	ACETONE	67-64-1	UG_L	2.5	UJ	s
8260C	BENZENE	71-43-2	UG_L	0.5	UJ	s
8260C	BROMODICHLOROMETHANE	75-27-4	UG_L	0.5	UJ	s
8260C	BROMOFORM	75-25-2	UG_L	0.5	UJ	s
8260C	BROMOMETHANE	74-83-9	UG_L	1	UJ	s
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	UJ	s
8260C	CARBON TETRACHLORIDE	56-23-5	UG_L	2	J	s
8260C	CHLOROBENZENE	108-90-7	UG_L	0.5	UJ	s
8260C	CHLOROETHANE	75-00-3	UG_L	1	UJ	s
8260C	CHLOROFORM	67-66-3	UG_L	1.6	J	s
8260C	CHLOROMETHANE	74-87-3	UG_L	1	UJ	s
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	4.9	J	s
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG_L	0.5	UJ	s
8260C	CYCLOHEXANE	110-82-7	UG_L	0.5	UJ	s
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG_L	0.5	UJ	s
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	0.36	J	s,c
8260C	ETHYLBENZENE	100-41-4	UG_L	0.5	UJ	s
8260C	ISOPROPYLBENZENE	98-82-8	UG_L	0.5	UJ	s
8260C	M- AND P-XYLENE	108-38-3/106-42	UG_L	1	UJ	s
8260C	METHYL ACETATE	79-20-9	UG_L	0.75	UJ	s
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	UJ	s
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	UJ	s
8260C	METHYLENE CHLORIDE	75-09-2	UG_L	2.5	UJ	s
8260C	O-XYLENE	95-47-6	UG_L	0.5	UJ	s
8260C	STYRENE	100-42-5	UG_L	0.5	UJ	s
8260C	TETRACHLOROETHENE	127-18-4	UG_L	2.6	J	s
8260C	TOLUENE	108-88-3	UG_L	0.5	UJ	s
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	UJ	s
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	UJ	s
8260C	TRICHLOROETHENE	79-01-6	UG_L	5200	J	s
8260C	TRICHLOROFUOROMETHANE	75-69-4	UG_L	1	UJ	s
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	UJ	s
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	UJ	s
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	17		



Sample Delivery Group				BETHPAGE-2		
Lab ID				SI7681-13		
Sample ID				RE122D3-GW-093015		
Sample Date				9/30/2015		
Sample Type				Groundwater		
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	UJ	s
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	UJ	s
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	0.5	UJ	s
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.5	UJ	s
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	0.5	UJ	s
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	0.5	UJ	s
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	UJ	s
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	UJ	s,c
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	UJ	s
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	UJ	s
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	UJ	s
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	1	UJ	s
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	UJ	s
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	UJ	s
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	UJ	s
8260C	2-BUTANONE	78-93-3	UG_L	2.5	UJ	s
8260C	2-HEXANONE	591-78-6	UG_L	2.5	UJ	s
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	UJ	s
8260C	ACETONE	67-64-1	UG_L	2.5	UJ	s
8260C	BENZENE	71-43-2	UG_L	0.5	UJ	s
8260C	BROMODICHLOROMETHANE	75-27-4	UG_L	0.5	UJ	s
8260C	BROMOFORM	75-25-2	UG_L	0.5	UJ	s
8260C	BROMOMETHANE	74-83-9	UG_L	1	UJ	s
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	UJ	s
8260C	CARBON TETRACHLORIDE	56-23-5	UG_L	0.5	UJ	s
8260C	CHLOROBENZENE	108-90-7	UG_L	0.5	UJ	s
8260C	CHLOROETHANE	75-00-3	UG_L	1	UJ	s
8260C	CHLOROFORM	67-66-3	UG_L	0.5	UJ	s
8260C	CHLOROMETHANE	74-87-3	UG_L	1	UJ	s
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	0.5	UJ	s
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG_L	0.5	UJ	s
8260C	CYCLOHEXANE	110-82-7	UG_L	0.5	UJ	s
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG_L	0.5	UJ	s
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	1	UJ	s,c
8260C	ETHYLBENZENE	100-41-4	UG_L	0.5	UJ	s
8260C	ISOPROPYLBENZENE	98-82-8	UG_L	0.5	UJ	s
8260C	M- AND P-XYLENE	108-38-3/106-42	UG_L	1	UJ	s
8260C	METHYL ACETATE	79-20-9	UG_L	0.75	UJ	s
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	UJ	s
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	UJ	s
8260C	METHYLENE CHLORIDE	75-09-2	UG_L	2.5	UJ	s
8260C	O-XYLENE	95-47-6	UG_L	0.5	UJ	s
8260C	STYRENE	100-42-5	UG_L	0.5	UJ	s
8260C	TETRACHLOROETHENE	127-18-4	UG_L	0.5	UJ	s
8260C	TOLUENE	108-88-3	UG_L	0.5	UJ	s
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	UJ	s
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	UJ	s
8260C	TRICHLOROETHENE	79-01-6	UG_L	10	J	s
8260C	TRICHLOROFUOROMETHANE	75-69-4	UG_L	1	UJ	s
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	UJ	s
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	UJ	s
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	0.17	U	

Sample Delivery Group				BETHPAGE-2		
Lab ID				SI7681-14		
Sample ID				RE103D1-GW-093015		
Sample Date				9/30/2015		
Sample Type				Groundwater		
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	U	
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	U	
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	12		
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.52	J	
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	1		
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	8.4		
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	U	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	U	
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	U	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	U	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	3.6		
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	U	
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	U	
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	U	
8260C	2-BUTANONE	78-93-3	UG_L	2.5	UJ	c
8260C	2-HEXANONE	591-78-6	UG_L	2.5	UJ	c
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	UJ	c
8260C	ACETONE	67-64-1	UG_L	2.5	UJ	c
8260C	BENZENE	71-43-2	UG_L	0.5	U	
8260C	BROMODICHLOROMETHANE	75-27-4	UG_L	0.5	U	
8260C	BROMOFORM	75-25-2	UG_L	0.5	U	
8260C	BROMOMETHANE	74-83-9	UG_L	1	U	
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	UJ	c
8260C	CARBON TETRACHLORIDE	56-23-5	UG_L	0.31	J	
8260C	CHLOROBENZENE	108-90-7	UG_L	0.5	U	
8260C	CHLOROETHANE	75-00-3	UG_L	1	U	
8260C	CHLOROFORM	67-66-3	UG_L	0.61	J	
8260C	CHLOROMETHANE	74-87-3	UG_L	1	U	
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	3.6		
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG_L	0.5	U	
8260C	CYCLOHEXANE	110-82-7	UG_L	0.5	U	
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG_L	0.5	U	
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	0.26	J	
8260C	ETHYLBENZENE	100-41-4	UG_L	0.5	U	
8260C	ISOPROPYLBENZENE	98-82-8	UG_L	0.5	U	
8260C	M- AND P-XYLENE	108-38-3/106-42	UG_L	1	U	
8260C	METHYL ACETATE	79-20-9	UG_L	0.75	U	
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	UJ	c
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	U	
8260C	METHYLENE CHLORIDE	75-09-2	UG_L	2.5	U	
8260C	O-XYLENE	95-47-6	UG_L	0.5	UJ	c
8260C	STYRENE	100-42-5	UG_L	0.5	U	
8260C	TETRACHLOROETHENE	127-18-4	UG_L	6.7		
8260C	TOLUENE	108-88-3	UG_L	0.5	U	
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	U	
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	U	
8260C	TRICHLOROETHENE	79-01-6	UG_L	860		
8260C	TRICHLOROFUOROMETHANE	75-69-4	UG_L	1	U	
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	U	
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	U	
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	24		

Sample Delivery Group				BETHPAGE-2		
Lab ID				SI7681-15		
Sample ID				RE103D2-GW-093015		
Sample Date				9/30/2015		
Sample Type				Groundwater		
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	U	
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	U	
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	4.9		
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.54	J	
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	0.72	J	
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	1.3		
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	U	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	U	
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	U	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	U	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	1.5	J	
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	U	
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	U	
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	U	
8260C	2-BUTANONE	78-93-3	UG_L	2.5	UJ	c
8260C	2-HEXANONE	591-78-6	UG_L	2.5	UJ	c
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	UJ	c
8260C	ACETONE	67-64-1	UG_L	2.5	UJ	c
8260C	BENZENE	71-43-2	UG_L	0.5	U	
8260C	BROMODICHLOROMETHANE	75-27-4	UG_L	0.5	U	
8260C	BROMOFORM	75-25-2	UG_L	0.5	U	
8260C	BROMOMETHANE	74-83-9	UG_L	1	U	
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	UJ	c
8260C	CARBON TETRACHLORIDE	56-23-5	UG_L	0.46	J	
8260C	CHLOROBENZENE	108-90-7	UG_L	0.5	U	
8260C	CHLOROETHANE	75-00-3	UG_L	1	U	
8260C	CHLOROFORM	67-66-3	UG_L	1	U	
8260C	CHLOROMETHANE	74-87-3	UG_L	1	U	
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	1.5		
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG_L	0.5	U	
8260C	CYCLOHEXANE	110-82-7	UG_L	0.5	U	
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG_L	0.5	U	
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	1	U	
8260C	ETHYLBENZENE	100-41-4	UG_L	0.5	U	
8260C	ISOPROPYLBENZENE	98-82-8	UG_L	0.5	U	
8260C	M- AND P-XYLENE	108-38-3/106-42	UG_L	1	U	
8260C	METHYL ACETATE	79-20-9	UG_L	0.75	U	
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	UJ	c
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	U	
8260C	METHYLENE CHLORIDE	75-09-2	UG_L	2.5	U	
8260C	O-XYLENE	95-47-6	UG_L	0.5	UJ	c
8260C	STYRENE	100-42-5	UG_L	0.5	U	
8260C	TETRACHLOROETHENE	127-18-4	UG_L	4		
8260C	TOLUENE	108-88-3	UG_L	0.5	U	
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	U	
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	U	
8260C	TRICHLOROETHENE	79-01-6	UG_L	830		
8260C	TRICHLOROFUOROMETHANE	75-69-4	UG_L	1	U	
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	U	
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	U	
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	2.1		

Sample Delivery Group				BETHPAGE-2		
Lab ID				SI7681-16		
Sample ID				RE103D3-GW-093015		
Sample Date				9/30/2015		
Sample Type				Groundwater		
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	U	
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	U	
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	2.3	J	s
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.5	U	
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	0.5	J	s
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	0.64	J	s
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	U	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	U	
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	U	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	U	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	0.96	J	s
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	U	
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	U	
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	U	
8260C	2-BUTANONE	78-93-3	UG_L	2.5	UJ	c
8260C	2-HEXANONE	591-78-6	UG_L	2.5	UJ	c
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	UJ	c
8260C	ACETONE	67-64-1	UG_L	2.5	UJ	c
8260C	BENZENE	71-43-2	UG_L	0.5	U	
8260C	BROMODICHLOROMETHANE	75-27-4	UG_L	0.5	U	
8260C	BROMOFORM	75-25-2	UG_L	0.5	U	
8260C	BROMOMETHANE	74-83-9	UG_L	1	U	
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	UJ	c
8260C	CARBON TETRACHLORIDE	56-23-5	UG_L	0.5	U	
8260C	CHLOROBENZENE	108-90-7	UG_L	0.5	U	
8260C	CHLOROETHANE	75-00-3	UG_L	1	U	
8260C	CHLOROFORM	67-66-3	UG_L	0.74	J	s
8260C	CHLOROMETHANE	74-87-3	UG_L	1	U	
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	0.96	J	s
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG_L	0.5	U	
8260C	CYCLOHEXANE	110-82-7	UG_L	0.5	U	
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG_L	0.5	U	
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	1	U	
8260C	ETHYLBENZENE	100-41-4	UG_L	0.5	U	
8260C	ISOPROPYLBENZENE	98-82-8	UG_L	0.5	U	
8260C	M- AND P-XYLENE	108-38-3/106-42	UG_L	1	U	
8260C	METHYL ACETATE	79-20-9	UG_L	0.75	U	
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	UJ	c
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	U	
8260C	METHYLENE CHLORIDE	75-09-2	UG_L	2.5	U	
8260C	O-XYLENE	95-47-6	UG_L	0.5	UJ	c
8260C	STYRENE	100-42-5	UG_L	0.5	U	
8260C	TETRACHLOROETHENE	127-18-4	UG_L	3.6	J	s
8260C	TOLUENE	108-88-3	UG_L	0.5	U	
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	U	
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	U	
8260C	TRICHLOROETHENE	79-01-6	UG_L	470		
8260C	TRICHLOROFUOROMETHANE	75-69-4	UG_L	1	U	
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	U	
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	U	
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	1		

Sample Delivery Group				BETHPAGE-2		
Lab ID				SI7681-17		
Sample ID				TB01-093015		
Sample Date				9/30/2015		
Sample Type				Trip Blank		
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	U	
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	U	
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	0.5	U	
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.5	U	
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	0.5	U	
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	0.5	U	
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	U	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	U	
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	U	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	U	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	1	U	
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	U	
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	U	
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	U	
8260C	2-BUTANONE	78-93-3	UG_L	2.5	UJ	c
8260C	2-HEXANONE	591-78-6	UG_L	2.5	UJ	c
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	UJ	c
8260C	ACETONE	67-64-1	UG_L	2.5	UJ	c
8260C	BENZENE	71-43-2	UG_L	0.5	U	
8260C	BROMODICHLOROMETHANE	75-27-4	UG_L	0.5	U	
8260C	BROMOFORM	75-25-2	UG_L	0.5	U	
8260C	BROMOMETHANE	74-83-9	UG_L	2		
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	UJ	bl,c
8260C	CARBON TETRACHLORIDE	56-23-5	UG_L	0.5	U	
8260C	CHLOROBENZENE	108-90-7	UG_L	0.5	U	
8260C	CHLOROETHANE	75-00-3	UG_L	1	U	
8260C	CHLOROFORM	67-66-3	UG_L	0.5	U	
8260C	CHLOROMETHANE	74-87-3	UG_L	0.82	J	
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	0.5	U	
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG_L	0.5	U	
8260C	CYCLOHEXANE	110-82-7	UG_L	0.5	U	
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG_L	0.5	U	
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	1	U	
8260C	ETHYLBENZENE	100-41-4	UG_L	0.5	U	
8260C	ISOPROPYLBENZENE	98-82-8	UG_L	0.5	U	
8260C	M- AND P-XYLENE	108-38-3/106-42	UG_L	1	U	
8260C	METHYL ACETATE	79-20-9	UG_L	0.75	U	
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	UJ	c
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	U	
8260C	METHYLENE CHLORIDE	75-09-2	UG_L	2.5	U	
8260C	O-XYLENE	95-47-6	UG_L	0.5	UJ	c
8260C	STYRENE	100-42-5	UG_L	0.5	U	
8260C	TETRACHLOROETHENE	127-18-4	UG_L	0.5	U	
8260C	TOLUENE	108-88-3	UG_L	0.5	U	
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	U	
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	U	
8260C	TRICHLOROETHENE	79-01-6	UG_L	0.5	U	
8260C	TRICHLOROFUOROMETHANE	75-69-4	UG_L	1	U	
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	U	
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	U	
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	NA		

Sample Delivery Group				BETHPAGE-2		
Lab ID				SI7681-1RA		
Sample ID				RE123D1-GW-092915		
Sample Date				9/29/2015		
Sample Type				Groundwater		
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	U	
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	U	
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	0.5	U	
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.5	U	
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	0.5	U	
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	0.42	J	
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	U	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	U	
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	U	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	U	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	0.5	J	
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	U	
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	U	
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	U	
8260C	2-BUTANONE	78-93-3	UG_L	2.5	UJ	c
8260C	2-HEXANONE	591-78-6	UG_L	2.5	UJ	c
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	UJ	c
8260C	ACETONE	67-64-1	UG_L	5.4	J	c
8260C	BENZENE	71-43-2	UG_L	0.5	U	
8260C	BROMODICHLOROMETHANE	75-27-4	UG_L	0.5	U	
8260C	BROMOFORM	75-25-2	UG_L	0.5	U	
8260C	BROMOMETHANE	74-83-9	UG_L	1	U	
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	UJ	c
8260C	CARBON TETRACHLORIDE	56-23-5	UG_L	0.5	U	
8260C	CHLOROBENZENE	108-90-7	UG_L	0.5	U	
8260C	CHLOROETHANE	75-00-3	UG_L	1	U	
8260C	CHLOROFORM	67-66-3	UG_L	0.5	U	
8260C	CHLOROMETHANE	74-87-3	UG_L	1	U	
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	0.5	J	
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG_L	0.5	U	
8260C	CYCLOHEXANE	110-82-7	UG_L	0.5	U	
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG_L	0.5	U	
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	1	U	
8260C	ETHYLBENZENE	100-41-4	UG_L	0.5	U	
8260C	ISOPROPYLBENZENE	98-82-8	UG_L	0.5	U	
8260C	M- AND P-XYLENE	108-38-3/106-42	UG_L	1	U	
8260C	METHYL ACETATE	79-20-9	UG_L	0.75	U	
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	U	
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	U	
8260C	METHYLENE CHLORIDE	75-09-2	UG_L	2.5	U	
8260C	O-XYLENE	95-47-6	UG_L	0.5	UJ	c
8260C	STYRENE	100-42-5	UG_L	0.5	U	
8260C	TETRACHLOROETHENE	127-18-4	UG_L	3.8	J	c
8260C	TOLUENE	108-88-3	UG_L	0.5	U	
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	U	
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	U	
8260C	TRICHLOROETHENE	79-01-6	UG_L	12		
8260C	TRICHLOROFUOROMETHANE	75-69-4	UG_L	1	U	
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	U	
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	U	
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	6.6		

Sample Delivery Group				BETHPAGE-2		
Lab ID				SI7681-2RA		
Sample ID				RE123D2-GW-092915		
Sample Date				9/29/2015		
Sample Type				Groundwater		
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	U	
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	U	
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	0.5	U	
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.5	U	
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	0.5	U	
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	0.5	U	
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	U	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	U	
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	U	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	U	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	1	U	
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	U	
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	U	
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	U	
8260C	2-BUTANONE	78-93-3	UG_L	2.5	UJ	c
8260C	2-HEXANONE	591-78-6	UG_L	2.5	UJ	c
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	UJ	c
8260C	ACETONE	67-64-1	UG_L	2.5	UJ	c
8260C	BENZENE	71-43-2	UG_L	0.5	U	
8260C	BROMODICHLOROMETHANE	75-27-4	UG_L	0.5	U	
8260C	BROMOFORM	75-25-2	UG_L	0.5	U	
8260C	BROMOMETHANE	74-83-9	UG_L	1	U	
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	UJ	c
8260C	CARBON TETRACHLORIDE	56-23-5	UG_L	0.5	U	
8260C	CHLOROBENZENE	108-90-7	UG_L	0.5	U	
8260C	CHLOROETHANE	75-00-3	UG_L	1	U	
8260C	CHLOROFORM	67-66-3	UG_L	0.5	U	
8260C	CHLOROMETHANE	74-87-3	UG_L	1	UJ	bt
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	0.5	U	
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG_L	0.5	U	
8260C	CYCLOHEXANE	110-82-7	UG_L	0.5	U	
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG_L	0.5	U	
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	1	U	
8260C	ETHYLBENZENE	100-41-4	UG_L	0.5	U	
8260C	ISOPROPYLBENZENE	98-82-8	UG_L	0.5	U	
8260C	M- AND P-XYLENE	108-38-3/106-42	UG_L	1	U	
8260C	METHYL ACETATE	79-20-9	UG_L	0.75	U	
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	U	
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	U	
8260C	METHYLENE CHLORIDE	75-09-2	UG_L	2.5	U	
8260C	O-XYLENE	95-47-6	UG_L	0.5	UJ	c
8260C	STYRENE	100-42-5	UG_L	0.5	U	
8260C	TETRACHLOROETHENE	127-18-4	UG_L	3.7	J	c
8260C	TOLUENE	108-88-3	UG_L	0.5	U	
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	U	
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	U	
8260C	TRICHLOROETHENE	79-01-6	UG_L	1.4		
8260C	TRICHLOROFUOROMETHANE	75-69-4	UG_L	1	U	
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	U	
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	U	
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	0.93		

Sample Delivery Group				BETHPAGE-2		
Lab ID				SI7681-3RA		
Sample ID				RE123D3-GW-092915		
Sample Date				9/29/2015		
Sample Type				Groundwater		
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	U	
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	U	
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	0.5	U	
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.5	U	
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	0.5	U	
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	0.5	U	
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	U	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	U	
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	U	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	U	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	1	U	
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	U	
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	U	
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	U	
8260C	2-BUTANONE	78-93-3	UG_L	2.5	UJ	c
8260C	2-HEXANONE	591-78-6	UG_L	2.5	UJ	c
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	UJ	c
8260C	ACETONE	67-64-1	UG_L	2.5	UJ	c
8260C	BENZENE	71-43-2	UG_L	0.5	U	
8260C	BROMODICHLOROMETHANE	75-27-4	UG_L	0.5	U	
8260C	BROMOFORM	75-25-2	UG_L	0.5	U	
8260C	BROMOMETHANE	74-83-9	UG_L	1	U	
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	UJ	c
8260C	CARBON TETRACHLORIDE	56-23-5	UG_L	0.5	U	
8260C	CHLOROBENZENE	108-90-7	UG_L	0.5	U	
8260C	CHLOROETHANE	75-00-3	UG_L	1	U	
8260C	CHLOROFORM	67-66-3	UG_L	0.5	U	
8260C	CHLOROMETHANE	74-87-3	UG_L	1	UJ	bt
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	0.5	U	
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG_L	0.5	U	
8260C	CYCLOHEXANE	110-82-7	UG_L	0.5	U	
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG_L	0.5	U	
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	1	U	
8260C	ETHYLBENZENE	100-41-4	UG_L	0.5	U	
8260C	ISOPROPYLBENZENE	98-82-8	UG_L	0.5	U	
8260C	M- AND P-XYLENE	108-38-3/106-42	UG_L	1	U	
8260C	METHYL ACETATE	79-20-9	UG_L	0.75	U	
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	U	
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	U	
8260C	METHYLENE CHLORIDE	75-09-2	UG_L	2.5	U	
8260C	O-XYLENE	95-47-6	UG_L	0.5	UJ	c
8260C	STYRENE	100-42-5	UG_L	0.5	U	
8260C	TETRACHLOROETHENE	127-18-4	UG_L	0.5	UJ	c
8260C	TOLUENE	108-88-3	UG_L	0.5	U	
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	U	
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	U	
8260C	TRICHLOROETHENE	79-01-6	UG_L	0.5	U	
8260C	TRICHLOROFUOROMETHANE	75-69-4	UG_L	1	U	
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	U	
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	U	
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	0.17	U	



Sample Delivery Group				BETHPAGE-2		
Lab ID				SI7681-4		
Sample ID				RE120D1-GW-092915		
Sample Date				9/29/2015		
Sample Type				Groundwater		
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	1.4		
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	U	
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	36		
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	1.6		
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	2.9		
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	20		
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	U	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	UJ	c
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	U	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	U	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	2.7		
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	U	
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	U	
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	U	
8260C	2-BUTANONE	78-93-3	UG_L	2.5	U	
8260C	2-HEXANONE	591-78-6	UG_L	2.5	U	
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	U	
8260C	ACETONE	67-64-1	UG_L	2.5	U	
8260C	BENZENE	71-43-2	UG_L	0.5	U	
8260C	BROMODICHLOROMETHANE	75-27-4	UG_L	0.5	U	
8260C	BROMOFORM	75-25-2	UG_L	0.5	U	
8260C	BROMOMETHANE	74-83-9	UG_L	1	U	
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	U	
8260C	CARBON TETRACHLORIDE	56-23-5	UG_L	0.77	J	
8260C	CHLOROBENZENE	108-90-7	UG_L	0.5	U	
8260C	CHLOROETHANE	75-00-3	UG_L	1	U	
8260C	CHLOROFORM	67-66-3	UG_L	0.5	U	
8260C	CHLOROMETHANE	74-87-3	UG_L	1	U	
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	2.7		
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG_L	0.5	U	
8260C	CYCLOHEXANE	110-82-7	UG_L	0.5	U	
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG_L	0.5	U	
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	1	UJ	c
8260C	ETHYLBENZENE	100-41-4	UG_L	0.5	U	
8260C	ISOPROPYLBENZENE	98-82-8	UG_L	0.5	U	
8260C	M- AND P-XYLENE	108-38-3/106-42	UG_L	1	U	
8260C	METHYL ACETATE	79-20-9	UG_L	0.75	U	
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	U	
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	U	
8260C	METHYLENE CHLORIDE	75-09-2	UG_L	2.5	U	
8260C	O-XYLENE	95-47-6	UG_L	0.5	U	
8260C	STYRENE	100-42-5	UG_L	0.5	U	
8260C	TETRACHLOROETHENE	127-18-4	UG_L	1.9		
8260C	TOLUENE	108-88-3	UG_L	0.5	U	
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	U	
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	U	
8260C	TRICHLOROETHENE	79-01-6	UG_L	1300	J	s
8260C	TRICHLOROFUOROMETHANE	75-69-4	UG_L	0.44	J	
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	U	
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	U	
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	25		

Sample Delivery Group				BETHPAGE-2		
Lab ID				SI7681-5		
Sample ID				RE120D2-GW-092915		
Sample Date				9/29/2015		
Sample Type				Groundwater		
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	U	
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	U	
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	21		
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.45	J	
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	0.5	U	
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	4.9		
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	U	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	UJ	c
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	U	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	U	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	2.7		
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	U	
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	U	
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	U	
8260C	2-BUTANONE	78-93-3	UG_L	2.5	U	
8260C	2-HEXANONE	591-78-6	UG_L	2.5	U	
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	U	
8260C	ACETONE	67-64-1	UG_L	2.5	U	
8260C	BENZENE	71-43-2	UG_L	0.5	U	
8260C	BROMODICHLOROMETHANE	75-27-4	UG_L	0.5	U	
8260C	BROMOFORM	75-25-2	UG_L	0.5	U	
8260C	BROMOMETHANE	74-83-9	UG_L	1	U	
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	U	
8260C	CARBON TETRACHLORIDE	56-23-5	UG_L	0.72	J	
8260C	CHLOROBENZENE	108-90-7	UG_L	0.5	U	
8260C	CHLOROETHANE	75-00-3	UG_L	1	U	
8260C	CHLOROFORM	67-66-3	UG_L	0.5	U	
8260C	CHLOROMETHANE	74-87-3	UG_L	1	U	
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	2.7		
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG_L	0.5	U	
8260C	CYCLOHEXANE	110-82-7	UG_L	0.5	U	
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG_L	0.5	U	
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	0.31	J	c
8260C	ETHYLBENZENE	100-41-4	UG_L	0.5	U	
8260C	ISOPROPYLBENZENE	98-82-8	UG_L	0.5	U	
8260C	M- AND P-XYLENE	108-38-3/106-42	UG_L	1	U	
8260C	METHYL ACETATE	79-20-9	UG_L	0.75	U	
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	U	
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	U	
8260C	METHYLENE CHLORIDE	75-09-2	UG_L	2.5	U	
8260C	O-XYLENE	95-47-6	UG_L	0.5	U	
8260C	STYRENE	100-42-5	UG_L	0.5	U	
8260C	TETRACHLOROETHENE	127-18-4	UG_L	3.1		
8260C	TOLUENE	108-88-3	UG_L	0.5	U	
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	U	
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	U	
8260C	TRICHLOROETHENE	79-01-6	UG_L	760	J	s
8260C	TRICHLOROFUOROMETHANE	75-69-4	UG_L	1	U	
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	U	
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	U	
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	16		

Sample Delivery Group				BETHPAGE-2		
Lab ID				SI7681-6		
Sample ID				RE120D3-GW-092915		
Sample Date				9/29/2015		
Sample Type				Groundwater		
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	U	
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	U	
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	2.9		
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.5	U	
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	0.5	U	
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	0.5	U	
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	U	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	UJ	c
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	U	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	U	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	1	U	
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	U	
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	U	
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	U	
8260C	2-BUTANONE	78-93-3	UG_L	2.5	U	
8260C	2-HEXANONE	591-78-6	UG_L	2.5	U	
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	U	
8260C	ACETONE	67-64-1	UG_L	2.5	U	
8260C	BENZENE	71-43-2	UG_L	0.5	U	
8260C	BROMODICHLOROMETHANE	75-27-4	UG_L	0.5	U	
8260C	BROMOFORM	75-25-2	UG_L	0.5	U	
8260C	BROMOMETHANE	74-83-9	UG_L	1	U	
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	U	
8260C	CARBON TETRACHLORIDE	56-23-5	UG_L	0.5	U	
8260C	CHLOROBENZENE	108-90-7	UG_L	0.5	U	
8260C	CHLOROETHANE	75-00-3	UG_L	1	U	
8260C	CHLOROFORM	67-66-3	UG_L	0.5	U	
8260C	CHLOROMETHANE	74-87-3	UG_L	1	U	
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	0.5	U	
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG_L	0.5	U	
8260C	CYCLOHEXANE	110-82-7	UG_L	0.5	U	
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG_L	0.5	U	
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	1	UJ	c
8260C	ETHYLBENZENE	100-41-4	UG_L	0.5	U	
8260C	ISOPROPYLBENZENE	98-82-8	UG_L	0.5	U	
8260C	M- AND P-XYLENE	108-38-3/106-42	UG_L	1	U	
8260C	METHYL ACETATE	79-20-9	UG_L	0.75	U	
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	U	
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	U	
8260C	METHYLENE CHLORIDE	75-09-2	UG_L	2.5	U	
8260C	O-XYLENE	95-47-6	UG_L	0.5	U	
8260C	STYRENE	100-42-5	UG_L	0.5	U	
8260C	TETRACHLOROETHENE	127-18-4	UG_L	0.5	U	
8260C	TOLUENE	108-88-3	UG_L	0.5	U	
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	U	
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	U	
8260C	TRICHLOROETHENE	79-01-6	UG_L	120		
8260C	TRICHLOROFUOROMETHANE	75-69-4	UG_L	1	U	
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	U	
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	U	
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	1.2		

Sample Delivery Group				BETHPAGE-2		
Lab ID				SI7681-7		
Sample ID				TT101D-GW-092915		
Sample Date				9/29/2015		
Sample Type				Groundwater		
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	U	
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	U	
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	14		
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.5	U	
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	0.43	J	
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	3.1		
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	U	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	UJ	c
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	U	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	U	
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	U	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	1.9	J	
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	U	
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	U	
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	U	
8260C	2-BUTANONE	78-93-3	UG_L	2.5	U	
8260C	2-HEXANONE	591-78-6	UG_L	2.5	U	
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	U	
8260C	ACETONE	67-64-1	UG_L	2.5	U	
8260C	BENZENE	71-43-2	UG_L	0.5	U	
8260C	BROMODICHLOROMETHANE	75-27-4	UG_L	0.5	U	
8260C	BROMOFORM	75-25-2	UG_L	0.5	U	
8260C	BROMOMETHANE	74-83-9	UG_L	1	U	
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	U	
8260C	CARBON TETRACHLORIDE	56-23-5	UG_L	0.5	U	
8260C	CHLOROBENZENE	108-90-7	UG_L	0.5	U	
8260C	CHLOROETHANE	75-00-3	UG_L	1	U	
8260C	CHLOROFORM	67-66-3	UG_L	0.5	U	
8260C	CHLOROMETHANE	74-87-3	UG_L	1	U	
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	1.9		
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG_L	0.5	U	
8260C	CYCLOHEXANE	110-82-7	UG_L	0.5	U	
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG_L	0.5	U	
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	2	J	c
8260C	ETHYLBENZENE	100-41-4	UG_L	0.5	U	
8260C	ISOPROPYLBENZENE	98-82-8	UG_L	0.5	U	
8260C	M- AND P-XYLENE	108-38-3/106-42	UG_L	1	U	
8260C	METHYL ACETATE	79-20-9	UG_L	0.75	U	
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	U	
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	U	
8260C	METHYLENE CHLORIDE	75-09-2	UG_L	2.5	U	
8260C	O-XYLENE	95-47-6	UG_L	0.5	U	
8260C	STYRENE	100-42-5	UG_L	0.5	U	
8260C	TETRACHLOROETHENE	127-18-4	UG_L	0.5	U	
8260C	TOLUENE	108-88-3	UG_L	0.5	U	
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	U	
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	U	
8260C	TRICHLOROETHENE	79-01-6	UG_L	67		
8260C	TRICHLOROFUOROMETHANE	75-69-4	UG_L	1	U	
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	U	
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	U	
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	13		

Sample Delivery Group				BETHPAGE-2		
Lab ID				SI7681-8		
Sample ID				TT101D1-GW-092915		
Sample Date				9/29/2015		
Sample Type				Groundwater		
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	UJ	s
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	UJ	s
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	13	J	s
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.49	J	s
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	0.5	UJ	s
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	4	J	s
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	UJ	s
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	UJ	s,c
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	UJ	s
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	UJ	s
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	UJ	s
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	0.92	J	s
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	UJ	s
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	UJ	s
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	UJ	s
8260C	2-BUTANONE	78-93-3	UG_L	2.5	UJ	s
8260C	2-HEXANONE	591-78-6	UG_L	2.5	UJ	s
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	UJ	s
8260C	ACETONE	67-64-1	UG_L	2.5	UJ	s
8260C	BENZENE	71-43-2	UG_L	0.5	UJ	s
8260C	BROMODICHLOROMETHANE	75-27-4	UG_L	0.5	UJ	s
8260C	BROMOFORM	75-25-2	UG_L	0.5	UJ	s
8260C	BROMOMETHANE	74-83-9	UG_L	1	UJ	s
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	UJ	s
8260C	CARBON TETRACHLORIDE	56-23-5	UG_L	1.6	J	s
8260C	CHLOROBENZENE	108-90-7	UG_L	0.5	UJ	s
8260C	CHLOROETHANE	75-00-3	UG_L	1	UJ	s
8260C	CHLOROFORM	67-66-3	UG_L	0.5	UJ	s
8260C	CHLOROMETHANE	74-87-3	UG_L	1	UJ	s
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	0.92	J	s
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG_L	0.5	UJ	s
8260C	CYCLOHEXANE	110-82-7	UG_L	0.5	UJ	s
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG_L	0.5	UJ	s
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	1.9	J	s,c
8260C	ETHYLBENZENE	100-41-4	UG_L	0.5	UJ	s
8260C	ISOPROPYLBENZENE	98-82-8	UG_L	0.5	UJ	s
8260C	M- AND P-XYLENE	108-38-3/106-42	UG_L	1	UJ	s
8260C	METHYL ACETATE	79-20-9	UG_L	0.75	UJ	s
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	UJ	s
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	UJ	s
8260C	METHYLENE CHLORIDE	75-09-2	UG_L	2.5	UJ	s
8260C	O-XYLENE	95-47-6	UG_L	0.5	UJ	s
8260C	STYRENE	100-42-5	UG_L	0.5	UJ	s
8260C	TETRACHLOROETHENE	127-18-4	UG_L	0.5	UJ	s
8260C	TOLUENE	108-88-3	UG_L	0.5	UJ	s
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	UJ	s
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	UJ	s
8260C	TRICHLOROETHENE	79-01-6	UG_L	170	J	s
8260C	TRICHLOROFUOROMETHANE	75-69-4	UG_L	1	UJ	s
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	UJ	s
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	UJ	s
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	11	J	s

Sample Delivery Group				BETHPAGE-2		
Lab ID				SI7681-9		
Sample ID				TT101D2-GW-092915		
Sample Date				9/29/2015		
Sample Type				Groundwater		
Method	Analyte	CAS No	Units	Result	Qual	RC
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG_L	0.5	UJ	s
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG_L	0.5	UJ	s
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG_L	19	J	s
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG_L	0.6	J	s
8260C	1,1-DICHLOROETHANE	75-34-3	UG_L	0.5	UJ	s
8260C	1,1-DICHLOROETHENE	75-35-4	UG_L	4.7	J	s
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG_L	0.5	UJ	s
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG_L	0.75	UJ	s,c
8260C	1,2-DIBROMOETHANE	106-93-4	UG_L	0.5	UJ	s
8260C	1,2-DICHLOROBENZENE	95-50-1	UG_L	0.5	UJ	s
8260C	1,2-DICHLOROETHANE	107-06-2	UG_L	0.5	UJ	s
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG_L	1	J	s
8260C	1,2-DICHLOROPROPANE	78-87-5	UG_L	0.5	UJ	s
8260C	1,3-DICHLOROBENZENE	541-73-1	UG_L	0.5	UJ	s
8260C	1,4-DICHLOROBENZENE	106-46-7	UG_L	0.5	UJ	s
8260C	2-BUTANONE	78-93-3	UG_L	2.5	UJ	s
8260C	2-HEXANONE	591-78-6	UG_L	2.5	UJ	s
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	UJ	s
8260C	ACETONE	67-64-1	UG_L	2.5	UJ	s
8260C	BENZENE	71-43-2	UG_L	0.5	UJ	s
8260C	BROMODICHLOROMETHANE	75-27-4	UG_L	0.5	UJ	s
8260C	BROMOFORM	75-25-2	UG_L	0.5	UJ	s
8260C	BROMOMETHANE	74-83-9	UG_L	1	UJ	s
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	UJ	s
8260C	CARBON TETRACHLORIDE	56-23-5	UG_L	1.4	J	s
8260C	CHLOROBENZENE	108-90-7	UG_L	0.5	UJ	s
8260C	CHLOROETHANE	75-00-3	UG_L	1	UJ	s
8260C	CHLOROFORM	67-66-3	UG_L	0.5	UJ	s
8260C	CHLOROMETHANE	74-87-3	UG_L	1	UJ	s
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG_L	1	J	s
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG_L	0.5	UJ	s
8260C	CYCLOHEXANE	110-82-7	UG_L	0.5	UJ	s
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG_L	0.5	UJ	s
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG_L	1	UJ	s,c
8260C	ETHYLBENZENE	100-41-4	UG_L	0.5	UJ	s
8260C	ISOPROPYLBENZENE	98-82-8	UG_L	0.5	UJ	s
8260C	M- AND P-XYLENE	108-38-3/106-42	UG_L	1	UJ	s
8260C	METHYL ACETATE	79-20-9	UG_L	0.75	UJ	s
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	UJ	s
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	UJ	s
8260C	METHYLENE CHLORIDE	75-09-2	UG_L	2.5	UJ	s
8260C	O-XYLENE	95-47-6	UG_L	0.5	UJ	s
8260C	STYRENE	100-42-5	UG_L	0.5	UJ	s
8260C	TETRACHLOROETHENE	127-18-4	UG_L	0.84	J	s
8260C	TOLUENE	108-88-3	UG_L	0.5	UJ	s
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	UJ	s
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	UJ	s
8260C	TRICHLOROETHENE	79-01-6	UG_L	640	J	s,m
8260C	TRICHLOROFLUOROMETHANE	75-69-4	UG_L	1	UJ	s
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	UJ	s
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	UJ	s
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	3.8		

**Notes:**

UG\_L = Micrograms per liter  
NA = Not analyzed  
Qual = Final qualifiers (See Attachment A)  
RC = Reason codes (See Attachment B)

**Appendix C**  
**Analytical Data Validation – ARCADIS**

## **Northrop Grumman Corporation- Operable Unit 2**

### **Data Review**

BETHPAGE, NEW YORK

Volatile and Semivolatile Analyses

SDGs #JC2751, JC2857 and JC2989

Analyses Performed By:  
Accutest Laboratories  
Dayton, New Jersey

Report #24352R  
September 29, 2015  
Review Level: Tier II  
Project #NY001496.1514.NAVI4



## SUMMARY

This data quality assessment summarizes the review of Sample Delivery Groups (SDGs) #JC2751, JC2857 and JC2989 for samples collected in association with the Northrop Grumman-Bethpage Site. The review was conducted as a Tier II evaluation and included review of data package completeness. Only analytical data associated with constituents of concern were reviewed for this validation. Field documentation was not included in this review. Included with this assessment are the validation annotated sample result sheets, and chain of custody. Analyses were performed on the following samples:

SDGs	Sample ID	Lab ID	Matrix	Sample Collection Date	Parent Sample	Analysis				
						VOC	SVOC	PCB	MET	MISC
JC2751	FB083115BT1	JC2751-1	Water	08/31/2015		X	X			
	TB083115BT1	JC2751-2	Water	08/31/2015		X				
	BPOW 5-1	JC2751-3	Water	08/31/2015		X	X			
JC2857	TB090115BT1	JC2857-1	Water	09/01/2015		X				
	FB090115BT1	JC2857-2	Water	09/01/2015		X	X			
	BPOW 5-2	JC2857-3	Water	09/01/2015		X	X			
JC2989	FB090215BT1	JC2989-1	Water	09/02/2015		X	X			
	TB090215BT1	JC2989-2	Water	09/02/2015		X				
	BPOW 5-3	JC2989-3	Water	09/02/2015		X	X			

Notes:

1. Semivolatile (1,4-Dioxane) analysis was performed at Accutest Laboratories, Marlborough, MA (Accutest-NE) facility.

## ANALYTICAL DATA PACKAGE DOCUMENTATION

### GENERAL INFORMATION

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

QA - Quality Assurance

## VOLATILE ANALYSIS INTRODUCTION

Analyses were performed according to United States Environmental Protection Agency (USEPA), Methods 524.2 and 8270D-Selected Ion Monitoring (SIM). Data were reviewed in accordance with USEPA National Functional Guidelines of October 1999.

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

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

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

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

## VOLATILE ORGANIC COMPOUND (VOC) ANALYSES

### 1. Holding Times

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

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

s.u. Standard units

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

### 2. Blank Contamination

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

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

Compounds were detected in the associated QA blanks; however, the associated sample results were greater than the BAL and/or were non-detect. No qualification of the sample results was required.

### 3. Surrogates/System Monitoring Compounds

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

All surrogate recoveries were within control limits.

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

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

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

A MS/MSD analysis was not performed on a sample location associated with these SDGs.

## **5. Laboratory Control Sample (LCS) Analysis**

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

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

## **6. Field Duplicate Analysis**

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

A field duplicate was not collected on a sample location associated with these SDGs.

## **7. Laboratory Duplicate Analysis**

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

A laboratory duplicate sample was not performed on a sample location associated with these SDGs.

## **8. System Performance and Overall Assessment**

Tentatively identified compounds (TICs) were not identified in any of the sample locations.

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

## DATA VALIDATION CHECKLIST FOR VOCs

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

%R Percent Recovery    RPD Relative Percent Difference

## SEMIVOLATILE ORGANIC COMPOUND (SVOC) ANALYSES

### 1. Holding Times

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

Method	Matrix	Holding Time	Preservation
SW-846 8270D SIM	Water	7 days from collection to extraction and 40 days from extraction to analysis	Cooled to <6°C

All samples were analyzed within the specified holding time criteria.

### 2. Blank Contamination

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

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

Compounds were not detected above the MDL in the associated blanks; therefore detected sample results were not associated with blank contamination.

### 3. Surrogates/System Monitoring Compounds

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

All surrogate recoveries were within control limits.

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

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

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

A MS/MSD analysis was not performed on a sample location associated with these SDGs.

## **5. Laboratory Control Sample/Laboratory Control Sample Duplicate (LCS/LCSD) Analysis**

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

All compounds associated with the LCS/LCSD analysis exhibited recoveries and RPD within the control limits.

## **6. Field Duplicate Analysis**

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

A field duplicate was not collected with a sample location associated with these SDGs.

## **7. System Performance and Overall Assessment**

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



## DATA VALIDATION CHECKLIST FOR SVOCs

SVOCs: SW-846 8270D-SIM	Reported		Performance Acceptable		Not Required
	No	Yes	No	Yes	
GAS CHROMATOGRAPHY/MASS SPECTROMETRY (GC/MS)					
<b>Tier II Validation</b>					
Holding times		X		X	
Reporting limits (units)		X		X	
Blanks					
A. Method blanks		X		X	
B. Equipment blanks		X		X	
Laboratory Control Sample (LCS) %R		X		X	
Laboratory Control Sample Duplicate(LCSD) %R		X		X	
LCS/LCSD Precision (RPD)		X		X	
Matrix Spike (MS) %R					X
Matrix Spike Duplicate(MSD) %R					X
MS/MSD Precision (RPD)					X
Field/Lab Duplicate (RPD)					X
Surrogate Spike Recoveries		X		X	
Dilution Factor		X		X	
Moisture Content					X

%R    Percent recovery                      RPD    Relative percent difference

VALIDATION PERFORMED BY: Lisa Horton

SIGNATURE:   
\_\_\_\_\_

DATE: September 29, 2015

PEER REVIEW BY: Todd Church

DATE: October 5, 2015

**CHAIN OF CUSTODY/  
CORRECTED SAMPLE ANALYSIS DATA SHEETS**



CW  
PB  
WIB

### CHAIN OF CUSTODY

Accutest New Jersey/SPL Environmental

2235 Route 130, Dayton, NJ 08810

TEL: 732-329-0200 FAX: 732-329-3499/3480

www.accutest.com

Client / Reporting Information		Project Information										Requested Analysis (see TEST CODE sheet)										Matrix Codes		
Company Name <b>Arcadis</b>		Project Name: <b>AGMNYM62235 // OU2 Outpost Wells Northrop Grumman OU2 Hydro</b>										FED-EX Tracking # <b>#4</b>										Bottle Order Control #		
Street Address <b>2 Huntington Quad, Suite 1S10</b>		Street <b>Bethpage NY</b>										Billing Information (if different from Report to)										Accutest Quote #		
City State Zip <b>Melville NY 11747</b>		City State <b>Bethpage NY</b>										Company Name <b>Arcadis, U.S., Inc. Attn: Accts Payable</b>										Accutest Job # <b>JC2751</b>		
Project Contact <b>Soma Das, soma.das@arcadis-us.com</b>		Project # <b>NY001496.314I.NAVIS</b>										Street Address <b>630 Plaza Drive, Suite 600</b>										DW - Drinking Water GW - Ground Water WW - Water SW - Surface Water SO - Soil SL - Sludge SED - Sediment OI - Oil LIQ - Other Liquid AIR - Air SOL - Other Solid WP - Wipe FB - Field Blank EB - Equipment Blank RB - Rinse Blank TB - Trip Blank		
Phone # <b>631-391-5247</b>		Client Purchase Order # <b>NY001496_2015</b>										City State Zip <b>Highlands Ranch, CO 80129</b>												
Fax # <b>631-249-7610</b>		Work Authorization # <b>NY001496_2015</b>										Attention: <b>Soma Das</b>												
Sampler(s) Name(s) <b>PATRICIA PRZORSKI, BRENDAN TONG</b>		Project Manager <b>Carlo San Giovanni</b>																						
Phone # <b>516-247-6247</b>																								
Turnaround Time (Business days)		Data Deliverable Information										Comments / Special Instructions										LAB USE ONLY		
<input type="checkbox"/> Std. 15 Business Days <input checked="" type="checkbox"/> Std. 10 Business Days (by Contract only) <input type="checkbox"/> 10 Day RUSH <input type="checkbox"/> 5 Day RUSH <input type="checkbox"/> 3 Day EMERGENCY <input type="checkbox"/> 2 Day EMERGENCY <input type="checkbox"/> 1 Day EMERGENCY Emergency & Rush T/A data available VIA Lablink		Approved By (Accutest PM): / Date: <input type="checkbox"/> Commercial "A" (Level 1) <input type="checkbox"/> Commercial "B" (Level 2) <input type="checkbox"/> FULLT1 (Level 3+4) <input type="checkbox"/> NJ Reduced <input type="checkbox"/> Commercial "C" Commercial "A" = Results Only Commercial "B" = Results + QC Summary NJ Reduced = Results + QC Summary + Partial Raw data										NYASP Category A NYASP Category B State Forms EDD Format <input checked="" type="checkbox"/> Other <b>COMMC+</b>										OU2 Hydro Analyte List (V5242NG14OW+40) plus 1,4-Dioxane (B8270SIM14DIOX) Please use "CP" for MS/MSD QA/QC Sample.		
Reinquished by Sampler: <b>1 Patricia Przeroski</b> <b>3 Brendan Tong</b> <b>5</b>		Date Time: <b>8/31/15 1900</b> Received By: <b>Chris Paul</b> <b>1</b> <b>3</b> <b>5</b>										Date Time: <b>9/1/15 1000</b> Received By: <b>Chris Paul</b> <b>2</b> <b>4</b>										Date Time: <b>9/1/15/1710</b> Received By: <b>JK</b> <b>2</b> <b>4</b>	Initial Assessment: <b>JK</b> Label Verification: <b>JK</b>	
Custody Seal #		Intact										Preserved where applicable										Cooler Temp. <b>2.7°C</b>		

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JC2751: Chain of Custody

Page 1 of 4

Client / Reporting Information		Project Information		Requested Analysis (see TEST CODE sheet)		Matrix Codes	
Company Name <b>Arcadis</b> Street Address <b>2 Huntington Quad, Suite 1S10</b> City State Zip <b>Melville NY 11747</b> Project Contact <b>Soma Das, soma.das@arcadis-us.com</b> Phone # Fax # <b>631-391-5247 631-249-7610</b> Same(s) Name(s) <b>PATRICIA PREZORSKI, BRENDAN TONG</b>		Project Name <b>AGMNYM62235 // OU2 Outpost Wells</b> <b>Northrop Grumman OU2 Hydro</b> Billing Information (if different from Report to) Company Name <b>Arcadis, U.S., Inc. Attn: Accts Payable</b> Street Address <b>630 Plaza Drive, Suite 600</b> City State Zip <b>Highlands Ranch, CO 80129</b> Client Purchase Order # <b>NY001496.314I.NAVI3</b> Work Authorization #: NY001496_2015 Project Manager <b>Carlo San Giovanni</b> Attention: <b>Soma Das</b>		Accutest Tracking # <b>#4</b> Accutest Quote # Bottle Order Contact # <b>JC2751</b>		DW - Drinking Water GW - Ground Water WW - Water SW - Surface Water SO - Soil SL - Sludge SED - Sediment OI - Oil LIQ - Other Liquid AIR - Air SOL - Other Solid WP - Wipe FB-Field Blank EB-Equipment Blank RB-Rinse Blank TB-Trip Blank	
Turnaround Time (Business days) <input type="checkbox"/> Std. 15 Business Days <input checked="" type="checkbox"/> Std. 10 Business Days (by Contract only) <input type="checkbox"/> 10 Day RUSH <input type="checkbox"/> 3 Day RUSH <input type="checkbox"/> 3 Day EMERGENCY <input type="checkbox"/> 2 Day EMERGENCY <input type="checkbox"/> 1 Day EMERGENCY Emergency & Rush TIA data available VIA Lablink		Approved By (Accutest PMP) / Date: _____ _____		Data Deliverable Information <input type="checkbox"/> Commercial "A" (Level 1) <input type="checkbox"/> NYASP Category A <input type="checkbox"/> Commercial "B" (Level 2) <input type="checkbox"/> NYASP Category B <input type="checkbox"/> FULLT1 (Level 3+4) <input type="checkbox"/> State Forms <input type="checkbox"/> NJ Reduced <input type="checkbox"/> EDD Format <input type="checkbox"/> Commercial "C" <input checked="" type="checkbox"/> Other <b>COMMC+</b> Commercial "A" = Results Only Commercial "B" = Results + QC Summary NJ Reduced = Results + QC Summary + Partial Raw data		Comments / Special Instructions <b>OU2 Hydro Analyte List (V5242NG14OW+40) plus 1,4-Dioxane (B8270SIM14DIOX)</b> Please use " <i>MS</i> " for MS/MSD QA/QC Sample.	
Retransmitted by Sample # <b>1 Patricia Prezorski</b> Date/Time: <b>8/15/15 16:00</b>		Received By: <b>FX</b> Date/Time: <b>9/1/15</b>		Retransmitted by: <b>FX</b> Date/Time: <b>9/1/15 9:30</b>		Received By: <b>Quincy</b>	
Retransmitted by Sample # <b>3 Brendan Tong</b> Date/Time: _____		Received By: _____ Date/Time: _____		Retransmitted by: _____ Date/Time: _____		Received By: _____	
Retransmitted by Sample # <b>5</b> Date/Time: _____		Received By: _____ Date/Time: _____		Retransmitted by: _____ Date/Time: _____		Received By: _____	

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## Report of Analysis

Client Sample ID: FB083115BT1	Date Sampled: 08/31/15
Lab Sample ID: JC2751-1	Date Received: 09/01/15
Matrix: AQ - Field Blank Water	Percent Solids: n/a
Method: EPA 524.2 REV 4.1	
Project: Northrop Grumman, OU2 Hydro, Bethpage, NY	

Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B99183.D	1	09/08/15	MD	n/a	n/a	V1B4703
Run #2							

Run #	Purge Volume
Run #1	5.0 ml
Run #2	

## Special VOA List

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	1.2	5.0	0.91	ug/l	J
78-93-3	2-Butanone	ND	5.0	0.57	ug/l	
71-43-2	Benzene	ND	0.50	0.057	ug/l	
75-27-4	Bromodichloromethane	ND	0.50	0.082	ug/l	
75-25-2	Bromoform	ND	0.50	0.046	ug/l	
74-83-9	Bromomethane	ND	0.50	0.077	ug/l	
75-15-0	Carbon disulfide	ND	0.50	0.028	ug/l	
108-90-7	Chlorobenzene	ND	0.50	0.027	ug/l	
75-00-3	Chloroethane	ND	0.50	0.037	ug/l	
67-66-3	Chloroform	ND	0.50	0.031	ug/l	
74-87-3	Chloromethane	ND	0.50	0.044	ug/l	
56-23-5	Carbon tetrachloride	ND	0.50	0.074	ug/l	
75-34-3	1,1-Dichloroethane	ND	0.50	0.039	ug/l	
75-35-4	1,1-Dichloroethylene	ND	0.50	0.054	ug/l	
107-06-2	1,2-Dichloroethane	ND	0.50	0.034	ug/l	
78-87-5	1,2-Dichloropropane	ND	0.50	0.082	ug/l	
124-48-1	Dibromochloromethane	ND	0.50	0.042	ug/l	
156-60-5	trans-1,2-Dichloroethylene	ND	0.50	0.039	ug/l	
156-59-2	cis-1,2-Dichloroethylene	ND	0.50	0.081	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.033	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.063	ug/l	
100-41-4	Ethylbenzene	ND	0.50	0.033	ug/l	
76-13-1	Freon 113	ND	1.0	0.10	ug/l	
591-78-6	2-Hexanone	ND	2.0	0.084	ug/l	
75-09-2	Methylene chloride	ND	0.50	0.047	ug/l	
108-10-1	4-Methyl-2-pentanone	ND	2.0	0.27	ug/l	
100-42-5	Styrene	ND	0.50	0.028	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	0.50	0.050	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.035	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	0.50	0.052	ug/l	
127-18-4	Tetrachloroethylene	ND	0.50	0.091	ug/l	
108-88-3	Toluene	ND	0.50	0.044	ug/l	

ND = Not detected

MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

## Report of Analysis

Client Sample ID: FB083115BT1		Date Sampled: 08/31/15
Lab Sample ID: JC2751-1		Date Received: 09/01/15
Matrix: AQ - Field Blank Water		Percent Solids: n/a
Method: EPA 524.2 REV 4.1		
Project: Northrop Grumman, OU2 Hydro, Bethpage, NY		

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	98%		78-114%
460-00-4	4-Bromofluorobenzene	91%		77-115%

CAS No.	Tentatively Identified Compounds	R.T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

---

ND = Not detected      MDL = Method Detection Limit      J = Indicates an estimated value  
 RL = Reporting Limit      B = Indicates analyte found in associated method blank  
 E = Indicates value exceeds calibration range      N = Indicates presumptive evidence of a compound

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## Report of Analysis

<b>Client Sample ID:</b> FB083115BT1 <b>Lab Sample ID:</b> JC2751-1 <b>Matrix:</b> AQ - Field Blank Water <b>Method:</b> SW846 8270D BY SIM SW846 3510C <b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY	<b>Date Sampled:</b> 08/31/15 <b>Date Received:</b> 09/01/15 <b>Percent Solids:</b> n/a
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Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 <sup>a</sup>	I97642.D	1	09/10/15	AMA	09/02/15	M:OP44462	M:MSI3652
Run #2							

Run #	Initial Volume	Final Volume
Run #1	980 ml	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
123-91-1	1,4-Dioxane	ND	0.20	0.078	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	72%		26-121%		
321-60-8	2-Fluorobiphenyl	66%		28-107%		
1718-51-0	Terphenyl-d14	86%		29-129%		

(a) Analysis performed at Accutest Laboratories, Marlborough, MA.

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ND = Not detected RL = Reporting Limit E = Indicates value exceeds calibration range	MDL = Method Detection Limit J = Indicates an estimated value B = Indicates analyte found in associated method blank N = Indicates presumptive evidence of a compound
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## Report of Analysis

Client Sample ID: TB083115BT1	Date Sampled: 08/31/15
Lab Sample ID: JC2751-2	Date Received: 09/01/15
Matrix: AQ - Trip Blank Water	Percent Solids: n/a
Method: EPA 524.2 REV 4.1	
Project: Northrop Grumman, OU2 Hydro, Bethpage, NY	

Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B99184.D	1	09/08/15	MD	n/a	n/a	V1B4703
Run #2							

Run #	Purge Volume
Run #1	5.0 ml
Run #2	

## Special VOA List

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

ND = Not detected

MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

## Report of Analysis

<b>Client Sample ID:</b> TB083115BT1		<b>Date Sampled:</b> 08/31/15
<b>Lab Sample ID:</b> JC2751-2		<b>Date Received:</b> 09/01/15
<b>Matrix:</b> AQ - Trip Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	98%		78-114%
460-00-4	4-Bromofluorobenzene	93%		77-115%

CAS No.	Tentatively Identified Compounds	R.T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

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ND = Not detected      MDL = Method Detection Limit      J = Indicates an estimated value  
 RL = Reporting Limit      B = Indicates analyte found in associated method blank  
 E = Indicates value exceeds calibration range      N = Indicates presumptive evidence of a compound

4.2  
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## Report of Analysis

Client Sample ID:	BPOW 5-1	Date Sampled:	08/31/15
Lab Sample ID:	JC2751-3	Date Received:	09/01/15
Matrix:	AQ - Ground Water	Percent Solids:	n/a
Method:	EPA 524.2 REV 4.1		
Project:	Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B99179.D	1	09/08/15	MD	n/a	n/a	V1B4703
Run #2							

Run #	Purge Volume
Run #1	5.0 ml
Run #2	

## Special VOA List

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

ND = Not detected

MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

## Report of Analysis

<b>Client Sample ID:</b> BPOW 5-1		<b>Date Sampled:</b> 08/31/15
<b>Lab Sample ID:</b> JC2751-3		<b>Date Received:</b> 09/01/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	99%		78-114%
460-00-4	4-Bromofluorobenzene	94%		77-115%

CAS No.	Tentatively Identified Compounds	R.T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

ND = Not detected      MDL = Method Detection Limit  
 RL = Reporting Limit  
 E = Indicates value exceeds calibration range

J = Indicates an estimated value  
 B = Indicates analyte found in associated method blank  
 N = Indicates presumptive evidence of a compound

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## Report of Analysis

<b>Client Sample ID:</b> BPOW 5-1 <b>Lab Sample ID:</b> JC2751-3 <b>Matrix:</b> AQ - Ground Water <b>Method:</b> SW846 8270D BY SIM SW846 3510C <b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY	<b>Date Sampled:</b> 08/31/15 <b>Date Received:</b> 09/01/15 <b>Percent Solids:</b> n/a
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Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 <sup>a</sup>	I97650.D	1	09/10/15	AMA	09/02/15	M:OP44462	M:MSI3652
Run #2							

Run #	Initial Volume	Final Volume
Run #1	950 ml	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
123-91-1	1,4-Dioxane	ND	0.21	0.080	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	79%		26-121%		
321-60-8	2-Fluorobiphenyl	70%		28-107%		
1718-51-0	Terphenyl-d14	82%		29-129%		

(a) Analysis performed at Accutest Laboratories, Marlborough, MA.

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ND = Not detected RL = Reporting Limit E = Indicates value exceeds calibration range	MDL = Method Detection Limit J = Indicates an estimated value B = Indicates analyte found in associated method blank N = Indicates presumptive evidence of a compound
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4.3  
4

GW  
FB  
WB

**CHAIN OF CUSTODY**

Accutest New Jersey/SPL Environmental  
2235 Route 130, Dayton, NJ 08810  
TEL: 732-329-0200 FAX: 732-329-3499/3480  
www.accutest.com

FED-EX Tracking #	#4	Bottle Order Control #	
Accutest Quote #		Accutest Job #	JC2857

Client / Reporting Information		Project Information		Requested Analysis ( see TEST CODE sheet)										Matrix Codes
<b>Company Name</b> Arcadis <b>Street Address</b> 2 Huntington Quad, Suite 1S10 <b>City State Zip</b> Melville NY 11747 <b>Project Contact</b> Soma Das, soma.das@arcadis-us.com <b>Phone #</b> 631-391-5247 <b>Fax #</b> 631-249-7610 <b>Sampler(s) Name(s)</b> PATRICIA PREZORSKI, BRENDAN TONG		<b>Project Name:</b> AGMNYM62235 // OU2 Outpost Wells Northrop Grumman OU2 Hydro <b>Street</b> <b>Billing Information ( If different from Report to)</b> <b>Company Name</b> Bethpage NY <b>Street Address</b> Arcadis, U.S., Inc. Attn: Accts Payable <b>Project #</b> NY001496.314I.NAVI3 <b>Client Purchase Order #</b> 630 Plaza Drive, Suite 600 <b>Work Authorization #:</b> NY001496_2015 <b>Project Manager</b> Carlo San Giovanni <b>Attention:</b> Soma Das		<b>Matrix Codes</b> DW - Drinking Water GW - Ground Water WW - Water SW - Surface Water SO - Soil SL - Sludge SED - Sediment OI - Oil LIQ - Other Liquid AIR - Air SOL - Other Solid WP - Wipe FB - Field Blank EB - Equipment Blank RB - Rinse Blank TB - Trip Blank										
<b>Turnaround Time ( Business days)</b> <input type="checkbox"/> Std. 15 Business Days <input checked="" type="checkbox"/> Std. 10 Business Days ( by Contract only) <input type="checkbox"/> 10 Day RUSH <input type="checkbox"/> 5 Day RUSH <input type="checkbox"/> 3 Day EMERGENCY <input type="checkbox"/> 2 Day EMERGENCY <input type="checkbox"/> 1 Day EMERGENCY Emergency & Rush T/A data available VIA Lablink		<b>Approved By (Accutest PM): I Date:</b> INITIAL ASSESSMENT An 31 LABEL VERIFICATION OB		<b>Data Deliverable Information</b> <input type="checkbox"/> Commercial "A" ( Level 1 ) <input type="checkbox"/> Commercial "B" ( Level 2 ) <input type="checkbox"/> FULLT1 ( Level 3+4 ) <input type="checkbox"/> NJ Reduced <input type="checkbox"/> Commercial "C" <input type="checkbox"/> NYASP Category A <input type="checkbox"/> NYASP Category B <input type="checkbox"/> State Forms <input type="checkbox"/> EDD Format <input checked="" type="checkbox"/> Other COMMC+ Commercial "A" = Results Only Commercial "B" = Results + QC Summary NJ Reduced = Results + QC Summary + Partial Raw data										<b>Comments / Special Instructions</b> OU2 Hydro Analyte List (V5242NG14OW+40) plus 1,4-Dioxane (B8270SIM14DIOX) Please use " " for MS/MSD QA/QC Sample. Full analyte list for 524,2
<b>Sample Custody must be documented below each time samples change possession, including courier delivery.</b>														
<b>Relinquished by Sampler:</b> 1 Patricia Prezorski <b>Date Time:</b> 9/11/15 2:00		<b>Received By:</b> 1 Chris Laed <b>Date Time:</b> 9/11/15 11:50		<b>Relinquished by:</b> 3 Brendan Tong <b>Date Time:</b> 		<b>Received By:</b> 3 <b>Date Time:</b> 		<b>Relinquished by:</b> 5 <b>Date Time:</b> 		<b>Received By:</b> 5 <b>Date Time:</b> 		<b>Custody Seal #</b> <input checked="" type="checkbox"/> Intact <input type="checkbox"/> Not Intact <b>Preserved where applicable</b> <input type="checkbox"/> <b>Cooler Temp.</b> 3.6°C		

JC2857: Chain of Custody

Page 1 of 3

## Report of Analysis

<b>Client Sample ID:</b> TB090115BT1		<b>Date Sampled:</b> 09/01/15
<b>Lab Sample ID:</b> JC2857-1		<b>Date Received:</b> 09/02/15
<b>Matrix:</b> AQ - Trip Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B99181.D	1	09/08/15	MD	n/a	n/a	V1B4703
Run #2							

Run #1	Purge Volume
Run #1	5.0 ml
Run #2	

## Special VOA List

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

ND = Not detected MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

## Report of Analysis

<b>Client Sample ID:</b> TB090115BT1		<b>Date Sampled:</b> 09/01/15
<b>Lab Sample ID:</b> JC2857-1		<b>Date Received:</b> 09/02/15
<b>Matrix:</b> AQ - Trip Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	99%		78-114%
460-00-4	4-Bromofluorobenzene	92%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

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ND = Not detected      MDL = Method Detection Limit      J = Indicates an estimated value  
 RL = Reporting Limit      B = Indicates analyte found in associated method blank  
 E = Indicates value exceeds calibration range      N = Indicates presumptive evidence of a compound

4.1  
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## Report of Analysis

<b>Client Sample ID:</b> FB090115BT1	<b>Date Sampled:</b> 09/01/15
<b>Lab Sample ID:</b> JC2857-2	<b>Date Received:</b> 09/02/15
<b>Matrix:</b> AQ - Field Blank Water	<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1	
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY	

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B99182.D	1	09/08/15	MD	n/a	n/a	V1B4703
Run #2							

Run #1	Purge Volume
Run #1	5.0 ml
Run #2	

## Special VOA List

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	1.6	5.0	0.91	ug/l	J
78-93-3	2-Butanone	ND	5.0	0.57	ug/l	
71-43-2	Benzene	ND	0.50	0.057	ug/l	
75-27-4	Bromodichloromethane	ND	0.50	0.082	ug/l	
75-25-2	Bromoform	ND	0.50	0.046	ug/l	
74-83-9	Bromomethane	ND	0.50	0.077	ug/l	
75-15-0	Carbon disulfide	ND	0.50	0.028	ug/l	
108-90-7	Chlorobenzene	ND	0.50	0.027	ug/l	
75-00-3	Chloroethane	ND	0.50	0.037	ug/l	
67-66-3	Chloroform	ND	0.50	0.031	ug/l	
74-87-3	Chloromethane	ND	0.50	0.044	ug/l	
56-23-5	Carbon tetrachloride	ND	0.50	0.074	ug/l	
75-34-3	1,1-Dichloroethane	ND	0.50	0.039	ug/l	
75-35-4	1,1-Dichloroethylene	ND	0.50	0.054	ug/l	
107-06-2	1,2-Dichloroethane	ND	0.50	0.034	ug/l	
78-87-5	1,2-Dichloropropane	ND	0.50	0.082	ug/l	
124-48-1	Dibromochloromethane	ND	0.50	0.042	ug/l	
156-60-5	trans-1,2-Dichloroethylene	ND	0.50	0.039	ug/l	
156-59-2	cis-1,2-Dichloroethylene	ND	0.50	0.081	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.033	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.063	ug/l	
100-41-4	Ethylbenzene	ND	0.50	0.033	ug/l	
76-13-1	Freon 113	ND	1.0	0.10	ug/l	
591-78-6	2-Hexanone	ND	2.0	0.084	ug/l	
75-09-2	Methylene chloride	ND	0.50	0.047	ug/l	
108-10-1	4-Methyl-2-pentanone	ND	2.0	0.27	ug/l	
100-42-5	Styrene	ND	0.50	0.028	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	0.50	0.050	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.035	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	0.50	0.052	ug/l	
127-18-4	Tetrachloroethylene	ND	0.50	0.091	ug/l	
108-88-3	Toluene	ND	0.50	0.044	ug/l	

ND = Not detected MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

## Report of Analysis

<b>Client Sample ID:</b> FB090115BT1		<b>Date Sampled:</b> 09/01/15
<b>Lab Sample ID:</b> JC2857-2		<b>Date Received:</b> 09/02/15
<b>Matrix:</b> AQ - Field Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	100%		78-114%
460-00-4	4-Bromofluorobenzene	92%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

ND = Not detected      MDL = Method Detection Limit  
 RL = Reporting Limit  
 E = Indicates value exceeds calibration range

J = Indicates an estimated value  
 B = Indicates analyte found in associated method blank  
 N = Indicates presumptive evidence of a compound

4.2  
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## Report of Analysis

<b>Client Sample ID:</b> FB090115BT1		<b>Date Sampled:</b> 09/01/15
<b>Lab Sample ID:</b> JC2857-2		<b>Date Received:</b> 09/02/15
<b>Matrix:</b> AQ - Field Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8270D BY SIM SW846 3510C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 <sup>a</sup>	I97654.D	1	09/10/15	AMA	09/03/15	M:OP44479	M:MSI3652
Run #2							

Run #	Initial Volume	Final Volume
Run #1	930 ml	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
123-91-1	1,4-Dioxane	ND	0.22	0.082	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	83%		26-121%		
321-60-8	2-Fluorobiphenyl	68%		28-107%		
1718-51-0	Terphenyl-d14	82%		29-129%		

(a) Analysis performed at Accutest Laboratories, Marlborough, MA.

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ND = Not detected	MDL = Method Detection Limit	J = Indicates an estimated value
RL = Reporting Limit		B = Indicates analyte found in associated method blank
E = Indicates value exceeds calibration range		N = Indicates presumptive evidence of a compound

4.2  
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**Report of Analysis**

<b>Client Sample ID:</b>	BPOW5-2	<b>Date Sampled:</b>	09/01/15
<b>Lab Sample ID:</b>	JC2857-3	<b>Date Received:</b>	09/02/15
<b>Matrix:</b>	AQ - Ground Water	<b>Percent Solids:</b>	n/a
<b>Method:</b>	EPA 524.2 REV 4.1		
<b>Project:</b>	Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B99180.D	1	09/08/15	MD	n/a	n/a	V1B4703
Run #2							

Run #1	Purge Volume
Run #1	5.0 ml
Run #2	

**Special VOA List**

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

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound

## Report of Analysis

<b>Client Sample ID:</b> BPOW5-2		<b>Date Sampled:</b> 09/01/15
<b>Lab Sample ID:</b> JC2857-3		<b>Date Received:</b> 09/02/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	98%		78-114%
460-00-4	4-Bromofluorobenzene	90%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

ND = Not detected      MDL = Method Detection Limit  
 RL = Reporting Limit  
 E = Indicates value exceeds calibration range

J = Indicates an estimated value  
 B = Indicates analyte found in associated method blank  
 N = Indicates presumptive evidence of a compound

4.3  
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## Report of Analysis

<b>Client Sample ID:</b> BPOW5-2		<b>Date Sampled:</b> 09/01/15
<b>Lab Sample ID:</b> JC2857-3		<b>Date Received:</b> 09/02/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8270D BY SIM SW846 3510C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 <sup>a</sup>	I97655.D	1	09/10/15	AMA	09/03/15	M:OP44479	M:MSI3652
Run #2							

Run #	Initial Volume	Final Volume
Run #1	950 ml	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
123-91-1	1,4-Dioxane	ND	0.21	0.080	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	82%		26-121%		
321-60-8	2-Fluorobiphenyl	64%		28-107%		
1718-51-0	Terphenyl-d14	85%		29-129%		

(a) Analysis performed at Accutest Laboratories, Marlborough, MA.

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ND = Not detected	MDL = Method Detection Limit	J = Indicates an estimated value
RL = Reporting Limit		B = Indicates analyte found in associated method blank
E = Indicates value exceeds calibration range		N = Indicates presumptive evidence of a compound

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GW  
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WIB

Client / Reporting Information		Project Information		Requested Analysis ( see TEST CODE sheet)		Matrix Codes	
Company Name <b>Arcadis</b>		Project Name: <b>AGMNYM62235 // OU2 Outpost Wells Northrop Grumman OU2 Hydro</b>		Requested Analysis ( see TEST CODE sheet)		Matrix Codes	
Street Address <b>2 Huntington Quad, Suite 1S10</b>		Street		Requested Analysis ( see TEST CODE sheet)		Matrix Codes	
City State Zip <b>Melville NY 11747</b>		City State <b>Bethpage NY</b>		Requested Analysis ( see TEST CODE sheet)		Matrix Codes	
Project Contact <b>Soma Das, soma.das@arcadis-us.com</b>		Project # <b>NY001496.314I.NAVI3</b>		Requested Analysis ( see TEST CODE sheet)		Matrix Codes	
Phone # <b>631-391-5247</b>		Client Purchase Order # <b>NY001496.314I.NAVI3</b>		Requested Analysis ( see TEST CODE sheet)		Matrix Codes	
Sample(s) Name(s) <b>PATRICIA PREZORSKI, BRENDAN TONG</b>		Work Authorization #: <b>NY001496_2015</b>		Requested Analysis ( see TEST CODE sheet)		Matrix Codes	
Phone # <b>516-249-7810</b>		Project Manager <b>Carlo San Giovanni</b>		Requested Analysis ( see TEST CODE sheet)		Matrix Codes	
Attention: <b>Soma Das</b>		City State Zip <b>Highlands Ranch, CO 80129</b>		Requested Analysis ( see TEST CODE sheet)		Matrix Codes	
Collection		Number of preserved Bottles		Requested Analysis ( see TEST CODE sheet)		Matrix Codes	
Account Sample #	Field ID / Point of Collection	MECH/DOI Val #	Date	Time	Sampled by	Matrix	# of bottles
1	FB090215BTL		9/21/15	0940	PP	FB	2
2	TB090215BTL		9/21/15	0930	PP	TB	2
3	BR&W 5-3		9/21/15	1305	PP	GW	3
Turnaround Time ( Business days)		Data Deliverable Information		Comments / Special Instructions			
<input type="checkbox"/> Std. 15 Business Days <input checked="" type="checkbox"/> Std. 10 Business Days ( by Contract only) <input type="checkbox"/> 10 Day RUSH <input type="checkbox"/> 5 Day RUSH <input type="checkbox"/> 3 Day EMERGENCY <input type="checkbox"/> 2 Day EMERGENCY <input type="checkbox"/> 1 Day EMERGENCY Emergency & Rush TIA data available VIA Lablink		Approved By (Accutest PM): / Date: INITIAL ASSESSMENT <b>QB OB</b> LABEL VERIFICATION <b>JK</b>		<input type="checkbox"/> Commercial "A" ( Level 1) <input type="checkbox"/> Commercial "B" ( Level 2) <input type="checkbox"/> FULLT1 ( Level 3+4) <input type="checkbox"/> NJ Reduced <input type="checkbox"/> Commercial "C" Commercial "A" = Results Only Commercial "B" = Results + QC Summary NJ Reduced = Results + QC Summary + Partial Raw data		OU2 Hydro Analyte List (V5242NG14OW+40) plus 1,4-Dioxane (B8270SIM14DIOX) Please use " <b>PP</b> " for MS/MSD QA/QC Sample. Full analyte list for 524.2 <b>1814</b>	
Emergency & Rush TIA data available VIA Lablink							
Chain of Custody must be documented below each time samples change possession, including courier delivery.							
Relinquished by Sampler:	Date Time:	Received By:	Date Time:	Relinquished By:	Date Time:	Received By:	Date Time:
1 Patricia Prezorski	9/21/15 19:15	PP	9/21/15 12:00	2 PP	9/21/15 18:11	3	9/21/15 18:11
3 Brendan Tong		3		4		4	
5		5					
Custody Seal #		<input type="checkbox"/> Intact <input type="checkbox"/> Not Intact		Preserved where applicable <input type="checkbox"/>		On Ice <input type="checkbox"/>	

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FED-EX Tracking # *934* Bottle Order Control # *JC2989*  
Accutest Order # *107* Accutest Job # *MC1127*

Client / Reporting Information		Project Information		Requested Analysis (see TEST CODE sheet)										Matrix Codes												
Company Name <b>Arcadis</b>		Project Name: <b>AGMNYM62235 // OU2 Outpost Wells Northrop Grumman OU2 Hydro</b>		<i>Full List</i> V5242NG440W448 B8270SIM14DIOX										DW - Drinking Water GW - Ground Water WW - Water SW - Surface Water SO - Soil SL - Sludge SED - Sediment O - Oil LIQ - Other Liquid AIR - Air SOL - Other Solid WP - Wipe FB - Field Blank EB - Equipment Blank RB - Rinse Blank TB - Trip Blank												
Street Address <b>2 Huntington Quad, Suite 1S10</b>		Street <b>Bethpage NY</b>																								
City State Zip <b>Melville NY 11747</b>		City State <b>Bethpage NY</b>																								
Project Contact <b>Soma Das, soma.das@arcadis-us.com</b>		Billing Information (if different from Report to) Company Name <b>Arcadis, U.S., Inc. Attn: Accts Payable</b>																								
Phone # <b>631-391-5247</b>		Project # <b>NY001496.3141.NAVI3</b>																								
Fax # <b>631-249-7810</b>		Client Purchase Order # <b>NY001496.2015</b>																								
Sample(s) Name(s) <b>PATRICIA PRZOROSKI, BRENDAN TONG</b>		Project Manager <b>Carlo San Giovanni</b>																								
Phone # <b>516-287-6442</b>		Attention: <b>Soma Das</b>																								
Accutest Sample #		Field ID / Point of Collection		MECHMDI Val #		Date		Time		Sampled by		Matrix		# of bottles		Number of preserved bottles										LAB USE ONLY
<i>14</i>		<i>FB090215B72 -1</i>				<i>9/2/15</i>		<i>0840</i>		<i>PP FB</i>		<i>2</i>												<i>-1</i>		
<i>15</i>		<i>BRW 5-3 -3</i>				<i>9/2/15</i>		<i>1305</i>		<i>PP GW</i>		<i>2</i>												<i>-3</i>		
																								<i>16D</i>		

Turnaround Time (Business days)		Data Deliverable Information				Comments / Special Instructions	
<input type="checkbox"/> Std. 15 Business Days <input checked="" type="checkbox"/> Std. 18 Business Days (by Contract only) <input type="checkbox"/> 10 Day RUSH <input type="checkbox"/> 5 Day RUSH <input type="checkbox"/> 3 Day EMERGENCY <input type="checkbox"/> 2 Day EMERGENCY <input type="checkbox"/> 1 Day EMERGENCY Emergency & Rush TIA data available VIA Lablink		Approved By (Account Mgr / Date): _____ _____ _____ _____ _____				Commercial "A" (Level 1) <input type="checkbox"/> Commercial "B" (Level 2) <input type="checkbox"/> FULLT1 (Level 3+4) <input type="checkbox"/> NJ Reduced <input type="checkbox"/> Commercial "C" NYASP Category A <input type="checkbox"/> NYASP Category B State Forms <input type="checkbox"/> EDD Format <input checked="" type="checkbox"/> Other <b>COMMC+</b> Commercial "A" = Results Only Commercial "B" = Results + QC Summary NJ Reduced = Results + QC Summary + Partial Raw data	
						OU2 Hydro Analysis (V5242NG440W448) (Full List) 1:4-Dioxane (B8270SIM14DIOX) Please use " <i>PP</i> " for MS/MSD QA/QC Sample.	
Requisitioned by: <i>Patricia Przoroski</i> Date/Time: <i>9/2/15 1700</i>		Received By: <i>FX</i> Date/Time: <i>9/2/15 9:30</i>		Requisitioned by: <i>FX</i> Date/Time: <i>9/2/15 9:30</i>		Received By: <i>Patricia Przoroski</i> Date/Time: <i>9/2/15 9:30</i>	
Requisitioned by: <i>Brendan Tong</i> Date/Time: _____		Received By: _____ Date/Time: _____		Requisitioned by: _____ Date/Time: _____		Received By: _____ Date/Time: _____	
Requisitioned by: _____ Date/Time: _____		Received By: _____ Date/Time: _____		Requisitioned by: _____ Date/Time: _____		Received By: _____ Date/Time: _____	
		Custody Seal # _____		<input type="checkbox"/> Intact <input type="checkbox"/> Not Intact		Preserved where applicable: _____ On Ice: <i>15</i> (F)	

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## Report of Analysis

<b>Client Sample ID:</b> FB090215BT1		
<b>Lab Sample ID:</b> JC2989-1		<b>Date Sampled:</b> 09/02/15
<b>Matrix:</b> AQ - Field Blank Water		<b>Date Received:</b> 09/03/15
<b>Method:</b> EPA 524.2 REV 4.1		<b>Percent Solids:</b> n/a
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B99193.D	1	09/08/15	MD	n/a	n/a	V1B4703
Run #2							

Run #1	Purge Volume
Run #1	5.0 ml
Run #2	

## Special VOA List

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

ND = Not detected

MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

## Report of Analysis

<b>Client Sample ID:</b> FB090215BT1		<b>Date Sampled:</b> 09/02/15
<b>Lab Sample ID:</b> JC2989-1		<b>Date Received:</b> 09/03/15
<b>Matrix:</b> AQ - Field Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	98%		78-114%
460-00-4	4-Bromofluorobenzene	91%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

ND = Not detected      MDL = Method Detection Limit  
 RL = Reporting Limit  
 E = Indicates value exceeds calibration range

J = Indicates an estimated value  
 B = Indicates analyte found in associated method blank  
 N = Indicates presumptive evidence of a compound

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## Report of Analysis

<b>Client Sample ID:</b> FB090215BT1	<b>Date Sampled:</b> 09/02/15
<b>Lab Sample ID:</b> JC2989-1	<b>Date Received:</b> 09/03/15
<b>Matrix:</b> AQ - Field Blank Water	<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8270D BY SIM SW846 3510C	
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY	

Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 <sup>a</sup>	I97759.D	1	09/16/15	AMA	09/08/15	M:OP44541	M:MSI3656
Run #2							

Run #	Initial Volume	Final Volume
Run #1	980 ml	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
123-91-1	1,4-Dioxane	ND	0.20	0.078	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	73%		26-121%		
321-60-8	2-Fluorobiphenyl	71%		28-107%		
1718-51-0	Terphenyl-d14	94%		29-129%		

(a) Analysis performed at Accutest Laboratories, Marlborough, MA.

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ND = Not detected	MDL = Method Detection Limit	J = Indicates an estimated value
RL = Reporting Limit		B = Indicates analyte found in associated method blank
E = Indicates value exceeds calibration range		N = Indicates presumptive evidence of a compound

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## Report of Analysis

<b>Client Sample ID:</b> TB090215BT1	<b>Date Sampled:</b> 09/02/15
<b>Lab Sample ID:</b> JC2989-2	<b>Date Received:</b> 09/03/15
<b>Matrix:</b> AQ - Trip Blank Water	<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1	
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY	

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B99194.D	1	09/08/15	MD	n/a	n/a	V1B4703
Run #2							

Run #1	Purge Volume
Run #1	5.0 ml
Run #2	

## Special VOA List

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

ND = Not detected MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

## Report of Analysis

<b>Client Sample ID:</b> TB090215BT1		<b>Date Sampled:</b> 09/02/15
<b>Lab Sample ID:</b> JC2989-2		<b>Date Received:</b> 09/03/15
<b>Matrix:</b> AQ - Trip Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	99%		78-114%
460-00-4	4-Bromofluorobenzene	90%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

ND = Not detected      MDL = Method Detection Limit  
 RL = Reporting Limit  
 E = Indicates value exceeds calibration range

J = Indicates an estimated value  
 B = Indicates analyte found in associated method blank  
 N = Indicates presumptive evidence of a compound

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## Report of Analysis

<b>Client Sample ID:</b>	BPOW 5-3	<b>Date Sampled:</b>	09/02/15
<b>Lab Sample ID:</b>	JC2989-3	<b>Date Received:</b>	09/03/15
<b>Matrix:</b>	AQ - Ground Water	<b>Percent Solids:</b>	n/a
<b>Method:</b>	EPA 524.2 REV 4.1		
<b>Project:</b>	Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B99202.D	1	09/09/15	MD	n/a	n/a	V1B4703
Run #2							

Run #1	Purge Volume
Run #1	5.0 ml
Run #2	

## Special VOA List

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

ND = Not detected MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

## Report of Analysis

<b>Client Sample ID:</b> BPOW 5-3		<b>Date Sampled:</b> 09/02/15
<b>Lab Sample ID:</b> JC2989-3		<b>Date Received:</b> 09/03/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	99%		78-114%
460-00-4	4-Bromofluorobenzene	90%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

ND = Not detected      MDL = Method Detection Limit  
 RL = Reporting Limit  
 E = Indicates value exceeds calibration range

J = Indicates an estimated value  
 B = Indicates analyte found in associated method blank  
 N = Indicates presumptive evidence of a compound

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## Report of Analysis

<b>Client Sample ID:</b> BPOW 5-3	<b>Date Sampled:</b> 09/02/15
<b>Lab Sample ID:</b> JC2989-3	<b>Date Received:</b> 09/03/15
<b>Matrix:</b> AQ - Ground Water	<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8270D BY SIM SW846 3510C	
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY	

Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 <sup>a</sup>	I97760.D	1	09/16/15	AMA	09/08/15	M:OP44541	M:MSI3656
Run #2							

Run #	Initial Volume	Final Volume
Run #1	960 ml	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
123-91-1	1,4-Dioxane	0.39	0.21	0.079	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	82%		26-121%		
321-60-8	2-Fluorobiphenyl	77%		28-107%		
1718-51-0	Terphenyl-d14	91%		29-129%		

(a) Analysis performed at Accutest Laboratories, Marlborough, MA.

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ND = Not detected	MDL = Method Detection Limit	J = Indicates an estimated value
RL = Reporting Limit		B = Indicates analyte found in associated method blank
E = Indicates value exceeds calibration range		N = Indicates presumptive evidence of a compound

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## **Northrop Grumman Corporation- Operable Unit 2**

### **Data Review**

BETHPAGE, NEW YORK

Volatile and Semivolatile Analyses

SDGs #JC3811, JC3916 and JC4092

Analyses Performed By:  
Accutest Laboratories  
Dayton, New Jersey

Report #24355R  
October 8, 2015  
Review Level: Tier II  
Project #NY001496.1514.NAVI4

## SUMMARY

This data quality assessment summarizes the review of Sample Delivery Groups (SDGs) #JC3811, JC3916 and JC4092 for samples collected in association with the Northrop Grumman-Bethpage Site. The review was conducted as a Tier II evaluation and included review of data package completeness. Only analytical data associated with constituents of concern were reviewed for this validation. Field documentation was not included in this review. Included with this assessment are the validation annotated sample result sheets, and chain of custody. Analyses were performed on the following samples:

SDGs	Sample ID	Lab ID	Matrix	Sample Collection Date	Parent Sample	Analysis				
						VOC	SVOC	PCB	MET	MISC
JC3811	TB091415BT1	JC3811-1	Water	09/14/2015		X				
	FB091415BT1	JC3811-2	Water	09/14/2015		X	X			
	BPOW 5-4	JC3811-3	Water	09/14/2015		X	X			
JC3916	TB091515BT1	JC3916-1	Water	09/15/2015		X				
	FB091515BT1	JC3916-2	Water	09/15/2015		X	X			
	BPOW 5-5	JC3916-3	Water	09/15/2015		X	X			
JC4092	TB091615BT1	JC4092-1	Water	09/16/2015		X				
	FB09165BT1	JC4092-2	Water	09/16/2015		X	X			
	BPOW 5-6	JC4092-3	Water	09/16/2015		X	X			

Notes:

1. Semivolatile (1,4-Dioxane) analysis was performed at Accutest Laboratories, Marlborough, MA (Accutest-NE) facility.
2. The matrix spike (MS) analysis was performed on sample location BPOW 5-5 for VOC and SVOC.

## ANALYTICAL DATA PACKAGE DOCUMENTATION

### GENERAL INFORMATION

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

QA - Quality Assurance

## VOLATILE ANALYSIS INTRODUCTION

Analyses were performed according to United States Environmental Protection Agency (USEPA), Methods 524.2 and 8270D-Selected Ion Monitoring (SIM). Data were reviewed in accordance with USEPA National Functional Guidelines of October 1999.

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

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

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

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

## VOLATILE ORGANIC COMPOUND (VOC) ANALYSES

### 1. Holding Times

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

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

s.u. Standard units

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

### 2. Blank Contamination

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

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

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

Sample Locations	Analytes	Sample Result	Qualification
BPOW 5-4	Acetone (Field Blank)	Detected sample results <RL and <BAL	"UB" at the RL
BPOW 5-5			
BPOW 5-6			

RL Reporting limit

### 3. Surrogates/System Monitoring Compounds

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

All surrogate recoveries were within control limits.

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

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

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

Sample locations associated with the MS/MSD exhibiting recoveries outside of the control limits are presented in the following table.

Sample Locations	Compound	MS Recovery	MSD Recovery
BPOW 5-5	2-Butanone	AC	>UL

AC Acceptable

The criteria used to evaluate the MS/MSD recoveries are presented in the following table. In the case of an MS/MSD deviation, the sample results are qualified as documented in the table below.

Control Limit	Sample Result	Qualification
> the upper control limit (UL)	Non-detect	No Action
	Detect	J
< the lower control limit (LL) but > 10%	Non-detect	UJ
	Detect	J
< 10%	Non-detect	R
	Detect	J
Parent sample concentration > four times the MS/MSD spiking solution concentration.	Detect	No Action
	Non-detect	

Sample locations associated with MS/MSD recoveries exhibiting an RPD greater than of the control limit presented in the following table.

Sample Locations	Compound
BPOW 5-5	1,1-Dichloroethane
	cis-1,3-Dichloropropene

The criteria used to evaluate the RPD between the MS/MSD recoveries are presented in the following table. In the case of an RPD deviation, the sample results are qualified as documented in the table below.

Control Limit	Sample Result	Qualification
> UL	Non-detect	UJ
	Detect	J

## **5. Laboratory Control Sample (LCS) Analysis**

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

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

## **6. Field Duplicate Analysis**

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

A field duplicate was not collected on a sample location associated with these SDGs.

## **7. Laboratory Duplicate Analysis**

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

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

## **8. System Performance and Overall Assessment**

Tentatively identified compounds (TICs) were not identified in any of the sample locations.

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

## DATA VALIDATION CHECKLIST FOR VOCs

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

%R Percent Recovery    RPD Relative Percent Difference



## SEMIVOLATILE ORGANIC COMPOUND (SVOC) ANALYSES

### 1. Holding Times

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

Method	Matrix	Holding Time	Preservation
SW-846 8270D SIM	Water	7 days from collection to extraction and 40 days from extraction to analysis	Cooled to <6°C

All samples were analyzed within the specified holding time criteria.

### 2. Blank Contamination

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

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

Compounds were not detected above the MDL in the associated blanks; therefore detected sample results were not associated with blank contamination.

### 3. Surrogates/System Monitoring Compounds

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

All surrogate recoveries were within control limits.

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

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

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

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

## **5. Laboratory Control Sample/Laboratory Control Sample Duplicate (LCS/LCSD) Analysis**

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

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

## **6. Field Duplicate Analysis**

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

A field duplicate was not collected with a sample location associated with these SDGs.

## **7. System Performance and Overall Assessment**

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

## DATA VALIDATION CHECKLIST FOR SVOCs

SVOCs: SW-846 8270D-SIM	Reported		Performance Acceptable		Not Required
	No	Yes	No	Yes	
GAS CHROMATOGRAPHY/MASS SPECTROMETRY (GC/MS)					
<b>Tier II Validation</b>					
Holding times		X		X	
Reporting limits (units)		X		X	
Blanks					
A. Method blanks		X		X	
B. Equipment blanks		X		X	
Laboratory Control Sample (LCS) %R		X		X	
Laboratory Control Sample Duplicate(LCSD) %R		X		X	
LCS/LCSD Precision (RPD)		X		X	
Matrix Spike (MS) %R		X		X	
Matrix Spike Duplicate(MSD) %R		X		X	
MS/MSD Precision (RPD)		X		X	
Field/Lab Duplicate (RPD)					X
Surrogate Spike Recoveries		X		X	
Dilution Factor		X		X	
Moisture Content					X

%R    Percent recovery                      RPD    Relative percent difference

VALIDATION PERFORMED BY: Lisa Horton

SIGNATURE:   
\_\_\_\_\_

DATE: October 8, 2015

PEER REVIEW BY: Todd Church

DATE: October 11, 2015

**CHAIN OF CUSTODY/  
CORRECTED SAMPLE ANALYSIS DATA SHEETS**

GW  
FB  
W/B

<b>FED-EX Tracking #</b> #4 <b>Accutest Quote #</b>		<b>Bottle Order Control #</b> <b>Accutest Job #</b> 503811															
<b>Client / Reporting Information</b> Company Name: <b>Arcadis</b> Street Address: <b>2 Huntington Quad, Suite 1S10</b> City: <b>Melville NY</b> Zip: <b>11747</b> Project Contact: <b>Soma Das, soma.das@arcadis-us.com</b> Phone #: <b>631-391-5247</b> Fax #: <b>631-249-7610</b> Sampler(s) Name(s): <b>PATRICIA PREZORSKI, BRENDAN TONG</b> Phone #: <b>516-247-6247</b>		<b>Project Information</b> Project Name: <b>AGMNYM62235 // OU2 Outpost Wells Northrop Grumman OU2 Hydro</b> Street: _____ Billing Information (if different from Report to): Company Name: <b>Bethpage NY</b> Street Address: <b>Arcadis, U.S., Inc. Attn: Accts Payable</b> Project #: <b>NY001496.1514.NAVI3</b> Client Purchase Order #: _____ Work Authorization #: <b>NY001496_2015</b> Project Manager: <b>Carlo San Giovanni</b> City: <b>Highlands Ranch, CO</b> State: <b>80129</b> Attention: <b>Soma Das</b>															
<b>Requested Analysis (see TEST CODE sheet)</b> <b>V5242NG FULL LIST</b> <b>B8270SIM14DIOX</b>		<b>Matrix Codes</b> DW - Drinking Water GW - Ground Water WW - Water SW - Surface Water SO - Soil SL - Sludge SED - Sediment OI - Oil LIQ - Other Liquid AIR - Air SOL - Other Solid WP - Wipe FB - Field Blank EB - Equipment Blank RB - Rinse Blank TB - Trip Blank															
<b>Collection</b> Accutest Sample #   Field ID / Point of Collection   MECHDI Val #   Date   Time   Sampled by   Matrix   # of bottles   HCl   HNO3   H2SO4   NONE   D/Water   MESH   ENCORE   LAB USE ONLY																	
1		TB091415BT2		9/14/15		1030		-		TB		2		2		✓	
2		FB091415BT2		9/14/15		1045		PP		FB		2		2		✓	
3		BPOW 5-4		9/14/15		1334		PP		GW		3		3		✓	
<b>Turnaround Time (Business days)</b> <input type="checkbox"/> Std. 15 Business Days <input checked="" type="checkbox"/> Std. 10 Business Days (by Contract only) <input type="checkbox"/> 10 Day RUSH <input type="checkbox"/> 5 Day RUSH <input type="checkbox"/> 3 Day EMERGENCY <input type="checkbox"/> 2 Day EMERGENCY <input type="checkbox"/> 1 Day EMERGENCY Emergency & Rush T/A data available VIA Lablink		<b>Approved By (Accutest PM): / Date:</b> _____ _____ _____		<b>Data Deliverable Information</b> <input type="checkbox"/> Commercial "A" (Level 1) <input type="checkbox"/> NYASP Category A <input type="checkbox"/> Commercial "B" (Level 2) <input type="checkbox"/> NYASP Category B <input type="checkbox"/> FULLT1 (Level 3+4) <input type="checkbox"/> State Forms <input type="checkbox"/> NJ Reduced <input type="checkbox"/> EDD Format <input type="checkbox"/> Commercial "C" <input checked="" type="checkbox"/> Other COMMC+ Commercial "A" = Results Only Commercial "B" = Results + QC Summary NJ Reduced = Results + QC Summary + Partial Raw data		<b>Comments / Special Instructions</b> OU2 Hydro Analyte List (V5242NG - FULL LIST) plus 1,4-Dioxane (B8270SIM14DIOX) Please use " <i>PP</i> " for MS/MSD QA/QC Sample. INITIAL ASSESSMENT <i>AC/PS</i> LABEL VERIFICATION <i>JK</i>											
<b>Sample Custody must be documented below each time samples change possession, including courier delivery.</b>																	
Relinquished by Sampler: <i>Patricia Prezorski</i> Date Time: <i>9/14/15 1745</i>		Received By: <i>Chris Paul</i> Date Time: <i>9/15/15 10:15</i>		Relinquished By: <i>Chris Paul</i> Date Time: <i>9/15/15 1650</i>		Received By: _____ Date Time: _____											
Relinquished by Sampler: _____ Date Time: _____		Received By: _____ Date Time: _____		Relinquished By: _____ Date Time: _____		Received By: _____ Date Time: _____											
Relinquished by: _____ Date Time: _____		Received By: _____ Date Time: _____		Custody Seal # _____ <input checked="" type="checkbox"/> Intact    Preserved where applicable <input type="checkbox"/> Not Intact <input type="checkbox"/>		On Ice <i>A</i> Cooler Temp. <i>1.3°C</i>											

5.1  
5



**CHAIN OF CUSTODY**  
 Accutest New Jersey/SPL Environmental  
 2235 Route 130, Dayton, NJ 08810  
 TEL: 732-329-0200 FAX: 732-329-3499/3480  
 www.accutest.com

FED-EX Tracking #	#4	Bole Order Control #																																																									
Accutest Quota #		Accutest Job #	JC3811																																																								
<b>Client / Reporting Information</b> Company Name: <b>Arcadis</b> Street Address: <b>2 Huntington Quad, Suite 1S10</b> City: <b>Melville NY</b> State: <b>NY</b> Zip: <b>11747</b>		<b>Project Information</b> Project Name: <b>AGMNYM62235 // OU2 Outpost Wells Northrop Grumman OU2 Hydro</b> Billing Information (if different from Report to): Company Name: <b>Arcadis, U.S., Inc. Attn: Accts Payable</b> Project #: <b>NY001496.1514.NAVI3</b> Street Address: <b>630 Plaza Drive, Suite 600</b> City: <b>Highlands Ranch, CO</b> State: <b>CO</b> Zip: <b>80129</b> Client Purchase Order #: <b>NY001496_2015</b> Work Authorization #: <b>NY001496_2015</b> Project Manager: <b>Carlo San Giovanni</b> Attention: <b>Soma Das</b>																																																									
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<table border="1"> <thead> <tr> <th>Accutest Sample #</th> <th>Field ID / Point of Collection</th> <th>MECHDI Vial #</th> <th>Date</th> <th>Time</th> <th>Sampled by</th> <th>Metric</th> <th># of bottles</th> <th>HCl</th> <th>NaOH</th> <th>HHO3</th> <th>H2SO4</th> <th>NO3</th> <th>DI Water</th> <th>USOP</th> <th>USOP2</th> <th>BRIDGE</th> <th>LAB USE ONLY</th> </tr> </thead> <tbody> <tr> <td>2</td> <td>FB091415B11</td> <td></td> <td>9/14/15</td> <td>10:45</td> <td>PD</td> <td>FB</td> <td>2</td> <td></td> <td></td> <td></td> <td></td> <td>3</td> <td></td> <td></td> <td></td> <td></td> <td></td> </tr> <tr> <td>3</td> <td>BPOL 5-4</td> <td></td> <td>9/14/15</td> <td>13:34</td> <td>PD</td> <td>GL</td> <td>2</td> <td></td> <td></td> <td></td> <td></td> <td>2</td> <td></td> <td></td> <td></td> <td></td> <td></td> </tr> </tbody> </table>	Accutest Sample #	Field ID / Point of Collection	MECHDI Vial #	Date	Time	Sampled by	Metric	# of bottles	HCl	NaOH	HHO3	H2SO4	NO3	DI Water	USOP	USOP2	BRIDGE	LAB USE ONLY	2	FB091415B11		9/14/15	10:45	PD	FB	2					3						3	BPOL 5-4		9/14/15	13:34	PD	GL	2					2						<b>Turnaround Time (Business days)</b> <input type="checkbox"/> Std. 15 Business Days <input checked="" type="checkbox"/> Std. 10 Business Days (by Contract only) <input type="checkbox"/> 10 Day RUSH <input type="checkbox"/> 5 Day RUSH <input type="checkbox"/> 3 Day EMERGENCY <input type="checkbox"/> 2 Day EMERGENCY <input type="checkbox"/> 1 Day EMERGENCY Emergency & Rush TIA data available VIA Lablink		<b>Data Deliverable Information</b> Approved By (Accutest PM): / Date: <input type="checkbox"/> Commercial "A" (Level 1) <input type="checkbox"/> Commercial "B" (Level 2) <input type="checkbox"/> FULLY1 (Level 3+4) <input type="checkbox"/> NJ Reduced <input type="checkbox"/> Commercial "C" <input type="checkbox"/> NYASP Category A <input type="checkbox"/> NYASP Category B <input type="checkbox"/> State Forms <input type="checkbox"/> EDD Format <input checked="" type="checkbox"/> Other <b>COMM+</b> Commercial "A" = Results Only Commercial "B" = Results + QC Summary NJ Reduced = Results + QC Summary + Partial Raw data		<b>Comments / Special Instructions</b> OU2 Hydro-Analyte List (V5242NG-FUEL LIST) plus 3,4-Dioxane (B8270SIM14DIOX) Please use "B" for MSMSD QA/QC Sample.
Accutest Sample #	Field ID / Point of Collection	MECHDI Vial #	Date	Time	Sampled by	Metric	# of bottles	HCl	NaOH	HHO3	H2SO4	NO3	DI Water	USOP	USOP2	BRIDGE	LAB USE ONLY																																										
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<b>Relinquished by Sampler:</b> Patricia Prezorski Date Time: 9/14/15 17:45 Received By: Date Time: Relinquished By: Date Time: Received By:																																																											
<b>Relinquished by Sampler:</b> Brendan Tong Date Time: 7-15-15 Received By: Date Time: Relinquished By: Date Time: Received By:																																																											
<b>Relinquished by:</b> Date Time: 7-15-15 Received By: Date Time: Relinquished By: Date Time: Received By:																																																											

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## Report of Analysis

<b>Client Sample ID:</b> TB091415BT1		<b>Date Sampled:</b> 09/14/15
<b>Lab Sample ID:</b> JC3811-1		<b>Date Received:</b> 09/15/15
<b>Matrix:</b> AQ - Trip Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B99424.D	1	09/17/15	MD	n/a	n/a	V1B4712
Run #2							

Run #1	Purge Volume
Run #1	5.0 ml
Run #2	

## Special VOA List

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

ND = Not detected MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound



## Report of Analysis

<b>Client Sample ID:</b> TB091415BT1		<b>Date Sampled:</b> 09/14/15
<b>Lab Sample ID:</b> JC3811-1		<b>Date Received:</b> 09/15/15
<b>Matrix:</b> AQ - Trip Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	104%		78-114%
460-00-4	4-Bromofluorobenzene	86%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

ND = Not detected      MDL = Method Detection Limit  
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 E = Indicates value exceeds calibration range

J = Indicates an estimated value  
 B = Indicates analyte found in associated method blank  
 N = Indicates presumptive evidence of a compound

4.1  
4

## Report of Analysis

<b>Client Sample ID:</b> FB091415BT1		
<b>Lab Sample ID:</b> JC3811-2		<b>Date Sampled:</b> 09/14/15
<b>Matrix:</b> AQ - Field Blank Water		<b>Date Received:</b> 09/15/15
<b>Method:</b> EPA 524.2 REV 4.1		<b>Percent Solids:</b> n/a
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B99425.D	1	09/17/15	MD	n/a	n/a	V1B4712
Run #2							

Run #1	Purge Volume
Run #1	5.0 ml
Run #2	

## Special VOA List

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

ND = Not detected

MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

## Report of Analysis

<b>Client Sample ID:</b> FB091415BT1	<b>Date Sampled:</b> 09/14/15
<b>Lab Sample ID:</b> JC3811-2	<b>Date Received:</b> 09/15/15
<b>Matrix:</b> AQ - Field Blank Water	<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1	
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY	

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	101%		78-114%
460-00-4	4-Bromofluorobenzene	86%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

ND = Not detected      MDL = Method Detection Limit  
 RL = Reporting Limit  
 E = Indicates value exceeds calibration range

J = Indicates an estimated value  
 B = Indicates analyte found in associated method blank  
 N = Indicates presumptive evidence of a compound

4.2  
4

## Report of Analysis

<b>Client Sample ID:</b> FB091415BT1		<b>Date Sampled:</b> 09/14/15
<b>Lab Sample ID:</b> JC3811-2		<b>Date Received:</b> 09/15/15
<b>Matrix:</b> AQ - Field Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8270D BY SIM SW846 3510C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 <sup>a</sup>	I97933.D	1	09/24/15	AMA	09/17/15	M:OP44660	M:MSI3662
Run #2							

Run #	Initial Volume	Final Volume
Run #1	900 ml	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
123-91-1	1,4-Dioxane	ND	0.22	0.084	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	77%		26-121%		
321-60-8	2-Fluorobiphenyl	67%		28-107%		
1718-51-0	Terphenyl-d14	89%		29-129%		

(a) Analysis performed at Accutest Laboratories, Marlborough, MA.

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ND = Not detected	MDL = Method Detection Limit	J = Indicates an estimated value
RL = Reporting Limit		B = Indicates analyte found in associated method blank
E = Indicates value exceeds calibration range		N = Indicates presumptive evidence of a compound

4.2  
4

## Report of Analysis

<b>Client Sample ID:</b> BPOW 5-4		<b>Date Sampled:</b> 09/14/15
<b>Lab Sample ID:</b> JC3811-3		<b>Date Received:</b> 09/15/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #2	1B99426.D	1	09/17/15	MD	n/a	n/a	V1B4712

Run #1	Purge Volume
Run #2	5.0 ml

## Special VOA List

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

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound

## Report of Analysis

<b>Client Sample ID:</b> BPOW 5-4		<b>Date Sampled:</b> 09/14/15
<b>Lab Sample ID:</b> JC3811-3		<b>Date Received:</b> 09/15/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	102%		78-114%
460-00-4	4-Bromofluorobenzene	84%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

ND = Not detected      MDL = Method Detection Limit  
 RL = Reporting Limit  
 E = Indicates value exceeds calibration range

J = Indicates an estimated value  
 B = Indicates analyte found in associated method blank  
 N = Indicates presumptive evidence of a compound

4.3  
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## Report of Analysis

<b>Client Sample ID:</b> BPOW 5-4	<b>Date Sampled:</b> 09/14/15
<b>Lab Sample ID:</b> JC3811-3	<b>Date Received:</b> 09/15/15
<b>Matrix:</b> AQ - Ground Water	<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8270D BY SIM SW846 3510C	
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY	

Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 <sup>a</sup>	I97934.D	1	09/24/15	AMA	09/17/15	M:OP44660	M:MSI3662
Run #2							

Run #	Initial Volume	Final Volume
Run #1	920 ml	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
123-91-1	1,4-Dioxane	0.28	0.22	0.083	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	77%		26-121%		
321-60-8	2-Fluorobiphenyl	75%		28-107%		
1718-51-0	Terphenyl-d14	89%		29-129%		

(a) Analysis performed at Accutest Laboratories, Marlborough, MA.

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ND = Not detected	MDL = Method Detection Limit	J = Indicates an estimated value
RL = Reporting Limit		B = Indicates analyte found in associated method blank
E = Indicates value exceeds calibration range		N = Indicates presumptive evidence of a compound

4.3  
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W  
FB  
WJB

# CHAIN OF CUSTODY

Accutest New Jersey/SPL Environmental  
2235 Route 130, Dayton, NJ 08810  
TEL: 732-329-0200 FAX: 732-329-3499/3480  
www.accutest.com

FED-EX Tracking # <b>#4</b>		Bottle Order Control #	
Accutest Quote #		Accutest Job # <b>JC3916</b>	

Client / Reporting Information		Project Information		Requested Analysis ( see TEST CODE sheet)		Matrix Codes	
Company Name <b>Arcadis</b>		Project Name: <b>AGMNYM62235 // OU2 Outpost Wells Northrop Grumman OU2 Hydro</b>		<b>V5242NG FULL LIST</b> <b>B8270SIM14DIOX</b>		DW - Drinking Water GW - Ground Water WW - Water SW - Surface Water SO - Soil SL - Sludge SED - Sediment OI - Oil LIQ - Other Liquid AIR - Air SOL - Other Solid WP - Wipe FB-Field Blank EB-Equipment Blank RB - Rinse Blank TB-Trip Blank	
Street Address <b>2 Huntington Quad, Suite 1S10</b>		Street <b>Bethpage NY</b>					
City State Zip <b>Melville NY 11747</b>		Billing Information ( If different from Report to) Company Name <b>Arcadis, U.S., Inc. Attn: Accts Payable</b>					
Project Contact <b>Soma Das, soma.das@arcadis-us.com</b>		Project # <b>NY001496.1514.NAVI3</b>					
Phone # Fax # <b>631-391-5247 631-249-7610</b>		Client Purchase Order # <b>Work Authorization #: NY001496_2015</b>					
Sampler(s) Name(s) <b>PATRICIA PREZORSKI, BRENDAN TONG</b>		City State Zip <b>Highlands Ranch, CO 80129</b>					
Project Manager <b>Carlo San Giovanni</b>		Attention: <b>Soma Das</b>					
Turnaround Time ( Business days) <input type="checkbox"/> Std. 15 Business Days <input checked="" type="checkbox"/> Std. 10 Business Days ( by Contract only) <input type="checkbox"/> 10 Day RUSH <input type="checkbox"/> 5 Day RUSH <input type="checkbox"/> 3 Day EMERGENCY <input type="checkbox"/> 2 Day EMERGENCY <input type="checkbox"/> 1 Day EMERGENCY Emergency & Rush TIA data available VIA Lablink		Data Deliverable Information <input type="checkbox"/> Commercial "A" ( Level 1) <input type="checkbox"/> Commercial "B" ( Level 2) <input type="checkbox"/> FULLT1 ( Level 3+4) <input type="checkbox"/> NJ Reduced <input type="checkbox"/> Commercial "C" Commercial "A" = Results Only Commercial "B" = Results + QC Summary NJ Reduced = Results + QC Summary + Partial Raw data					
Approved By (Accutest PM): / Date: <b>INITIAL ASSESSMENT HLL 3/3</b> <b>LABEL VERIFICATION Am</b>		Comments / Special Instructions <b>OU2 Hydro Analyte List (V5242NG -FULL LIST) plus 1,4-Dioxane (B8270SIM14DIOX)</b> Please use " <b>BPOW5-5</b> " for MS/MSD QA/QC Sample.					
Requisitioned by Sampler: <b>1 Patricia Prezorski</b>		Received By: <b>Chris Paul 9/15/15</b>					
Requisitioned by Sampler: <b>3 Brendan Tong</b>		Received By: <b>Chris Paul</b>					
Requisitioned by: <b>5</b>		Received By: <b>4</b>					

Account Sample #	Field ID / Point of Collection	MECH/DOI Vial #	Date	Time	Sampled by	Matrix	# of bottles	HCI	NH3	PHOS	PHOS3	H2SO4	NH4	DI Water	MEDI	ENCDRE	LAB USE ONLY
1	TB091515B72		9/15/15	1000	TB		2										
2	FB091515B72		9/15/15	1040	FB		2										V82
3	BPOW 5-5		9/15/15	1328	GW		9										

Custody Seal #	<input type="checkbox"/> Intact <input checked="" type="checkbox"/> Not Intact	Preserved where applicable <input type="checkbox"/>	On ice <input checked="" type="checkbox"/>	Cooler Temp. <b>2.6°C</b>
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JC3916: Chain of Custody

Page 1 of 3



Client / Reporting Information		Project Information		Requested Analysis ( see TEST CODE sheet)		Matrix Codes	
Company Name <b>Arcadis</b>		Project Name: <b>AGMNYM62235 // OU2 Outpost Wells</b>		FED-EX Tracking # <b>#4</b>		Bottle Order Control # <b>JC3916</b>	
Street Address <b>2 Huntington Quad, Suite 1S10</b>		Billing Information ( if different from Report to)		Accutest Quote #		Accutest Job # <b>JC3916</b>	
City <b>Melville NY</b>	State <b>NY</b>	Street <b>Bethpage NY</b>	City <b>NY</b>	Company Name <b>Arcadis, U.S., Inc. Attn: Accts Payable</b>		DW - Drinking Water GW - Ground Water WW - Water SW - Surface Water SO - Soil SL - Sludge SED - Sediment OI - Oil LIQ - Other Liquid AIR - Air SOL - Other Solid WP - Wipe FB - Field Blank EB - Equipment Blank RB - Rise Blank TB - Trip Blank	
Project Contact <b>Soma Das, soma.das@arcadis-us.com</b>	E-mail	Project # <b>NY001496.1514.NAVI3</b>	Company Name <b>630 Plaza Drive, Suite 600</b>	Street Address		LAB USE ONLY	
Phone # <b>631-391-5247</b>	Fax # <b>631-249-7610</b>	Client Purchase Order # <b>NY001496_2015</b>	City <b>Highlands Ranch, CO</b>	State <b>CO</b>	Zip <b>80129</b>		
Sampler(s) Name(s) <b>PATRICIA PRZORSKI, BRENDAN TONG</b>		Project Manager <b>Carlo San Giovanni</b>	Attention: <b>Soma Das</b>				
Turnaround Time ( Business days)		Data Deliverable Information		Comments / Special Instructions			
<input type="checkbox"/> Std. 15 Business Days <input checked="" type="checkbox"/> Std. 10 Business Days ( by Contract only) <input type="checkbox"/> 10 Day RUSH <input type="checkbox"/> 5 Day RUSH <input type="checkbox"/> 3 Day EMERGENCY <input type="checkbox"/> 2 Day EMERGENCY <input type="checkbox"/> 1 Day EMERGENCY Emergency & Rush TIA data available via Lablink		Approved By (Accutest PM) / Date: _____ _____		<input type="checkbox"/> Commercial "A" (Level 1) <input type="checkbox"/> Commercial "B" (Level 2) <input type="checkbox"/> FULLT1 (Level 3+4) <input type="checkbox"/> NJ Reduced <input type="checkbox"/> Commercial "C" Commercial "A" = Results Only Commercial "B" = Results + QC Summary NJ Reduced = Results + QC Summary + Partial Raw data		OU2 Hydro Analyte List (V5242NG-FULL LIST) plus 1,4-Dioxane (B8270SIM14DIOX) Please use "B8270SIM14DIOX" for MS/MSD QA/QC Sample.	
Sample Custody must be documented below each time samples change possession, including courier delivery.							
Relinquished by Sampler: <b>1 Patricia Przorski</b>	Date Time: <b>9/15/15 1830</b>	Received By: <b>F. Przek</b>	Received By: <b>F. Przek</b>	Relinquished By: <b>7</b>	Date Time: <b>9/16/15</b>	Received By: <b>7</b>	Received By: <b>7</b>
Relinquished by Sampler: <b>3 Brendan Tong</b>	Date Time:	Received By:	Received By:	Relinquished By:	Date Time:	Received By:	Received By:
Relinquished by: <b>5</b>	Date Time:	Received By:	Received By:	Custody Seal #	Intact <input type="checkbox"/> Hot Intact <input type="checkbox"/>	Preserved where applicable <input type="checkbox"/>	On Ice <input checked="" type="checkbox"/> Cooler Temp <b>0.4, 7.3 °C</b>

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## Report of Analysis

<b>Client Sample ID:</b> TB091515BT1	<b>Date Sampled:</b> 09/15/15
<b>Lab Sample ID:</b> JC3916-1	<b>Date Received:</b> 09/16/15
<b>Matrix:</b> AQ - Trip Blank Water	<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1	
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY	

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B99416.D	1	09/17/15	MD	n/a	n/a	V1B4712
Run #2							

Run #1	Purge Volume
Run #1	5.0 ml
Run #2	

## Special VOA List

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	0.94	5.0	0.91	ug/l	J
78-93-3	2-Butanone	ND	5.0	0.57	ug/l	
71-43-2	Benzene	ND	0.50	0.057	ug/l	
75-27-4	Bromodichloromethane	ND	0.50	0.082	ug/l	
75-25-2	Bromoform	ND	0.50	0.046	ug/l	
74-83-9	Bromomethane	ND	0.50	0.077	ug/l	
75-15-0	Carbon disulfide	ND	0.50	0.028	ug/l	
108-90-7	Chlorobenzene	ND	0.50	0.027	ug/l	
75-00-3	Chloroethane	ND	0.50	0.037	ug/l	
67-66-3	Chloroform	ND	0.50	0.031	ug/l	
74-87-3	Chloromethane	ND	0.50	0.044	ug/l	
56-23-5	Carbon tetrachloride	ND	0.50	0.074	ug/l	
75-34-3	1,1-Dichloroethane	ND	0.50	0.039	ug/l	
75-35-4	1,1-Dichloroethylene	ND	0.50	0.054	ug/l	
107-06-2	1,2-Dichloroethane	ND	0.50	0.034	ug/l	
78-87-5	1,2-Dichloropropane	ND	0.50	0.082	ug/l	
124-48-1	Dibromochloromethane	ND	0.50	0.042	ug/l	
156-60-5	trans-1,2-Dichloroethylene	ND	0.50	0.039	ug/l	
156-59-2	cis-1,2-Dichloroethylene	ND	0.50	0.081	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.033	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.063	ug/l	
100-41-4	Ethylbenzene	ND	0.50	0.033	ug/l	
76-13-1	Freon 113	ND	1.0	0.10	ug/l	
591-78-6	2-Hexanone	ND	2.0	0.084	ug/l	
75-09-2	Methylene chloride	ND	0.50	0.047	ug/l	
108-10-1	4-Methyl-2-pentanone	ND	2.0	0.27	ug/l	
100-42-5	Styrene	ND	0.50	0.028	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	0.50	0.050	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.035	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	0.50	0.052	ug/l	
127-18-4	Tetrachloroethylene	ND	0.50	0.091	ug/l	
108-88-3	Toluene	ND	0.50	0.044	ug/l	

ND = Not detected

MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

## Report of Analysis

<b>Client Sample ID:</b> TB091515BT1		<b>Date Sampled:</b> 09/15/15
<b>Lab Sample ID:</b> JC3916-1		<b>Date Received:</b> 09/16/15
<b>Matrix:</b> AQ - Trip Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	100%		78-114%
460-00-4	4-Bromofluorobenzene	85%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

ND = Not detected      MDL = Method Detection Limit  
 RL = Reporting Limit  
 E = Indicates value exceeds calibration range

J = Indicates an estimated value  
 B = Indicates analyte found in associated method blank  
 N = Indicates presumptive evidence of a compound

4.1  
4

## Report of Analysis

<b>Client Sample ID:</b> FB091515BT1		
<b>Lab Sample ID:</b> JC3916-2		<b>Date Sampled:</b> 09/15/15
<b>Matrix:</b> AQ - Field Blank Water		<b>Date Received:</b> 09/16/15
<b>Method:</b> EPA 524.2 REV 4.1		<b>Percent Solids:</b> n/a
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B99417.D	1	09/17/15	MD	n/a	n/a	V1B4712
Run #2							

Run #1	Purge Volume
Run #1	5.0 ml
Run #2	

## Special VOA List

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	1.2	5.0	0.91	ug/l	J
78-93-3	2-Butanone	ND	5.0	0.57	ug/l	
71-43-2	Benzene	ND	0.50	0.057	ug/l	
75-27-4	Bromodichloromethane	ND	0.50	0.082	ug/l	
75-25-2	Bromoform	ND	0.50	0.046	ug/l	
74-83-9	Bromomethane	ND	0.50	0.077	ug/l	
75-15-0	Carbon disulfide	ND	0.50	0.028	ug/l	
108-90-7	Chlorobenzene	ND	0.50	0.027	ug/l	
75-00-3	Chloroethane	ND	0.50	0.037	ug/l	
67-66-3	Chloroform	ND	0.50	0.031	ug/l	
74-87-3	Chloromethane	ND	0.50	0.044	ug/l	
56-23-5	Carbon tetrachloride	ND	0.50	0.074	ug/l	
75-34-3	1,1-Dichloroethane	ND	0.50	0.039	ug/l	
75-35-4	1,1-Dichloroethylene	ND	0.50	0.054	ug/l	
107-06-2	1,2-Dichloroethane	ND	0.50	0.034	ug/l	
78-87-5	1,2-Dichloropropane	ND	0.50	0.082	ug/l	
124-48-1	Dibromochloromethane	ND	0.50	0.042	ug/l	
156-60-5	trans-1,2-Dichloroethylene	ND	0.50	0.039	ug/l	
156-59-2	cis-1,2-Dichloroethylene	ND	0.50	0.081	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.033	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.063	ug/l	
100-41-4	Ethylbenzene	ND	0.50	0.033	ug/l	
76-13-1	Freon 113	ND	1.0	0.10	ug/l	
591-78-6	2-Hexanone	ND	2.0	0.084	ug/l	
75-09-2	Methylene chloride	ND	0.50	0.047	ug/l	
108-10-1	4-Methyl-2-pentanone	ND	2.0	0.27	ug/l	
100-42-5	Styrene	ND	0.50	0.028	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	0.50	0.050	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.035	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	0.50	0.052	ug/l	
127-18-4	Tetrachloroethylene	ND	0.50	0.091	ug/l	
108-88-3	Toluene	ND	0.50	0.044	ug/l	

ND = Not detected

MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

## Report of Analysis

<b>Client Sample ID:</b> FB091515BT1		<b>Date Sampled:</b> 09/15/15
<b>Lab Sample ID:</b> JC3916-2		<b>Date Received:</b> 09/16/15
<b>Matrix:</b> AQ - Field Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	102%		78-114%
460-00-4	4-Bromofluorobenzene	86%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

ND = Not detected      MDL = Method Detection Limit  
 RL = Reporting Limit  
 E = Indicates value exceeds calibration range

J = Indicates an estimated value  
 B = Indicates analyte found in associated method blank  
 N = Indicates presumptive evidence of a compound

4.2  
4

## Report of Analysis

<b>Client Sample ID:</b> FB091515BT1		<b>Date Sampled:</b> 09/15/15
<b>Lab Sample ID:</b> JC3916-2		<b>Date Received:</b> 09/16/15
<b>Matrix:</b> AQ - Field Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8270D BY SIM SW846 3510C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 <sup>a</sup>	I97911.D	1	09/23/15	AMA	09/19/15	M:OP44675	M:MSI3661
Run #2							

	Initial Volume	Final Volume
Run #1	820 ml	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
123-91-1	1,4-Dioxane	ND	0.24	0.093	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	69%		26-121%		
321-60-8	2-Fluorobiphenyl	64%		28-107%		
1718-51-0	Terphenyl-d14	74%		29-129%		

(a) Analysis performed at Accutest Laboratories, Marlborough, MA.

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ND = Not detected	MDL = Method Detection Limit	J = Indicates an estimated value
RL = Reporting Limit		B = Indicates analyte found in associated method blank
E = Indicates value exceeds calibration range		N = Indicates presumptive evidence of a compound

4.2  
4

## Report of Analysis

<b>Client Sample ID:</b> BPOW 5-5		<b>Date Sampled:</b> 09/15/15
<b>Lab Sample ID:</b> JC3916-3		<b>Date Received:</b> 09/16/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B99414.D	1	09/17/15	MD	n/a	n/a	V1B4712
Run #2							

Run #1	Purge Volume
Run #1	5.0 ml
Run #2	

## Special VOA List

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	5.0 <del>2.2</del>	5.0	0.91	ug/l	<del>J</del> UB
78-93-3	2-Butanone	ND	5.0	0.57	ug/l	
71-43-2	Benzene	ND	0.50	0.057	ug/l	
75-27-4	Bromodichloromethane	ND	0.50	0.082	ug/l	
75-25-2	Bromoform	ND	0.50	0.046	ug/l	
74-83-9	Bromomethane	ND	0.50	0.077	ug/l	
75-15-0	Carbon disulfide	ND	0.50	0.028	ug/l	
108-90-7	Chlorobenzene	ND	0.50	0.027	ug/l	
75-00-3	Chloroethane	ND	0.50	0.037	ug/l	
67-66-3	Chloroform	ND	0.50	0.031	ug/l	
74-87-3	Chloromethane	ND	0.50	0.044	ug/l	
56-23-5	Carbon tetrachloride	ND	0.50	0.074	ug/l	
75-34-3	1,1-Dichloroethane	ND	0.50	0.039	ug/l	J
75-35-4	1,1-Dichloroethylene	ND	0.50	0.054	ug/l	
107-06-2	1,2-Dichloroethane	ND	0.50	0.034	ug/l	
78-87-5	1,2-Dichloropropane	ND	0.50	0.082	ug/l	
124-48-1	Dibromochloromethane	ND	0.50	0.042	ug/l	
156-60-5	trans-1,2-Dichloroethylene	ND	0.50	0.039	ug/l	
156-59-2	cis-1,2-Dichloroethylene	ND	0.50	0.081	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.033	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.063	ug/l	J
100-41-4	Ethylbenzene	ND	0.50	0.033	ug/l	
76-13-1	Freon 113	ND	1.0	0.10	ug/l	
591-78-6	2-Hexanone	ND	2.0	0.084	ug/l	
75-09-2	Methylene chloride	ND	0.50	0.047	ug/l	
108-10-1	4-Methyl-2-pentanone	ND	2.0	0.27	ug/l	
100-42-5	Styrene	ND	0.50	0.028	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	0.50	0.050	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.035	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	0.50	0.052	ug/l	
127-18-4	Tetrachloroethylene	ND	0.50	0.091	ug/l	
108-88-3	Toluene	ND	0.50	0.044	ug/l	

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound

## Report of Analysis

<b>Client Sample ID:</b> BPOW 5-5		<b>Date Sampled:</b> 09/15/15
<b>Lab Sample ID:</b> JC3916-3		<b>Date Received:</b> 09/16/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	101%		78-114%
460-00-4	4-Bromofluorobenzene	86%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

ND = Not detected      MDL = Method Detection Limit  
 RL = Reporting Limit  
 E = Indicates value exceeds calibration range

J = Indicates an estimated value  
 B = Indicates analyte found in associated method blank  
 N = Indicates presumptive evidence of a compound

4.3  
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## Report of Analysis

<b>Client Sample ID:</b> BPOW 5-5	<b>Date Sampled:</b> 09/15/15
<b>Lab Sample ID:</b> JC3916-3	<b>Date Received:</b> 09/16/15
<b>Matrix:</b> AQ - Ground Water	<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8270D BY SIM SW846 3510C	
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY	

Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 <sup>a</sup>	I97909.D	1	09/23/15	AMA	09/19/15	M:OP44675	M:MSI3661
Run #2							

Run #	Initial Volume	Final Volume
Run #1	910 ml	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
123-91-1	1,4-Dioxane	0.39	0.22	0.084	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	68%		26-121%		
321-60-8	2-Fluorobiphenyl	63%		28-107%		
1718-51-0	Terphenyl-d14	78%		29-129%		

(a) Analysis performed at Accutest Laboratories, Marlborough, MA.

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ND = Not detected	MDL = Method Detection Limit	J = Indicates an estimated value
RL = Reporting Limit		B = Indicates analyte found in associated method blank
E = Indicates value exceeds calibration range		N = Indicates presumptive evidence of a compound

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GW  
FB  
WIB

FED-EX Tracking # #4  
Accutest Quote #  
Bottle Order Control #  
Accutest Job # JC4092

Client / Reporting Information		Project Information				Requested Analysis ( see TEST CODE sheet)										Matrix Codes	
Company Name <b>Arcadis</b>		Project Name: <b>AGMNYM62235 // OU2 Outpost Wells Northrop Grumman OU2 Hydro</b>				<b>V5242NG FULL LIST B8270SIM14DIOX</b>										DW - Drinking Water GW - Ground Water WW - Water SW - Surface Water SO - Soil SL - Sludge SED - Sediment OI - Oil LIQ - Other Liquid AIR - Air SOL - Other Solid WP - Wipe FB - Field Blank EB - Equipment Blank RB - Rinse Blank TB - Trip Blank	
Street Address <b>2 Huntington Quad, Suite 1S10</b>		Billing Information ( If different from Report to)															
City State Zip <b>Melville NY 11747</b>		Company Name <b>Bethpage NY</b>															
Project Contact <b>Soma Das, soma.das@arcadis-us.com</b>		Arcadis, U.S., Inc. Attn: Accts Payable															
Phone # Fax # <b>631-391-5247 631-249-7610</b>		Client Purchase Order # <b>NY001496.1514.NAVI3</b>				Street Address <b>630 Plaza Drive, Suite 600</b>										LAB USE ONLY	
Sampler(s) Name(s) <b>PATRICIA PREZORSKI, BRENDAN TONG</b>		Work Authorization #: NY001496 2015				City State Zip <b>Highlands Ranch, CO 80129</b>											
E-mail <b>516 287-6147</b>		Project Manager <b>Carlo San Giovanni</b>				Attention: <b>Soma Das</b>											
Collection		Number of preserved Bottles															
Accutest Sample #	Field ID / Point of Collection	MECH/DI Val #	Date	Time	Sampled by	Matrix	# of bottles	HCl	NACH	HNUS	HSO4	HSO3A	NO3E	DI Water	MEDH	ENCORE	
1	TR091615B7A		9/16/15	1000	JP	FB	2										
2	FB091615B7A		9/16/15	1010	JP	FB	2										
3	BR091615B7A		9/16/15	1259	JP	GL	3										

Turnaround Time ( Business days)		Approved By (Accutest PM) / Date:		Data Deliverable Information				Comments / Special Instructions	
<input type="checkbox"/> Std. 15 Business Days <input checked="" type="checkbox"/> Std. 10 Business Days ( by Contract only) <input type="checkbox"/> 10 Day RUSH <input type="checkbox"/> 5 Day RUSH <input type="checkbox"/> 3 Day EMERGENCY <input type="checkbox"/> 2 Day EMERGENCY <input type="checkbox"/> 1 Day EMERGENCY Emergency & Rush T/A data available VIA Lablink		INITIAL ASSESSMENT: ZGM LABEL VERIFICATION: JK		<input type="checkbox"/> Commercial "A" ( Level 1) <input type="checkbox"/> Commercial "B" ( Level 2) <input type="checkbox"/> FULLT1 ( Level 3+4 ) <input type="checkbox"/> NJ Reduced <input type="checkbox"/> Commercial "C" <input type="checkbox"/> NYASP Category A <input type="checkbox"/> NYASP Category B <input type="checkbox"/> State Forms <input type="checkbox"/> EDD Format <input checked="" type="checkbox"/> Other COMMC+				OU2 Hydro Analyte List (V5242NG -FULL LIST) plus 1,4-Dioxane (B8270SIM14DIOX) Please use " " for MS/MSD QA/QC Sample.	
Sample Custody must be documented below each time samples change possession, including courier delivery.									
Relinquished by Sampler: <b>1 Patricia Prezorski</b>		Date Time: <b>9/16/15 2000</b>		Received By: <b>[Signature]</b>		Date Time: <b>9/16/15 17:00</b>		Relinquished By: <b>2 [Signature]</b>	
Relinquished by Sampler: <b>3 Brendan Tong</b>		Date Time: <b>9/16/15 17:00</b>		Received By: <b>[Signature]</b>		Date Time: <b>9/16/15 17:00</b>		Relinquished By: <b>4 [Signature]</b>	
Relinquished by: <b>5</b>		Date Time: <b>9/16/15 17:00</b>		Received By: <b>[Signature]</b>		Date Time: <b>9/16/15 17:00</b>		Relinquished By: <b>4</b>	

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<b>Client / Reporting Information</b> Company Name: <b>Arcadis</b> Street Address: <b>2 Huntington Quad, Suite 1S10</b> City/State/Zip: <b>Melville NY 11747</b> Project Contact: <b>Soma Das, soma.das@arcadis-us.com</b> Phone/Fax: <b>631-391-5247 / 631-249-7610</b> Sample(s) Name(s): <b>PATRICIA PREZORSKI, BRENDAN TONG</b>		<b>Project Information</b> Project Name: <b>AGMNYM62235 // OU2 Outpost Wells Northrop Grumman OU2 Hydro</b> Billing Information (if different from Report to): Company Name: <b>Arcadis, U.S., Inc. Attn: Accts Payable</b> Street Address: <b>630 Plaza Drive, Suite 600</b> City/State/Zip: <b>Highlands Ranch, CO 80129</b> Project Manager: <b>Carlo San Giovanni</b>		<b>Requested Analysis (see TEST CODE sheet)</b> V5242NG FULL LIST B8270SIM14DIOX		<b>Matrix Codes</b> DW - Drinking Water GW - Ground Water WW - Water SW - Surface Water SO - Soil SL - Sludge SED - Sediment OI - Oil LIQ - Other Liquid AIR - Air SOL - Other Solid WP - Wipe FB - Field Blank EB - Equipment Blank RB - Rinse Blank TB - Trip Blank
<b>Turnaround Time (Business days)</b> <input type="checkbox"/> Std. 15 Business Days <input checked="" type="checkbox"/> Std. 10 Business Days (by Contract only) <input type="checkbox"/> 10 Day RUSH <input type="checkbox"/> 5 Day RUSH <input type="checkbox"/> 3 Day EMERGENCY <input type="checkbox"/> 2 Day EMERGENCY <input type="checkbox"/> 1 Day EMERGENCY Emergency & Rush TIA data available VIA Lablink		<b>Approved By (Account PM): (Date)</b> _____ _____ _____		<b>Data Deliverable Information</b> <input type="checkbox"/> Commercial "A" (Level 1) <input type="checkbox"/> Commercial "B" (Level 2) <input type="checkbox"/> FULLT1 (Level 3+4) <input type="checkbox"/> NJ Reduced <input type="checkbox"/> Commercial "C" <input type="checkbox"/> NYASP Category A <input type="checkbox"/> NYASP Category B <input type="checkbox"/> State Forms <input type="checkbox"/> EDD Format <input checked="" type="checkbox"/> Other <b>COMM/C+</b> Commercial "A" = Results Only Commercial "B" = Results + QC Summary NJ Reduced = Results + QC Summary + Partial Raw data		<b>Comments / Special Instructions</b> OU2 Hydro Analyte List (V5242NG - FULL LIST) plus 1,4-Dioxane (B8270SIM14DIOX) Please use " " for MS/MSD QA/QC Sample.
<b>Sample Custody must be documented below each time samples change possession, including courier delivery.</b>						
Relinquished by: <b>Patricia Prezorski</b> Date/Time: <b>9/16/15 18:00</b>	Received By: <b>1</b> Date/Time: _____	Relinquished by: <b>2</b> Date/Time: _____	Received By: <b>2</b> Date/Time: _____	Relinquished by: <b>3</b> Date/Time: _____	Received By: <b>3</b> Date/Time: _____	
Relinquished by: <b>Brendan Tong</b> Date/Time: _____	Received By: <b>4</b> Date/Time: _____	Relinquished by: <b>4</b> Date/Time: _____	Received By: <b>4</b> Date/Time: _____	Relinquished by: <b>5</b> Date/Time: <b>9-17-15 7:30</b>	Received By: <b>5</b> Date/Time: _____	
Custody Seal # _____ <input type="checkbox"/> Intact <input type="checkbox"/> Not Intact		Preserved where applicable <input type="checkbox"/>		On Ice <input checked="" type="checkbox"/> / Cooler Temp. <b>13°C</b>		

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## Report of Analysis

<b>Client Sample ID:</b> TB091615BT1	<b>Date Sampled:</b> 09/16/15
<b>Lab Sample ID:</b> JC4092-1	<b>Date Received:</b> 09/17/15
<b>Matrix:</b> AQ - Trip Blank Water	<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1	
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY	

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #2	1B99463.D	1	09/19/15	MD	n/a	n/a	V1B4714

Run #1	Purge Volume
Run #2	5.0 ml

## Special VOA List

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

ND = Not detected MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

## Report of Analysis

<b>Client Sample ID:</b> TB091615BT1		<b>Date Sampled:</b> 09/16/15
<b>Lab Sample ID:</b> JC4092-1		<b>Date Received:</b> 09/17/15
<b>Matrix:</b> AQ - Trip Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	102%		78-114%
460-00-4	4-Bromofluorobenzene	85%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

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ND = Not detected      MDL = Method Detection Limit      J = Indicates an estimated value  
 RL = Reporting Limit      B = Indicates analyte found in associated method blank  
 E = Indicates value exceeds calibration range      N = Indicates presumptive evidence of a compound

4.1  
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## Report of Analysis

<b>Client Sample ID:</b> FB091615BT1		
<b>Lab Sample ID:</b> JC4092-2		<b>Date Sampled:</b> 09/16/15
<b>Matrix:</b> AQ - Field Blank Water		<b>Date Received:</b> 09/17/15
<b>Method:</b> EPA 524.2 REV 4.1		<b>Percent Solids:</b> n/a
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B99464.D	1	09/19/15	MD	n/a	n/a	V1B4714
Run #2							

Run #1	Purge Volume
Run #1	5.0 ml
Run #2	

## Special VOA List

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

ND = Not detected MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

## Report of Analysis

<b>Client Sample ID:</b> FB091615BT1		<b>Date Sampled:</b> 09/16/15
<b>Lab Sample ID:</b> JC4092-2		<b>Date Received:</b> 09/17/15
<b>Matrix:</b> AQ - Field Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	103%		78-114%
460-00-4	4-Bromofluorobenzene	83%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

ND = Not detected      MDL = Method Detection Limit  
 RL = Reporting Limit  
 E = Indicates value exceeds calibration range

J = Indicates an estimated value  
 B = Indicates analyte found in associated method blank  
 N = Indicates presumptive evidence of a compound

4.2  
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## Report of Analysis

<b>Client Sample ID:</b> FB091615BT1		<b>Date Sampled:</b> 09/16/15
<b>Lab Sample ID:</b> JC4092-2		<b>Date Received:</b> 09/17/15
<b>Matrix:</b> AQ - Field Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8270D BY SIM SW846 3510C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 <sup>a</sup>	I98026.D	1	09/30/15	AMA	09/20/15	M:OP44688	M:MSI3665
Run #2							

	Initial Volume	Final Volume
Run #1	980 ml	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
123-91-1	1,4-Dioxane	ND	0.20	0.078	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	51%		26-121%		
321-60-8	2-Fluorobiphenyl	47%		28-107%		
1718-51-0	Terphenyl-d14	59%		29-129%		

(a) Analysis performed at Accutest Laboratories, Marlborough, MA.

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ND = Not detected	MDL = Method Detection Limit	J = Indicates an estimated value
RL = Reporting Limit		B = Indicates analyte found in associated method blank
E = Indicates value exceeds calibration range		N = Indicates presumptive evidence of a compound

4.2  
4



## Report of Analysis

<b>Client Sample ID:</b> BPOW 5-6		<b>Date Sampled:</b> 09/16/15
<b>Lab Sample ID:</b> JC4092-3		<b>Date Received:</b> 09/17/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B99465.D	1	09/19/15	MD	n/a	n/a	V1B4714
Run #2							

Run #1	Purge Volume
Run #1	5.0 ml
Run #2	

## Special VOA List

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

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound

## Report of Analysis

<b>Client Sample ID:</b> BPOW 5-6		<b>Date Sampled:</b> 09/16/15
<b>Lab Sample ID:</b> JC4092-3		<b>Date Received:</b> 09/17/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	101%		78-114%
460-00-4	4-Bromofluorobenzene	84%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

ND = Not detected      MDL = Method Detection Limit  
 RL = Reporting Limit  
 E = Indicates value exceeds calibration range

J = Indicates an estimated value  
 B = Indicates analyte found in associated method blank  
 N = Indicates presumptive evidence of a compound

4.3  
4

## Report of Analysis

<b>Client Sample ID:</b> BPOW 5-6		<b>Date Sampled:</b> 09/16/15
<b>Lab Sample ID:</b> JC4092-3		<b>Date Received:</b> 09/17/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8270D BY SIM SW846 3510C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 <sup>a</sup>	I97967.D	1	09/28/15	AMA	09/20/15	M:OP44688	M:MSI3663
Run #2							

	Initial Volume	Final Volume
Run #1	880 ml	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
123-91-1	1,4-Dioxane	ND	0.23	0.086	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	46%		26-121%		
321-60-8	2-Fluorobiphenyl	40%		28-107%		
1718-51-0	Terphenyl-d14	55%		29-129%		

(a) Analysis performed at Accutest Laboratories, Marlborough, MA.

---

ND = Not detected	MDL = Method Detection Limit	J = Indicates an estimated value
RL = Reporting Limit		B = Indicates analyte found in associated method blank
E = Indicates value exceeds calibration range		N = Indicates presumptive evidence of a compound

4.3  
4

## **Northrop Grumman Corporation- Operable Unit 2**

### **Data Review**

BETHPAGE, NEW YORK

Volatile and Semivolatile Analyses

SDGs #JC3119, JC3192 and JC3373

Analyses Performed By:  
Accutest Laboratories  
Dayton, New Jersey

Report #24353R  
September 29, 2015  
Review Level: Tier II  
Project #NY001496.1514.NAVI4

## SUMMARY

This data quality assessment summarizes the review of Sample Delivery Groups (SDGs) #JC3119, JC3192 and JC3373 for samples collected in association with the Northrop Grumman-Bethpage Site. The review was conducted as a Tier II evaluation and included review of data package completeness. Only analytical data associated with constituents of concern were reviewed for this validation. Field documentation was not included in this review. Included with this assessment are the validation annotated sample result sheets, and chain of custody. Analyses were performed on the following samples:

SDGs	Sample ID	Lab ID	Matrix	Sample Collection Date	Parent Sample	Analysis				
						VOC	SVOC	PCB	MET	MISC
JC3119	FB090315BT1	JC3119-1	Water	09/03/2015		X	X			
	TB090315BT1	JC3119-2	Water	09/03/2015		X				
	BPOW 6-1	JC3119-3	Water	09/03/2015		X	X			
JC3192	FB090415BT1	JC3192-1	Water	09/04/2015		X	X			
	TB090415BT1	JC3192-2	Water	09/04/2015		X				
	BPOW 6-2	JC3192-3	Water	09/04/2015		X	X			
JC3373	TB090815BT1	JC3373-1	Water	09/08/2015		X				
	FB090815BT1	JC3373-2	Water	09/08/2015		X	X			
	BPOW 6-3	JC3373-3	Water	09/08/2015		X	X			

Notes:

1. Semivolatile (1,4-Dioxane) analysis was performed at Accutest Laboratories, Marlborough, MA (Accutest-NE) facility.

## ANALYTICAL DATA PACKAGE DOCUMENTATION

### GENERAL INFORMATION

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

QA - Quality Assurance

## VOLATILE ANALYSIS INTRODUCTION

Analyses were performed according to United States Environmental Protection Agency (USEPA), Methods 524.2 and 8270D-Selected Ion Monitoring (SIM). Data were reviewed in accordance with USEPA National Functional Guidelines of October 1999.

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

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

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

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

## VOLATILE ORGANIC COMPOUND (VOC) ANALYSES

### 1. Holding Times

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

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

s.u. Standard units

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

### 2. Blank Contamination

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

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

Compounds were detected in the associated QA blanks; however, the associated sample results were greater than the BAL and/or were non-detect. No qualification of the sample results was required.

### 3. Surrogates/System Monitoring Compounds

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

All surrogate recoveries were within control limits.

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

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

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

A MS/MSD analysis was not performed on a sample location associated with these SDGs.



## **5. Laboratory Control Sample (LCS) Analysis**

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

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

## **6. Field Duplicate Analysis**

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

A field duplicate was not collected on a sample location associated with these SDGs.

## **7. Laboratory Duplicate Analysis**

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

A laboratory duplicate sample was not performed on a sample location associated with these SDGs.

## **8. System Performance and Overall Assessment**

Tentatively identified compounds (TICs) were not identified in any of the sample locations.

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

## DATA VALIDATION CHECKLIST FOR VOCs

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

%R Percent Recovery    RPD Relative Percent Difference

# SEMIVOLATILE ORGANIC COMPOUND (SVOC) ANALYSES

## 1. Holding Times

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

Method	Matrix	Holding Time	Preservation
SW-846 8270D SIM	Water	7 days from collection to extraction and 40 days from extraction to analysis	Cooled to <6°C

All samples were analyzed within the specified holding time criteria.

## 2. Blank Contamination

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

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

Compounds were not detected above the MDL in the associated blanks; therefore detected sample results were not associated with blank contamination.

## 3. Surrogates/System Monitoring Compounds

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

All surrogate recoveries were within control limits.

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

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

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

A MS/MSD analysis was not performed on a sample location associated with these SDGs.

## **5. Laboratory Control Sample/Laboratory Control Sample Duplicate (LCS/LCSD) Analysis**

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

All compounds associated with the LCS/LCSD analysis exhibited recoveries and RPD within the control limits.

## **6. Field Duplicate Analysis**

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

A field duplicate was not collected with a sample location associated with these SDGs.

## **7. System Performance and Overall Assessment**

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

## DATA VALIDATION CHECKLIST FOR SVOCs

SVOCs: SW-846 8270D-SIM	Reported		Performance Acceptable		Not Required
	No	Yes	No	Yes	
GAS CHROMATOGRAPHY/MASS SPECTROMETRY (GC/MS)					
<b>Tier II Validation</b>					
Holding times		X		X	
Reporting limits (units)		X		X	
Blanks					
A. Method blanks		X		X	
B. Equipment blanks		X		X	
Laboratory Control Sample (LCS) %R		X		X	
Laboratory Control Sample Duplicate(LCSD) %R		X		X	
LCS/LCSD Precision (RPD)		X		X	
Matrix Spike (MS) %R					X
Matrix Spike Duplicate(MSD) %R					X
MS/MSD Precision (RPD)					X
Field/Lab Duplicate (RPD)					X
Surrogate Spike Recoveries		X		X	
Dilution Factor		X		X	
Moisture Content					X

%R    Percent recovery                      RPD    Relative percent difference

VALIDATION PERFORMED BY: Lisa Horton

SIGNATURE:   
\_\_\_\_\_

DATE: September 29, 2015

PEER REVIEW BY: Todd Church

DATE: October 5, 2015

**CHAIN OF CUSTODY/  
CORRECTED SAMPLE ANALYSIS DATA SHEETS**

GU  
FB  
WIB

**CHAIN OF CUSTODY**  
Accutest New Jersey/SPL Environmental  
2235 Route 130, Dayton, NJ 08810  
TEL: 732-329-0200 FAX: 732-329-3499/3480  
www.accutest.com

Client / Reporting Information		Project Information		Requested Analysis (see TEST CODE sheet)		Matrix Codes	
Company Name <b>Arcadis</b>		Project Name <b>AGMNYM62235 // OU2 Outpost Wells Northrop Grumman OU2 Hydro</b>		FED-EX Tracking # <b>#4</b>		Bottle Order Control #	
Street Address <b>2 Huntington Quad, Suite 1S10</b>		Street		Accutest Quote #		Accutest Job # <b>JC3119</b>	
City State Zip <b>Melville NY 11747</b>		City State <b>Bethpage NY</b>		Billing Information (if different from Report to)		Matrix Codes DW - Drinking Water GW - Ground Water WW - Water SW - Surface Water SO - Soil SL - Sludge SED - Sediment OI - Oil LIQ - Other Liquid AR - Air SOL - Other Solid WP - Wipe FB - Field Blank EB - Equipment Blank RB - Rinse Blank TB - Trip Blank	
Project Contact <b>Soma Das, soma.das@arcadis-us.com</b>		Project # <b>NY001496.314I.NAVI3</b>		Company Name <b>Arcadis, U.S., Inc. Attn: Accts Payable</b>		Requested Analysis (see TEST CODE sheet)	
Phone # <b>631-391-5247</b>		Client Purchase Order # <b>NY001496_2015</b>		Street Address <b>630 Plaza Drive, Suite 600</b>			
FAX # <b>631-249-7610</b>		Work Authorization #		City State Zip <b>Highlands Ranch, CO 80129</b>		LAB USE ONLY	
Sampler(s) Name(s) <b>PATRICIA PREZORSKI, BRENDAN TONG</b>		Project Manager <b>Carlo San Giovanni</b>		Attention: <b>Soma Das</b>			
Field ID / Point of Collection		MECHDI Val #		Date		Time	
1 <b>FB090315B7A</b>				9/3/15		0850	
2 <b>TB090315B7A</b>				9/3/15		0830	
3 <b>BP090315G-1</b>				9/3/15		1208	
Turnaround Time (Business days)		Approved By (Accutest PM) / Date:		Data Deliverable Information		Comments / Special Instructions	
<input type="checkbox"/> Std. 15 Business Days <input checked="" type="checkbox"/> Std. 10 Business Days (by Contract only) <input type="checkbox"/> 10 Day RUSH <input type="checkbox"/> 5 Day RUSH <input type="checkbox"/> 3 Day EMERGENCY <input type="checkbox"/> 2 Day EMERGENCY <input type="checkbox"/> 1 Day EMERGENCY Emergency & Rush T/A data available VIA Lablink		Approved By: <b>Chris Law</b> INITIAL ASSESSMENT <b>2B OB</b> LABEL VERIFICATION <b>OB</b>		<input type="checkbox"/> Commercial "A" (Level 1) <input type="checkbox"/> Commercial "B" (Level 2) <input type="checkbox"/> FULLT1 (Level 3+4) <input type="checkbox"/> NJ Reduced <input type="checkbox"/> Commercial "C" Commercial "A" = Results Only Commercial "B" = Results + QC Summary NJ Reduced = Results + QC Summary + Partial Raw data		OU2 Hydro Analyte List (V5242NG14QW140) plus 1,4-Dioxane (B8270SIM14DIOX) <b>Full list</b> Please use " " for MS/MSD QA/QC Sample.	
Sample Custody must be documented below each time samples change possession, including courier delivery.							
Relinquished by Sampler:	Date/Time:	Received By:	Date/Time:	Relinquished by:	Date/Time:	Received By:	Date/Time:
1 Patricia Prezorski	9/3/15 1745	Chris Law	9/3/15 12:52	Chris Law	9/3/15/1653		
3 Brendan Tong							
5							
Custody Seal #		<input checked="" type="checkbox"/> Intact		Preserved where applicable		On Ice Cooler Temp. <b>2.1°C</b>	

5.1  
5



CHAIN OF CUSTODY  
 Accutest New Jersey/SPL Environmental  
 2335 Route 130, Dayton, NJ  
 TEL: 732-329-0200 FAX: 732-329-3499/3480

Jc3119 PAGE 1 OF 1



CHAIN OF CUSTODY  
 Accutest New Jersey/SPL Environmental  
 2235 Route 130, Dayton, NJ 08810  
 TEL: 732-329-0200 FAX: 732-329-3499/3480  
 www.acctest.com

FED-EX Tracking # #4  
 Accutest Quote #  
 Bottle Order Control #  
 Accutest Job #

**Client / Reporting Information**  
 Company Name: **Arcadis**  
 Street Address: **2 Huntington Quad, Suite 1S10**  
 City: **Melville** State: **NY** Zip: **11747**  
 Project Contact: **Soma Das**  
 Phone #: **631-391-5247** E-mail: **soma.das@arcadis-us.com**

**Project Information**  
 Project Name: **AGMNYM62235 // OU2 Outpost Wells Northrop Grumman OU2 Hydro**  
 Project #:  
 Billing Information (if different from Report to):  
 Company Name: **Bethpage** State: **NY**  
 Project #:  
 Client Purchase Order #: **NY001496.314I.NAVI3**  
 Work Authorization #: **NY001496\_2015**

**Requested Analysis (see TEST CODE sheet)**  
 V5242NG44GW40 Full List  
 B8270SIM14DIOX

**Matrix Codes**  
 DW - Drinking Water  
 GW - Ground Water  
 WW - Water  
 SW - Surface Water  
 SO - Soil  
 SL - Sludge  
 SED - Sediment  
 OL - Oil  
 LIQ - Other Liquid  
 AIR - Air  
 SOL - Other Solid  
 WP - Wipe  
 FB - Field Blank  
 EB - Equipment Blank  
 RB - Rinse Blank  
 TB - Trip Blank

**LAB USE ONLY**

Accession Sample #	Field ID / Point of Collection	MECHID/Val #	Collection		Sampled by	Matrix	# of bottles	Number of preserved bottles												
			Date	Time				INCH	INSD	INSPD	INQIC	INWIP	INWTR	INWED	INWRE					
	FB090315BT1		9/3/15	0850	SP	FB	2													
	BROW G-1		9/3/15	1208	SP	GL	2													

**Turnaround Time (Business days)**  
 Std. 15 Business Days  
 Std. 10 Business Days (by Contract only)  
 10 Day RUSH  
 5 Day RUSH  
 3 Day EMERGENCY  
 2 Day EMERGENCY  
 1 Day EMERGENCY  
 Policy & Rush TIA data available VIA LabLink

**Approved By (Accutest PM): / Date:**

**Data Deliverable Information**  
 Commercial "A" (Level 1)  NYASP Category A  
 Commercial "B" (Level 2)  NYASP Category B  
 FULLT1 (Level 3+4)  State Forms  
 NJ Reduced  EDD Format  
 Commercial "C"  Other **COMMC+**  
 Commercial "A" = Results Only  
 Commercial "B" = Results + QC Summary  
 NJ Reduced = Results + QC Summary + Partial Raw data

**Comments / Special Instructions**  
 OU2 Hydro Analyte List (V5242NG44GW40) plus 1,4-Dioxane (B8270SIM14DIOX) Full List  
 Please use \* for MS/MSD QA/QC Sample.

Sample Custody must be documented below each time samples change possession, including courier delivery.

Received By:	Date Time:	Received By:	Date Time:	Received By:	Date Time:
1	4/3/15 1600	2		3	
3		4		5	

Relinquished By: 2, 4

Custody Seal #  Intact  Preserved where applicable  On Ice  Cool Temp.

5.1  
5

## Report of Analysis

<b>Client Sample ID:</b> FB090315BT1		
<b>Lab Sample ID:</b> JC3119-1		<b>Date Sampled:</b> 09/03/15
<b>Matrix:</b> AQ - Field Blank Water		<b>Date Received:</b> 09/03/15
<b>Method:</b> EPA 524.2 REV 4.1		<b>Percent Solids:</b> n/a
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B99199.D	1	09/08/15	MD	n/a	n/a	V1B4703
Run #2							

Run #1	Purge Volume
Run #1	5.0 ml
Run #2	

## Special VOA List

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

ND = Not detected

MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

## Report of Analysis

<b>Client Sample ID:</b> FB090315BT1	<b>Date Sampled:</b> 09/03/15
<b>Lab Sample ID:</b> JC3119-1	<b>Date Received:</b> 09/03/15
<b>Matrix:</b> AQ - Field Blank Water	<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1	
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY	

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	100%		78-114%
460-00-4	4-Bromofluorobenzene	92%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

ND = Not detected      MDL = Method Detection Limit  
 RL = Reporting Limit  
 E = Indicates value exceeds calibration range

J = Indicates an estimated value  
 B = Indicates analyte found in associated method blank  
 N = Indicates presumptive evidence of a compound

4.1  
4

## Report of Analysis

<b>Client Sample ID:</b> FB090315BT1	<b>Date Sampled:</b> 09/03/15
<b>Lab Sample ID:</b> JC3119-1	<b>Date Received:</b> 09/03/15
<b>Matrix:</b> AQ - Field Blank Water	<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8270D BY SIM SW846 3510C	
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY	

Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 <sup>a</sup>	I97761.D	1	09/16/15	AMA	09/08/15	M:OP44541	M:MSI3656
Run #2							

Run #	Initial Volume	Final Volume
Run #1	900 ml	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
123-91-1	1,4-Dioxane	ND	0.22	0.084	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	79%		26-121%		
321-60-8	2-Fluorobiphenyl	75%		28-107%		
1718-51-0	Terphenyl-d14	93%		29-129%		

(a) Analysis performed at Accutest Laboratories, Marlborough, MA.

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ND = Not detected	MDL = Method Detection Limit	J = Indicates an estimated value
RL = Reporting Limit		B = Indicates analyte found in associated method blank
E = Indicates value exceeds calibration range		N = Indicates presumptive evidence of a compound

4.1  
4

## Report of Analysis

<b>Client Sample ID:</b> TB090315BT1		<b>Date Sampled:</b> 09/03/15
<b>Lab Sample ID:</b> JC3119-2		<b>Date Received:</b> 09/03/15
<b>Matrix:</b> AQ - Trip Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B99200.D	1	09/09/15	MD	n/a	n/a	V1B4703
Run #2							

Run #1	Purge Volume
Run #1	5.0 ml
Run #2	

## Special VOA List

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

ND = Not detected MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

## Report of Analysis

<b>Client Sample ID:</b> TB090315BT1		<b>Date Sampled:</b> 09/03/15
<b>Lab Sample ID:</b> JC3119-2		<b>Date Received:</b> 09/03/15
<b>Matrix:</b> AQ - Trip Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	98%		78-114%
460-00-4	4-Bromofluorobenzene	90%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

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ND = Not detected      MDL = Method Detection Limit      J = Indicates an estimated value  
 RL = Reporting Limit      B = Indicates analyte found in associated method blank  
 E = Indicates value exceeds calibration range      N = Indicates presumptive evidence of a compound

4.2  
4

## Report of Analysis

<b>Client Sample ID:</b> BPOW6-1		<b>Date Sampled:</b> 09/03/15
<b>Lab Sample ID:</b> JC3119-3		<b>Date Received:</b> 09/03/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B99201.D	1	09/09/15	MD	n/a	n/a	V1B4703
Run #2							

Run #1	Purge Volume
Run #1	5.0 ml
Run #2	

## Special VOA List

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

ND = Not detected MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

## Report of Analysis

<b>Client Sample ID:</b> BPOW6-1		<b>Date Sampled:</b> 09/03/15
<b>Lab Sample ID:</b> JC3119-3		<b>Date Received:</b> 09/03/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	100%		78-114%
460-00-4	4-Bromofluorobenzene	89%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

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ND = Not detected      MDL = Method Detection Limit      J = Indicates an estimated value  
 RL = Reporting Limit      B = Indicates analyte found in associated method blank  
 E = Indicates value exceeds calibration range      N = Indicates presumptive evidence of a compound

4.3  
4



## Report of Analysis

<b>Client Sample ID:</b> BPOW6-1		<b>Date Sampled:</b> 09/03/15
<b>Lab Sample ID:</b> JC3119-3		<b>Date Received:</b> 09/03/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8270D BY SIM SW846 3510C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 <sup>a</sup>	I97762.D	1	09/16/15	AMA	09/08/15	M:OP44541	M:MSI3656
Run #2							

	Initial Volume	Final Volume
Run #1	950 ml	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
123-91-1	1,4-Dioxane	ND	0.21	0.080	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	78%		26-121%		
321-60-8	2-Fluorobiphenyl	71%		28-107%		
1718-51-0	Terphenyl-d14	83%		29-129%		

(a) Analysis performed at Accutest Laboratories, Marlborough, MA.

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ND = Not detected	MDL = Method Detection Limit	J = Indicates an estimated value
RL = Reporting Limit		B = Indicates analyte found in associated method blank
E = Indicates value exceeds calibration range		N = Indicates presumptive evidence of a compound

4.3  
4





## Report of Analysis

<b>Client Sample ID:</b> FB090415BT1		<b>Date Sampled:</b> 09/04/15
<b>Lab Sample ID:</b> JC3192-1		<b>Date Received:</b> 09/04/15
<b>Matrix:</b> AQ - Field Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B99220.D	1	09/09/15	MD	n/a	n/a	V1B4705
Run #2							

Run #1	Purge Volume
Run #1	5.0 ml
Run #2	

## Special VOA List

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	1.2	5.0	0.91	ug/l	J
78-93-3	2-Butanone	ND	5.0	0.57	ug/l	
71-43-2	Benzene	ND	0.50	0.057	ug/l	
75-27-4	Bromodichloromethane	ND	0.50	0.082	ug/l	
75-25-2	Bromoform	ND	0.50	0.046	ug/l	
74-83-9	Bromomethane	ND	0.50	0.077	ug/l	
75-15-0	Carbon disulfide	ND	0.50	0.028	ug/l	
108-90-7	Chlorobenzene	ND	0.50	0.027	ug/l	
75-00-3	Chloroethane	ND	0.50	0.037	ug/l	
67-66-3	Chloroform	ND	0.50	0.031	ug/l	
74-87-3	Chloromethane	ND	0.50	0.044	ug/l	
56-23-5	Carbon tetrachloride	ND	0.50	0.074	ug/l	
75-34-3	1,1-Dichloroethane	ND	0.50	0.039	ug/l	
75-35-4	1,1-Dichloroethylene	ND	0.50	0.054	ug/l	
107-06-2	1,2-Dichloroethane	ND	0.50	0.034	ug/l	
78-87-5	1,2-Dichloropropane	ND	0.50	0.082	ug/l	
124-48-1	Dibromochloromethane	ND	0.50	0.042	ug/l	
156-60-5	trans-1,2-Dichloroethylene	ND	0.50	0.039	ug/l	
156-59-2	cis-1,2-Dichloroethylene	ND	0.50	0.081	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.033	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.063	ug/l	
100-41-4	Ethylbenzene	ND	0.50	0.033	ug/l	
76-13-1	Freon 113	ND	1.0	0.10	ug/l	
591-78-6	2-Hexanone	ND	2.0	0.084	ug/l	
75-09-2	Methylene chloride	ND	0.50	0.047	ug/l	
108-10-1	4-Methyl-2-pentanone	ND	2.0	0.27	ug/l	
100-42-5	Styrene	ND	0.50	0.028	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	0.50	0.050	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.035	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	0.50	0.052	ug/l	
127-18-4	Tetrachloroethylene	ND	0.50	0.091	ug/l	
108-88-3	Toluene	ND	0.50	0.044	ug/l	

ND = Not detected MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

## Report of Analysis

<b>Client Sample ID:</b> FB090415BT1	<b>Date Sampled:</b> 09/04/15
<b>Lab Sample ID:</b> JC3192-1	<b>Date Received:</b> 09/04/15
<b>Matrix:</b> AQ - Field Blank Water	<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1	
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY	

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	99%		78-114%
460-00-4	4-Bromofluorobenzene	86%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

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ND = Not detected      MDL = Method Detection Limit      J = Indicates an estimated value  
 RL = Reporting Limit      B = Indicates analyte found in associated method blank  
 E = Indicates value exceeds calibration range      N = Indicates presumptive evidence of a compound

4.1  
4

## Report of Analysis

<b>Client Sample ID:</b> FB090415BT1	<b>Date Sampled:</b> 09/04/15
<b>Lab Sample ID:</b> JC3192-1	<b>Date Received:</b> 09/04/15
<b>Matrix:</b> AQ - Field Blank Water	<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8270D BY SIM SW846 3510C	
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY	

Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 <sup>a</sup>	I97764.D	1	09/16/15	AMA	09/08/15	M:OP44541	M:MSI3656
Run #2							

Run #	Initial Volume	Final Volume
Run #1	900 ml	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
123-91-1	1,4-Dioxane	ND	0.22	0.084	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	85%		26-121%		
321-60-8	2-Fluorobiphenyl	73%		28-107%		
1718-51-0	Terphenyl-d14	84%		29-129%		

(a) Analysis performed at Accutest Laboratories, Marlborough, MA.

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ND = Not detected	MDL = Method Detection Limit	J = Indicates an estimated value
RL = Reporting Limit		B = Indicates analyte found in associated method blank
E = Indicates value exceeds calibration range		N = Indicates presumptive evidence of a compound

4.1  
4

## Report of Analysis

<b>Client Sample ID:</b> TB090415BT1		<b>Date Sampled:</b> 09/04/15
<b>Lab Sample ID:</b> JC3192-2		<b>Date Received:</b> 09/04/15
<b>Matrix:</b> AQ - Trip Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B99221.D	1	09/09/15	MD	n/a	n/a	V1B4705
Run #2							

Run #1	Purge Volume
Run #1	5.0 ml
Run #2	

## Special VOA List

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

ND = Not detected MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

## Report of Analysis

<b>Client Sample ID:</b> TB090415BT1		<b>Date Sampled:</b> 09/04/15
<b>Lab Sample ID:</b> JC3192-2		<b>Date Received:</b> 09/04/15
<b>Matrix:</b> AQ - Trip Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	98%		78-114%
460-00-4	4-Bromofluorobenzene	88%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

ND = Not detected      MDL = Method Detection Limit  
 RL = Reporting Limit  
 E = Indicates value exceeds calibration range

J = Indicates an estimated value  
 B = Indicates analyte found in associated method blank  
 N = Indicates presumptive evidence of a compound

4.2  
4



# Report of Analysis

<b>Client Sample ID:</b> BPOW 6-2		<b>Date Sampled:</b> 09/04/15
<b>Lab Sample ID:</b> JC3192-3		<b>Date Received:</b> 09/04/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B99222.D	1	09/09/15	MD	n/a	n/a	V1B4705
Run #2							

Run #1	Purge Volume
Run #1	5.0 ml
Run #2	

### Special VOA List

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

ND = Not detected      MDL = Method Detection Limit      J = Indicates an estimated value  
 RL = Reporting Limit      B = Indicates analyte found in associated method blank  
 E = Indicates value exceeds calibration range      N = Indicates presumptive evidence of a compound

4.3  
 4

## Report of Analysis

<b>Client Sample ID:</b> BPOW 6-2		<b>Date Sampled:</b> 09/04/15
<b>Lab Sample ID:</b> JC3192-3		<b>Date Received:</b> 09/04/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	99%		78-114%
460-00-4	4-Bromofluorobenzene	89%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

ND = Not detected      MDL = Method Detection Limit  
 RL = Reporting Limit  
 E = Indicates value exceeds calibration range

J = Indicates an estimated value  
 B = Indicates analyte found in associated method blank  
 N = Indicates presumptive evidence of a compound

4.3  
4

## Report of Analysis

<b>Client Sample ID:</b> BPOW 6-2	<b>Date Sampled:</b> 09/04/15
<b>Lab Sample ID:</b> JC3192-3	<b>Date Received:</b> 09/04/15
<b>Matrix:</b> AQ - Ground Water	<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8270D BY SIM SW846 3510C	
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY	

Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 <sup>a</sup>	I97763.D	1	09/16/15	AMA	09/08/15	M:OP44541	M:MSI3656
Run #2							

Run #	Initial Volume	Final Volume
Run #1	700 ml	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
123-91-1	1,4-Dioxane	ND	0.29	0.11	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	79%		26-121%		
321-60-8	2-Fluorobiphenyl	71%		28-107%		
1718-51-0	Terphenyl-d14	85%		29-129%		

(a) Analysis performed at Accutest Laboratories, Marlborough, MA.

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ND = Not detected	MDL = Method Detection Limit	J = Indicates an estimated value
RL = Reporting Limit		B = Indicates analyte found in associated method blank
E = Indicates value exceeds calibration range		N = Indicates presumptive evidence of a compound

4.3  
4

GW  
FB  
W/B

**CHAIN OF CUSTODY**  
Accutest New Jersey/SPL Environmental  
2235 Route 130, Dayton, NJ 08810  
TEL: 732-329-0200 FAX: 732-329-3499/3480  
www.accutest.com

FED-EX Tracking # **#4** Bottle Order Control # **JC3373**  
Accutest Quote # **JC3411** Accutest Job # **09/15**

Client / Reporting Information		Project Information		Requested Analysis (see TEST CODE sheet)										Matrix Codes		
<b>Company Name</b> Arcadis <b>Street Address</b> 2 Huntington Quad, Suite 1S10 <b>City</b> Melville <b>State</b> NY <b>Zip</b> 11747 <b>Project Contact</b> Soma Das, soma.das@arcadis-us.com <b>Phone #</b> 631-391-5247 <b>Fax #</b> 631-249-7610 <b>Sampler(s) Name(s)</b> PATRICIA PREZORSKI, BRENDAN TONG		<b>Project Name</b> AGMNYM62235 // OU2 Outpost Wells <b>Street</b> Northrop Grumman OU2 Hydro <b>City</b> Bethpage <b>State</b> NY <b>Project #</b> 1514 <b>Client Purchase Order #</b> NY001496-514-NAV13 <b>Work Authorization #</b> NY001496_2015 <b>Project Manager</b> Carlo San Giovanni		<b>Billing Information (if different from Report to)</b> <b>Company Name</b> Arcadis, U.S., Inc. Attn: Accts Payable <b>Street Address</b> 630 Plaza Drive, Suite 600 <b>City</b> Highlands Ranch, <b>State</b> CO <b>Zip</b> 80129 <b>Attention:</b> Soma Das										DW - Drinking Water GW - Ground Water WW - Water SW - Surface Water SO - Soil SL - Sludge SED - Sediment OI - Oil LIQ - Other Liquid AIR - Air SOL - Other Solid WP - Wipe FB - Field Blank EB - Equipment Blank RB - Rinse Blank TB - Trip Blank		
Account Sample #	Field ID / Point of Collection	MECH/DI Val #	Date	Time	Sampled by	Matrix	# of bottles	HCl	INCH	INCO	PRECO	NONE	DI Water	MEPH	ENDURE	LAB USE ONLY
1	TB090815BT1		9/8/15	0830	TB	2	2									
2	FB090815BT2		9/8/15	0855	FB	2	2									V1208
3	BBW 6-3		9/8/15	1205	GW	3	3									

V5242NG+6W+40 plus 14-Dioxane (B270SIM14DIOX)

INITIAL ASSESSMENT *[Signature]*  
LABEL VERIFICATION *[Signature]*

Turnaround Time (Business days)		Data Deliverable Information		Comments / Special Instructions	
<input type="checkbox"/> Std. 15 Business Days <input checked="" type="checkbox"/> Std. 10 Business Days (by Contract only) <input type="checkbox"/> 10 Day RUSH <input type="checkbox"/> 5 Day RUSH <input type="checkbox"/> 3 Day EMERGENCY <input type="checkbox"/> 2 Day EMERGENCY <input type="checkbox"/> 1 Day EMERGENCY Emergency & Rush T/A data available VIA Lablink	<b>Approved By (Accutest PM): / Date:</b> _____	<input type="checkbox"/> Commercial "A" (Level 1) <input type="checkbox"/> Commercial "B" (Level 2) <input type="checkbox"/> FULLT1 (Level 3+4) <input type="checkbox"/> NJ Reduced <input type="checkbox"/> Commercial "C" Commercial "A" = Results Only Commercial "B" = Results + QC Summary NJ Reduced = Results + QC Summary + Partial Raw data	<input type="checkbox"/> NYASP Category A <input type="checkbox"/> NYASP Category B <input type="checkbox"/> State Forms <input type="checkbox"/> EDD Format <input checked="" type="checkbox"/> Other COMM-C+	<b>OU2 Hydro Analyte List (V5242NG+6W+40) plus 14-Dioxane (B270SIM14DIOX)</b> Please use " " for MS/MSD QA/QC Sample.	
<b>Sample Custody must be documented below each time samples change possession, including courier delivery.</b>					
<b>Relinquished by Sampler:</b> 1 Patricia Prezorski <b>Date Time:</b> 9/8/15 2030	<b>Received By:</b> Chris Jaul <b>Date Time:</b> 9/8/15 10:00	<b>Relinquished by Sampler:</b> 3 Brendan Tong <b>Date Time:</b>	<b>Received By:</b> Chris Jaul <b>Date Time:</b> 9/8/15 1430	<b>Relinquished by Sampler:</b> 5	<b>Received By:</b> <b>Date Time:</b>
<b>Emergency &amp; Rush T/A data available VIA Lablink</b>		<b>Preserved where applicable</b> <input checked="" type="checkbox"/> Intact <input type="checkbox"/> Not Intact		<b>On Ice</b> <input checked="" type="checkbox"/> <b>Cooler Temp.</b> 1.9°C	

5.1  
5

Client / Reporting Information			Project Information			Requested Analysis ( see TEST CODE sheet)										Matrix Codes
<b>Company Name</b> Arcadis			<b>Project Name:</b> AGMNYM62235 // OU2 Outpost Wells Northrop Grumman OU2 Hydro			FED-EX Tracking # <i>1145</i> Bottle Order Control # <i>JC3373</i>										DW - Drinking Water GW - Ground Water WW - Water SW - Surface Water SO - Soil SL - Sludge SED - Sediment OI - Oil LIQ - Other Liquid AIR - Air SOL - Other Solid WP - Wipe FB - Field Blank EB - Equipment Blank RB - Rinse Blank TB - Trip Blank
<b>Street Address</b> 2 Huntington Quad, Suite 1S10 City: <i>Molville</i> NY State: <i>NY</i> Zip: <i>11747</i>			<b>Street</b> Bethpage NY State: <i>NY</i> Zip: <i>11714</i>			<b>Billing Information ( If different from Report to )</b> Company Name: <i>Arcadis, U.S., Inc. Attn: Accts Payable</i> Street Address: <i>630 Plaza Drive, Suite 600</i> City: <i>Highlands Ranch, CO</i> State: <i>CO</i> Zip: <i>80129</i>										
<b>Project Contact</b> Soma Das, soma.das@arcadis-us.com Phone # <i>631-391-5247</i> Fax # <i>631-249-7610</i>			<b>Client Purchase Order #</b> <i>1514</i> <b>Work Authorization #:</b> NY001496_2015			<b>Requested Analysis</b> <i>V5242NG 14QWA-40</i> <i>B8270SIM14DIOX</i>										
<b>Sample(s) Name(s)</b> PATRICIA PRZORSKI, BRENDAN TONG			<b>Project Manager</b> Carlo San Giovanni			<b>Attention:</b> <i>Soma Das</i>										<b>LAB USE ONLY</b>
Sample #	Field ID / Point of Collection	MEQHD1 Val #	Date	Time	Sampled by	Matrix	# of bottles	LCI	MCN	MGD	USDB	HOHE	DI Water	MEQH	ENCLOSURE	
-2	FB090915BT1		9/8/15	0955	PP	FB	2									
-3	BPOW G-3		9/8/15	1205	PP	GW	2									
<b>Turnaround Time ( Business days )</b> <input type="checkbox"/> Std. 15 Business Days <input checked="" type="checkbox"/> Std. 10 Business Days ( by Contract only ) <input type="checkbox"/> 10 Day RUSH <input type="checkbox"/> 5 Day RUSH <input type="checkbox"/> 3 Day EMERGENCY <input type="checkbox"/> 2 Day EMERGENCY <input type="checkbox"/> 1 Day EMERGENCY			<b>Approved By ( Accutest Pkg. / Date )</b> _____			<b>Data Deliverable Information</b> <input type="checkbox"/> Commercial "A" ( Level 1 ) <input type="checkbox"/> Commercial "B" ( Level 2 ) <input type="checkbox"/> FULLT1 ( Level 3+4 ) <input type="checkbox"/> NJ Reduced <input type="checkbox"/> Commercial "C" <input type="checkbox"/> MYASP Category A <input type="checkbox"/> MYASP Category B <input type="checkbox"/> State Forms <input type="checkbox"/> EDD Format <input checked="" type="checkbox"/> Other <b>COMMCT</b>					<b>Comments / Special Instructions</b> OU2 Hydro Analysis ( V5242NG ) 14-Dioxane ( B8270SIM14DIOX ) Please use <i>PP</i> for MS/MSD QA/QC Sample.					
<b>Emergency &amp; Rush TIA data available VIA Lablink</b>																
<b>Sample Custody must be documented below each time samples change possession, including courier delivery.</b>																
<b>Relinquished by Sample:</b> 1 Patricia Przorski	<b>Date Time:</b> <i>9/8/15 1600</i>		<b>Received By:</b> <i>FX</i> <b>Date Time:</b>				<b>Relinquished by:</b> <i>FX</i>				<b>Date Time:</b> <i>9/8/15 1000</i>		<b>Received By:</b> <i>[Signature]</i>			
<b>Relinquished by Sampler:</b> 3 Brendan Tong	<b>Date Time:</b>		<b>Received By:</b>				<b>Relinquished by:</b>				<b>Date Time:</b>		<b>Received By:</b>			
<b>Relinquished by:</b> 5	<b>Date Time:</b>		<b>Received By:</b>				<b>Custody Seal #</b> <input type="checkbox"/> Intact <input type="checkbox"/> Not Intact				<b>Preserved where applicable</b> <input type="checkbox"/>		<b>Cooler Temp.</b> <i>1.5°C</i>			

51  
5

16F

## Report of Analysis

<b>Client Sample ID:</b> TB090815BT1		<b>Date Sampled:</b> 09/08/15
<b>Lab Sample ID:</b> JC3373-1		<b>Date Received:</b> 09/09/15
<b>Matrix:</b> AQ - Trip Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B99315.D	1	09/14/15	MD	n/a	n/a	V1B4710
Run #2							

Run #1	Purge Volume
Run #1	5.0 ml
Run #2	

## Special VOA List

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

ND = Not detected MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

## Report of Analysis

<b>Client Sample ID:</b> TB090815BT1		<b>Date Sampled:</b> 09/08/15
<b>Lab Sample ID:</b> JC3373-1		<b>Date Received:</b> 09/09/15
<b>Matrix:</b> AQ - Trip Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	101%		78-114%
460-00-4	4-Bromofluorobenzene	87%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

ND = Not detected      MDL = Method Detection Limit  
 RL = Reporting Limit  
 E = Indicates value exceeds calibration range

J = Indicates an estimated value  
 B = Indicates analyte found in associated method blank  
 N = Indicates presumptive evidence of a compound

4.1  
4

## Report of Analysis

<b>Client Sample ID:</b> FB090815BT1	<b>Date Sampled:</b> 09/08/15
<b>Lab Sample ID:</b> JC3373-2	<b>Date Received:</b> 09/09/15
<b>Matrix:</b> AQ - Field Blank Water	<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1	
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY	

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B99316.D	1	09/14/15	MD	n/a	n/a	V1B4710
Run #2							

Run #1	Purge Volume
Run #1	5.0 ml
Run #2	

## Special VOA List

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

ND = Not detected

MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound



## Report of Analysis

<b>Client Sample ID:</b> FB090815BT1		<b>Date Sampled:</b> 09/08/15
<b>Lab Sample ID:</b> JC3373-2		<b>Date Received:</b> 09/09/15
<b>Matrix:</b> AQ - Field Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	102%		78-114%
460-00-4	4-Bromofluorobenzene	88%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

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ND = Not detected      MDL = Method Detection Limit      J = Indicates an estimated value  
 RL = Reporting Limit      B = Indicates analyte found in associated method blank  
 E = Indicates value exceeds calibration range      N = Indicates presumptive evidence of a compound

4.2  
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## Report of Analysis

<b>Client Sample ID:</b> FB090815BT1	<b>Date Sampled:</b> 09/08/15
<b>Lab Sample ID:</b> JC3373-2	<b>Date Received:</b> 09/09/15
<b>Matrix:</b> AQ - Field Blank Water	<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8270D BY SIM SW846 3510C	
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY	

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 <sup>a</sup>	I97870.D	1	09/22/15	AMA	09/11/15	M:OP44578	M:MSI3660
Run #2							

	Initial Volume	Final Volume
Run #1	880 ml	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
123-91-1	1,4-Dioxane	ND	0.23	0.086	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	62%		26-121%		
321-60-8	2-Fluorobiphenyl	53%		28-107%		
1718-51-0	Terphenyl-d14	84%		29-129%		

(a) Analysis performed at Accutest Laboratories, Marlborough, MA.

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ND = Not detected	MDL = Method Detection Limit	J = Indicates an estimated value
RL = Reporting Limit		B = Indicates analyte found in associated method blank
E = Indicates value exceeds calibration range		N = Indicates presumptive evidence of a compound

4.2  
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## Report of Analysis

<b>Client Sample ID:</b> BPOW 6-3	<b>Date Sampled:</b> 09/08/15
<b>Lab Sample ID:</b> JC3373-3	<b>Date Received:</b> 09/09/15
<b>Matrix:</b> AQ - Ground Water	<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1	
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY	

Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B99317.D	1	09/14/15	MD	n/a	n/a	V1B4710
Run #2							

Run #	Purge Volume
Run #1	5.0 ml
Run #2	

## Special VOA List

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

ND = Not detected MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

## Report of Analysis

<b>Client Sample ID:</b> BPOW 6-3		<b>Date Sampled:</b> 09/08/15
<b>Lab Sample ID:</b> JC3373-3		<b>Date Received:</b> 09/09/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	101%		78-114%
460-00-4	4-Bromofluorobenzene	85%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

ND = Not detected      MDL = Method Detection Limit  
 RL = Reporting Limit  
 E = Indicates value exceeds calibration range

J = Indicates an estimated value  
 B = Indicates analyte found in associated method blank  
 N = Indicates presumptive evidence of a compound

4.3  
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## Report of Analysis

<b>Client Sample ID:</b> BPOW 6-3	<b>Date Sampled:</b> 09/08/15
<b>Lab Sample ID:</b> JC3373-3	<b>Date Received:</b> 09/09/15
<b>Matrix:</b> AQ - Ground Water	<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8270D BY SIM SW846 3510C	
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY	

Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 <sup>a</sup>	I97871.D	1	09/22/15	AMA	09/11/15	M:OP44578	M:MSI3660
Run #2							

Run #	Initial Volume	Final Volume
Run #1	910 ml	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
123-91-1	1,4-Dioxane	ND	0.22	0.084	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	69%		26-121%		
321-60-8	2-Fluorobiphenyl	54%		28-107%		
1718-51-0	Terphenyl-d14	75%		29-129%		

(a) Analysis performed at Accutest Laboratories, Marlborough, MA.

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ND = Not detected	MDL = Method Detection Limit	J = Indicates an estimated value
RL = Reporting Limit		B = Indicates analyte found in associated method blank
E = Indicates value exceeds calibration range		N = Indicates presumptive evidence of a compound

4.3  
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## **Northrop Grumman Corporation- Operable Unit 2**

### **Data Review**

BETHPAGE, NEW YORK

Volatile and Semivolatile Analyses

SDGs #JC3500, JC3649 and JC3705

Analyses Performed By:  
Accutest Laboratories  
Dayton, New Jersey

Report #24354R  
October 8, 2015  
Review Level: Tier II  
Project #NY001496.1514.NAVI4

## SUMMARY

This data quality assessment summarizes the review of Sample Delivery Groups (SDGs) #JC3500, JC3649 and JC3705 for samples collected in association with the Northrop Grumman-Bethpage Site. The review was conducted as a Tier II evaluation and included review of data package completeness. Only analytical data associated with constituents of concern were reviewed for this validation. Field documentation was not included in this review. Included with this assessment are the validation annotated sample result sheets, and chain of custody. Analyses were performed on the following samples:

SDGs	Sample ID	Lab ID	Matrix	Sample Collection Date	Parent Sample	Analysis				
						VOC	SVOC	PCB	MET	MISC
JC3500	TB090915BT1	JC3500-1	Water	09/09/2015		X				
	FB090915BT1	JC3500-2	Water	09/09/2015		X	X			
	BPOW 6-4	JC3500-3	Water	09/09/2015		X	X			
	BPOW 6-4R	JC3500-4	Water	09/09/2015		X	X			
JC3649	TB091015BT1	JC3649-1	Water	09/10/2015		X				
	FB091015BT1	JC3649-2	Water	09/10/2015		X	X			
	BPOW 6-5	JC3649-3	Water	09/10/2015		X	X			
JC3705	FB091115BT1	JC3705-1	Water	09/11/2015		X	X			
	TB091115BT1	JC3705-2	Water	09/11/2015		X				
	BPOW 6-6	JC3705-3	Water	09/11/2015		X	X			

Notes:

1. Semivolatile (1,4-Dioxane) analysis was performed at Accutest Laboratories, Marlborough, MA (Accutest-NE) facility.
2. The matrix spike (MS) analysis was performed on sample location BPOW 6-4 for VOC.

## ANALYTICAL DATA PACKAGE DOCUMENTATION

### GENERAL INFORMATION

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

QA - Quality Assurance



## VOLATILE ANALYSIS INTRODUCTION

Analyses were performed according to United States Environmental Protection Agency (USEPA), Methods 524.2 and 8270D-Selected Ion Monitoring (SIM). Data were reviewed in accordance with USEPA National Functional Guidelines of October 1999.

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

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

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

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

## VOLATILE ORGANIC COMPOUND (VOC) ANALYSES

### 1. Holding Times

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

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

s.u. Standard units

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

### 2. Blank Contamination

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

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

Compounds were detected in the associated QA blanks; however, the associated sample results were greater than the BAL and/or were non-detect. No qualification of the sample results was required.

### 3. Surrogates/System Monitoring Compounds

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

All surrogate recoveries were within control limits.

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

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

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

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

## **5. Laboratory Control Sample (LCS) Analysis**

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

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

## **6. Field Duplicate Analysis**

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

A field duplicate was not collected on a sample location associated with these SDGs.

## **7. Laboratory Duplicate Analysis**

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

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

## **8. System Performance and Overall Assessment**

Tentatively identified compounds (TICs) were not identified in any of the sample locations.

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

## DATA VALIDATION CHECKLIST FOR VOCs

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

%R Percent Recovery    RPD Relative Percent Difference

## SEMIVOLATILE ORGANIC COMPOUND (SVOC) ANALYSES

### 1. Holding Times

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

Method	Matrix	Holding Time	Preservation
SW-846 8270D SIM	Water	7 days from collection to extraction and 40 days from extraction to analysis	Cooled to <6°C

All samples were analyzed within the specified holding time criteria.

### 2. Blank Contamination

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

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

Compounds were not detected above the MDL in the associated blanks; therefore detected sample results were not associated with blank contamination.

### 3. Surrogates/System Monitoring Compounds

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

All surrogate recoveries were within control limits.

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

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

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

A MS/MSD analysis was not performed on a sample location associated with these SDGs.

## **5. Laboratory Control Sample/Laboratory Control Sample Duplicate (LCS/LCSD) Analysis**

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

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

## **6. Field Duplicate Analysis**

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

A field duplicate was not collected with a sample location associated with these SDGs.

## **7. System Performance and Overall Assessment**

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

## DATA VALIDATION CHECKLIST FOR SVOCs

SVOCs: SW-846 8270D-SIM	Reported		Performance Acceptable		Not Required
	No	Yes	No	Yes	
GAS CHROMATOGRAPHY/MASS SPECTROMETRY (GC/MS)					
<b>Tier II Validation</b>					
Holding times		X		X	
Reporting limits (units)		X		X	
Blanks					
A. Method blanks		X		X	
B. Equipment blanks		X		X	
Laboratory Control Sample (LCS) %R		X		X	
Laboratory Control Sample Duplicate(LCSD) %R		X		X	
LCS/LCSD Precision (RPD)		X		X	
Matrix Spike (MS) %R					X
Matrix Spike Duplicate(MSD) %R					X
MS/MSD Precision (RPD)					X
Field/Lab Duplicate (RPD)					X
Surrogate Spike Recoveries		X		X	
Dilution Factor		X		X	
Moisture Content					X

%R    Percent recovery                      RPD    Relative percent difference

VALIDATION PERFORMED BY: Lisa Horton

SIGNATURE:   
\_\_\_\_\_

DATE: October 8, 2015

PEER REVIEW BY: Todd Church

DATE: October 11, 2015



**CHAIN OF CUSTODY/  
CORRECTED SAMPLE ANALYSIS DATA SHEETS**

GW  
FB  
WTB

**CHAIN OF CUSTODY**  
Accutest New Jersey/SPL Environmental  
2235 Route 130, Dayton, NJ 08810  
TEL: 732-329-0200 FAX: 732-329-3499/3480  
www.accutest.com

<b>FED-EX Tracking #</b> #4 <b>Accutest Quote #</b>		<b>Boiler Order Control #</b> Accutest Job # <b>JC3500</b>	
<b>Client / Reporting Information</b> Company Name: <b>Arcadis</b> Street Address: <b>2 Huntington Quad, Suite 1S10</b> City: <b>Melville NY</b> State: <b>NY</b> Zip: <b>11747</b> Project Contact: <b>Soma Das, soma.das@arcadis-us.com</b> Phone #: <b>631-391-5247</b> Fax #: <b>631-249-7610</b> Samplers (Name(s)): <b>PATRICIA PREZORSKI, BRENDAN TONG</b> Phone #: <b>516-287-6247</b>		<b>Project Information</b> Project Name: <b>AGMNYM62235 // OU2 Outpost Wells Northrop Grumman OU2 Hydro</b> Street: _____ Billing Information (if different from Report to): Company Name: <b>Arcadis, U.S., Inc. Attn: Accts Payable</b> Street Address: <b>630 Plaza Drive, Suite 600</b> City: <b>Highlands Ranch, CO</b> State: _____ Zip: <b>80129</b> Client Purchase Order #: _____ Work Authorization #: <b>NY001496_2015</b> Project Manager: <b>Carlo San Giovanni</b> Attention: <b>Soma Das</b>	
<b>Requested Analysis (see TEST CODE sheet)</b> V5242NG FULL LIST B8270SIM14DIOX		<b>Matrix Codes</b> DW - Drinking Water GW - Ground Water WW - Water SW - Surface Water SO - Soil SL - Sludge SED - Sediment OI - Oil LIQ - Other Liquid AIR - Air SOL - Other Solid WP - Wipe FB - Field Blank EB - Equipment Blank RB - Rinse Blank TB - Trip Blank	
<b>Accutest Sample #</b> 1 2 3 4	<b>Field ID / Point of Collection</b> TB090915BT1 FB090915BT1 BROW G-4 BROW G-4R	<b>Collection</b> MECH/DI Vial # Date Time Sampled by Matrix # of bottles HCl HNO3 H2O2 P2O5 NONE DI Water MEOH ENDOXINE	<b>Number of preserved bottles</b> HCl HNO3 H2O2 P2O5 NONE DI Water MEOH ENDOXINE
<b>Turnaround Time (Business days)</b> <input type="checkbox"/> Std. 15 Business Days <input checked="" type="checkbox"/> Std. 10 Business Days (by Contract only) <input type="checkbox"/> 10 Day RUSH <input type="checkbox"/> 5 Day RUSH <input type="checkbox"/> 3 Day EMERGENCY <input type="checkbox"/> 2 Day EMERGENCY <input type="checkbox"/> 1 Day EMERGENCY Emergency & Rush TIA data available VIA Lablink		<b>Data Deliverable Information</b> <input type="checkbox"/> Commercial "A" (Level 1) <input type="checkbox"/> NYASP Category A <input type="checkbox"/> Commercial "B" (Level 2) <input type="checkbox"/> NYASP Category B <input type="checkbox"/> FULLT1 (Level 3+4) <input type="checkbox"/> State Forms <input type="checkbox"/> NJ Reduced <input type="checkbox"/> EDD Format <input type="checkbox"/> Commercial "C" <input checked="" type="checkbox"/> Other COMMC+ Commercial "A" = Results Only Commercial "B" = Results + QC Summary NJ Reduced = Results + QC Summary + Partial Raw data	
<b>Approved By (Accutest PM): / Date:</b> _____ INITIAL ASSESSMENT <b>AL IA</b> LABEL VERIFICATION <b>OB</b>		<b>Comments / Special Instructions</b> OU2 Hydro Analyte List (V5242NG -FULL LIST) plus 1,4-Dioxane (B8270SIM14DIOX) Please use "_____" for MS/MSD QA/QC Sample.	
<b>Sample Custody must be documented below each time samples change possession, including courier delivery.</b>			
Relinquished by Sampler: 1 Patricia Prezorski 3 Brendan Tong 5 Relinquished by:	Date Time: 9/9/15 2:00 9/9/15 11:30	Received By: 1 [Signature] 2 [Signature] 3 [Signature] 4 [Signature] 5 [Signature]	Date Time: 9/10/15 19:00 9/10/15 19:00
Custody Seal # _____ <input type="checkbox"/> Intact <input type="checkbox"/> Not Intact		Preserved where applicable <input type="checkbox"/> On Ice <input type="checkbox"/>	

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JC3500: Chain of Custody

Page 1 of 3

**Report of Analysis**

<b>Client Sample ID:</b> TB090915BT1		<b>Date Sampled:</b> 09/09/15
<b>Lab Sample ID:</b> JC3500-1		<b>Date Received:</b> 09/10/15
<b>Matrix:</b> AQ - Trip Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

	<b>File ID</b>	<b>DF</b>	<b>Analyzed</b>	<b>By</b>	<b>Prep Date</b>	<b>Prep Batch</b>	<b>Analytical Batch</b>
Run #1	1B99313.D	1	09/14/15	MD	n/a	n/a	V1B4710
Run #2							

	<b>Purge Volume</b>
Run #1	5.0 ml
Run #2	

**Special VOA List**

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

ND = Not detected

MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

## Report of Analysis

<b>Client Sample ID:</b> TB090915BT1		<b>Date Sampled:</b> 09/09/15
<b>Lab Sample ID:</b> JC3500-1		<b>Date Received:</b> 09/10/15
<b>Matrix:</b> AQ - Trip Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	99%		78-114%
460-00-4	4-Bromofluorobenzene	88%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

ND = Not detected      MDL = Method Detection Limit  
 RL = Reporting Limit  
 E = Indicates value exceeds calibration range

J = Indicates an estimated value  
 B = Indicates analyte found in associated method blank  
 N = Indicates presumptive evidence of a compound

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4

## Report of Analysis

<b>Client Sample ID:</b> FB090915BT1		<b>Date Sampled:</b> 09/09/15
<b>Lab Sample ID:</b> JC3500-2		<b>Date Received:</b> 09/10/15
<b>Matrix:</b> AQ - Field Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B99314.D	1	09/14/15	MD	n/a	n/a	V1B4710
Run #2							

Run #1	Purge Volume
Run #1	5.0 ml
Run #2	

## Special VOA List

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

ND = Not detected

MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

## Report of Analysis

<b>Client Sample ID:</b> FB090915BT1		<b>Date Sampled:</b> 09/09/15
<b>Lab Sample ID:</b> JC3500-2		<b>Date Received:</b> 09/10/15
<b>Matrix:</b> AQ - Field Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	104%		78-114%
460-00-4	4-Bromofluorobenzene	88%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

---

ND = Not detected      MDL = Method Detection Limit      J = Indicates an estimated value  
 RL = Reporting Limit      B = Indicates analyte found in associated method blank  
 E = Indicates value exceeds calibration range      N = Indicates presumptive evidence of a compound

4.2  
4

## Report of Analysis

<b>Client Sample ID:</b> FB090915BT1	<b>Date Sampled:</b> 09/09/15
<b>Lab Sample ID:</b> JC3500-2	<b>Date Received:</b> 09/10/15
<b>Matrix:</b> AQ - Field Blank Water	<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8270D BY SIM SW846 3510C	
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY	

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 <sup>a</sup>	I97890.D	1	09/22/15	AMA	09/12/15	M:OP44598	M:MSI3660
Run #2							

	Initial Volume	Final Volume
Run #1	850 ml	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
123-91-1	1,4-Dioxane	ND	0.24	0.089	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	74%		26-121%		
321-60-8	2-Fluorobiphenyl	70%		28-107%		
1718-51-0	Terphenyl-d14	86%		29-129%		

(a) Analysis performed at Accutest Laboratories, Marlborough, MA.

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ND = Not detected	MDL = Method Detection Limit	J = Indicates an estimated value
RL = Reporting Limit		B = Indicates analyte found in associated method blank
E = Indicates value exceeds calibration range		N = Indicates presumptive evidence of a compound

4.2  
4

## Report of Analysis

<b>Client Sample ID:</b>	BPOW 6-4	<b>Date Sampled:</b>	09/09/15
<b>Lab Sample ID:</b>	JC3500-3	<b>Date Received:</b>	09/10/15
<b>Matrix:</b>	AQ - Ground Water	<b>Percent Solids:</b>	n/a
<b>Method:</b>	EPA 524.2 REV 4.1		
<b>Project:</b>	Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B99311.D	1	09/14/15	MD	n/a	n/a	V1B4710
Run #2							

Run #	Purge Volume
Run #1	5.0 ml
Run #2	

## Special VOA List

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

ND = Not detected

MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound



## Report of Analysis

<b>Client Sample ID:</b> BPOW 6-4		<b>Date Sampled:</b> 09/09/15
<b>Lab Sample ID:</b> JC3500-3		<b>Date Received:</b> 09/10/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	104%		78-114%
460-00-4	4-Bromofluorobenzene	91%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

ND = Not detected      MDL = Method Detection Limit  
 RL = Reporting Limit  
 E = Indicates value exceeds calibration range

J = Indicates an estimated value  
 B = Indicates analyte found in associated method blank  
 N = Indicates presumptive evidence of a compound

4.3  
4

## Report of Analysis

<b>Client Sample ID:</b> BPOW 6-4	<b>Date Sampled:</b> 09/09/15
<b>Lab Sample ID:</b> JC3500-3	<b>Date Received:</b> 09/10/15
<b>Matrix:</b> AQ - Ground Water	<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8270D BY SIM SW846 3510C	
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY	

Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 <sup>a</sup>	I97891.D	1	09/22/15	AMA	09/12/15	M:OP44598	M:MSI3660
Run #2							

Run #	Initial Volume	Final Volume
Run #1	910 ml	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
123-91-1	1,4-Dioxane	ND	0.22	0.084	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	81%		26-121%		
321-60-8	2-Fluorobiphenyl	69%		28-107%		
1718-51-0	Terphenyl-d14	81%		29-129%		

(a) Analysis performed at Accutest Laboratories, Marlborough, MA.

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ND = Not detected	MDL = Method Detection Limit	J = Indicates an estimated value
RL = Reporting Limit		B = Indicates analyte found in associated method blank
E = Indicates value exceeds calibration range		N = Indicates presumptive evidence of a compound

4.3  
4

## Report of Analysis

<b>Client Sample ID:</b> BPOW 6-4R		<b>Date Sampled:</b> 09/09/15
<b>Lab Sample ID:</b> JC3500-4		<b>Date Received:</b> 09/10/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #2	1B99312.D	1	09/14/15	MD	n/a	n/a	V1B4710

Run #1	Purge Volume
Run #2	5.0 ml

## Special VOA List

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

ND = Not detected      MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound

## Report of Analysis

<b>Client Sample ID:</b> BPOW 6-4R		<b>Date Sampled:</b> 09/09/15
<b>Lab Sample ID:</b> JC3500-4		<b>Date Received:</b> 09/10/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	101%		78-114%
460-00-4	4-Bromofluorobenzene	87%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

ND = Not detected      MDL = Method Detection Limit  
 RL = Reporting Limit  
 E = Indicates value exceeds calibration range

J = Indicates an estimated value  
 B = Indicates analyte found in associated method blank  
 N = Indicates presumptive evidence of a compound

4.4  
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## Report of Analysis

<b>Client Sample ID:</b> BPOW 6-4R		<b>Date Sampled:</b> 09/09/15
<b>Lab Sample ID:</b> JC3500-4		<b>Date Received:</b> 09/10/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8270D BY SIM SW846 3510C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 <sup>a</sup>	I97892.D	1	09/22/15	AMA	09/12/15	M:OP44598	M:MSI3660
Run #2							

	Initial Volume	Final Volume
Run #1	880 ml	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
123-91-1	1,4-Dioxane	ND	0.23	0.086	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	78%		26-121%		
321-60-8	2-Fluorobiphenyl	71%		28-107%		
1718-51-0	Terphenyl-d14	84%		29-129%		

(a) Analysis performed at Accutest Laboratories, Marlborough, MA.

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ND = Not detected	MDL = Method Detection Limit	J = Indicates an estimated value
RL = Reporting Limit		B = Indicates analyte found in associated method blank
E = Indicates value exceeds calibration range		N = Indicates presumptive evidence of a compound

4.4  
4

GW  
FB  
WTB

CHAIN OF CUSTODY

Accutest New Jersey/SPL Environmental  
2235 Route 130, Dayton, NJ 08810  
TEL: 732-329-0200 FAX: 732-329-3499/3480  
www.accutest.com

WPNR

Client / Reporting Information		Project Information					Requested Analysis ( see TEST CODE sheet)										Matrix Codes	
<b>Company Name:</b> Arcadis <b>Street Address:</b> 2 Huntington Quad, Suite 1S10 City: Melville NY 11747 <b>Project Contact:</b> Soma Das, soma.das@arcadis-us.com Phone #: 631-391-5247 Fax #: 631-249-7610 Email: 516-497-6247 Sampler(s) Name(s): PATRICIA PREZORSKI, BRENDAN TONG		<b>Project Name:</b> AGMNYM62235 // OU2 Outpost Wells Northrop Grumman OU2 Hydro <b>Street:</b> <b>City:</b> Bethpage NY <b>Project #:</b> NY001496.1514.NAVI3 <b>Client Purchase Order #:</b> NY001496.2015 <b>Work Authorization #:</b> NY001496.2015 <b>Project Manager:</b> Carlo San Giovanni Attention: Soma Das					<b>Accutest Tracking #:</b> #4 <b>Accutest Quote #:</b> <b>Bottle Order Control #:</b> JC3649 <b>Accutest Job #:</b>										<b>Matrix Codes:</b> DW - Drinking Water GW - Ground Water WW - Water SW - Surface Water SO - Soil SL - Sludge SED - Sediment OI - Oil LIQ - Other Liquid AIR - Air SOL - Other Solid WP - Wipe FB - Field Blank EB - Equipment Blank RB - Rinse Blank TB - Trip Blank	
<b>Billing Information ( If different from Report to )</b>																		
<b>Company Name:</b> Arcadis, U.S., Inc. Attn: Accts Payable																		
<b>Street Address:</b> 630 Plaza Drive, Suite 600																		
<b>City:</b> Highlands Ranch, CO 80129																		
<b>State:</b> CO																		
<b>Zip:</b> 80129																		
<b>Attention:</b> Soma Das																		
<b>Collection</b>																		
<b>Field ID / Point of Collection</b>		<b>MECH/DOI Val #</b>		<b>Date</b>		<b>Time</b>		<b>Sampled by</b>		<b>Matrix</b>		<b># of bottles</b>		<b>Number of preserved bottles</b>		<b>LAB USE ONLY</b>		
1 TB09101587A				9/10/15		0830		- TB		2		2		V5242NG FULL LIST				
2 FB09101587A				9/10/15		0900		FB		2		2		B82705IM14DIOX				
3 BROW G-5				9/10/15		1225		GW		3		3				V30		
<b>Turnaround Time ( Business days )</b>		<b>Approved By (Accutest PM): / Date:</b>					<b>Data Deliverable Information</b>					<b>Comments / Special Instructions</b>						
<input type="checkbox"/> Std. 15 Business Days <input checked="" type="checkbox"/> Std. 10 Business Days ( by Contract only ) <input type="checkbox"/> 10 Day RUSH <input type="checkbox"/> 5 Day RUSH <input type="checkbox"/> 3 Day EMERGENCY <input type="checkbox"/> 2 Day EMERGENCY <input type="checkbox"/> 1 Day EMERGENCY Emergency & Rush T/A data available VIA Lablink		<b>INITIAL ASSESSMENT 2AOB</b> <b>LABEL VERIFICATION JK</b>					<input type="checkbox"/> Commercial "A" ( Level 1 ) <input type="checkbox"/> Commercial "B" ( Level 2 ) <input type="checkbox"/> FULLT1 ( Level 3+4 ) <input type="checkbox"/> NJ Reduced <input type="checkbox"/> Commercial "C" <input type="checkbox"/> NYASP Category A <input type="checkbox"/> NYASP Category B <input type="checkbox"/> State Forms <input type="checkbox"/> EDD Format <input checked="" type="checkbox"/> Other COMM+ Commercial "A" = Results Only Commercial "B" = Results + QC Summary NJ Reduced = Results + QC Summary + Partial Raw data					<b>OU2 Hydro Analyte List (V5242NG -FULL LIST) plus 1,4-Dioxane (B82705IM14DIOX)</b> Please use " " for MS/MSD QA/QC Sample.						
<b>Sample Custody must be documented below each time samples change possession, including courier delivery.</b>																		
<b>Relinquished by Sampler:</b>		<b>Date Time:</b>		<b>Received By:</b>		<b>Date Time:</b>		<b>Relinquished By:</b>		<b>Date Time:</b>		<b>Received By:</b>		<b>Relinquished By:</b>		<b>Date Time:</b>		
1 Patricia Prezorski		9/10/15 1545		[Signature]		[Signature]		2 [Signature]		4:25		2 [Signature]		[Signature]		[Signature]		
3 Brendan Tong				3 [Signature]				4 [Signature]				4 [Signature]						
5 [Signature]				5 [Signature]				5 [Signature]				5 [Signature]						
<input type="checkbox"/> Intact Preserved where applicable <input type="checkbox"/> Not Intact Cooler Temp. 42																		

5.1  
5

# Report of Analysis

<b>Client Sample ID:</b> TB091015BT1		<b>Date Sampled:</b> 09/10/15
<b>Lab Sample ID:</b> JC3649-1		<b>Date Received:</b> 09/11/15
<b>Matrix:</b> AQ - Trip Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B99372.D	1	09/15/15	MD	n/a	n/a	V1B4711
Run #2							

Run #1	Purge Volume
Run #1	5.0 ml
Run #2	

### Special VOA List

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

ND = Not detected      MDL = Method Detection Limit      J = Indicates an estimated value  
 RL = Reporting Limit      B = Indicates analyte found in associated method blank  
 E = Indicates value exceeds calibration range      N = Indicates presumptive evidence of a compound

4.1  
 4

## Report of Analysis

<b>Client Sample ID:</b> TB091015BT1		<b>Date Sampled:</b> 09/10/15
<b>Lab Sample ID:</b> JC3649-1		<b>Date Received:</b> 09/11/15
<b>Matrix:</b> AQ - Trip Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	103%		78-114%
460-00-4	4-Bromofluorobenzene	88%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

ND = Not detected      MDL = Method Detection Limit  
 RL = Reporting Limit  
 E = Indicates value exceeds calibration range

J = Indicates an estimated value  
 B = Indicates analyte found in associated method blank  
 N = Indicates presumptive evidence of a compound

4.1  
4



## Report of Analysis

<b>Client Sample ID:</b> FB091015BT1	
<b>Lab Sample ID:</b> JC3649-2	<b>Date Sampled:</b> 09/10/15
<b>Matrix:</b> AQ - Field Blank Water	<b>Date Received:</b> 09/11/15
<b>Method:</b> EPA 524.2 REV 4.1	<b>Percent Solids:</b> n/a
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY	

Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B99420.D	1	09/17/15	MD	n/a	n/a	V1B4712
Run #2							

Run #	Purge Volume
Run #1	5.0 ml
Run #2	

## Special VOA List

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

ND = Not detected MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

## Report of Analysis

<b>Client Sample ID:</b> FB091015BT1		<b>Date Sampled:</b> 09/10/15
<b>Lab Sample ID:</b> JC3649-2		<b>Date Received:</b> 09/11/15
<b>Matrix:</b> AQ - Field Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	100%		78-114%
460-00-4	4-Bromofluorobenzene	86%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

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ND = Not detected      MDL = Method Detection Limit      J = Indicates an estimated value  
 RL = Reporting Limit      B = Indicates analyte found in associated method blank  
 E = Indicates value exceeds calibration range      N = Indicates presumptive evidence of a compound

4.2  
4

## Report of Analysis

<b>Client Sample ID:</b> FB091015BT1		<b>Date Sampled:</b> 09/10/15
<b>Lab Sample ID:</b> JC3649-2		<b>Date Received:</b> 09/11/15
<b>Matrix:</b> AQ - Field Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8270D BY SIM SW846 3510C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 <sup>a</sup>	I97783.D	1	09/17/15	AMA	09/13/15	M:OP44602	M:MSI3657
Run #2							

	Initial Volume	Final Volume
Run #1	800 ml	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
123-91-1	1,4-Dioxane	ND	0.25	0.095	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	87%		26-121%		
321-60-8	2-Fluorobiphenyl	73%		28-107%		
1718-51-0	Terphenyl-d14	90%		29-129%		

(a) Analysis performed at Accutest Laboratories, Marlborough, MA.

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ND = Not detected	MDL = Method Detection Limit	J = Indicates an estimated value
RL = Reporting Limit		B = Indicates analyte found in associated method blank
E = Indicates value exceeds calibration range		N = Indicates presumptive evidence of a compound

4.2  
4

## Report of Analysis

<b>Client Sample ID:</b> BPOW6-5	<b>Date Sampled:</b> 09/10/15
<b>Lab Sample ID:</b> JC3649-3	<b>Date Received:</b> 09/11/15
<b>Matrix:</b> AQ - Ground Water	<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1	
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY	

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B99421.D	1	09/17/15	MD	n/a	n/a	V1B4712
Run #2							

Run #1	Purge Volume
Run #1	5.0 ml
Run #2	

## Special VOA List

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	5.0	0.91	ug/l	
78-93-3	2-Butanone	ND	5.0	0.57	ug/l	
71-43-2	Benzene	ND	0.50	0.057	ug/l	
75-27-4	Bromodichloromethane	ND	0.50	0.082	ug/l	
75-25-2	Bromoform	ND	0.50	0.046	ug/l	
74-83-9	Bromomethane	ND	0.50	0.077	ug/l	
75-15-0	Carbon disulfide	0.24	0.50	0.028	ug/l	J
108-90-7	Chlorobenzene	ND	0.50	0.027	ug/l	
75-00-3	Chloroethane	ND	0.50	0.037	ug/l	
67-66-3	Chloroform	ND	0.50	0.031	ug/l	
74-87-3	Chloromethane	ND	0.50	0.044	ug/l	
56-23-5	Carbon tetrachloride	ND	0.50	0.074	ug/l	
75-34-3	1,1-Dichloroethane	ND	0.50	0.039	ug/l	
75-35-4	1,1-Dichloroethylene	ND	0.50	0.054	ug/l	
107-06-2	1,2-Dichloroethane	ND	0.50	0.034	ug/l	
78-87-5	1,2-Dichloropropane	ND	0.50	0.082	ug/l	
124-48-1	Dibromochloromethane	ND	0.50	0.042	ug/l	
156-60-5	trans-1,2-Dichloroethylene	ND	0.50	0.039	ug/l	
156-59-2	cis-1,2-Dichloroethylene	ND	0.50	0.081	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.033	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.063	ug/l	
100-41-4	Ethylbenzene	ND	0.50	0.033	ug/l	
76-13-1	Freon 113	ND	1.0	0.10	ug/l	
591-78-6	2-Hexanone	ND	2.0	0.084	ug/l	
75-09-2	Methylene chloride	ND	0.50	0.047	ug/l	
108-10-1	4-Methyl-2-pentanone	ND	2.0	0.27	ug/l	
100-42-5	Styrene	ND	0.50	0.028	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	0.50	0.050	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.035	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	0.50	0.052	ug/l	
127-18-4	Tetrachloroethylene	ND	0.50	0.091	ug/l	
108-88-3	Toluene	ND	0.50	0.044	ug/l	

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound

## Report of Analysis

<b>Client Sample ID:</b> BPOW6-5		<b>Date Sampled:</b> 09/10/15
<b>Lab Sample ID:</b> JC3649-3		<b>Date Received:</b> 09/11/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	102%		78-114%
460-00-4	4-Bromofluorobenzene	87%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

ND = Not detected      MDL = Method Detection Limit  
 RL = Reporting Limit  
 E = Indicates value exceeds calibration range

J = Indicates an estimated value  
 B = Indicates analyte found in associated method blank  
 N = Indicates presumptive evidence of a compound

4.3  
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## Report of Analysis

<b>Client Sample ID:</b> BPOW6-5		<b>Date Sampled:</b> 09/10/15
<b>Lab Sample ID:</b> JC3649-3		<b>Date Received:</b> 09/11/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8270D BY SIM SW846 3510C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 <sup>a</sup>	I97784.D	1	09/17/15	AMA	09/13/15	M:OP44602	M:MSI3657
Run #2							

	Initial Volume	Final Volume
Run #1	920 ml	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
123-91-1	1,4-Dioxane	ND	0.22	0.083	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	70%		26-121%		
321-60-8	2-Fluorobiphenyl	63%		28-107%		
1718-51-0	Terphenyl-d14	88%		29-129%		

(a) Analysis performed at Accutest Laboratories, Marlborough, MA.

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ND = Not detected	MDL = Method Detection Limit	J = Indicates an estimated value
RL = Reporting Limit		B = Indicates analyte found in associated method blank
E = Indicates value exceeds calibration range		N = Indicates presumptive evidence of a compound

4.3  
4

GW  
FB  
WB

<b>Client / Reporting Information</b> Company Name: <b>Arcadis</b> Street Address: <b>2 Huntington Quad, Suite 1S10</b> City: <b>Melville NY 11747</b> Project Contact: <b>Soma Das, soma.das@arcadis-us.com</b> Phone #: <b>631-391-5247</b> Fax #: <b>631-249-7610</b> Sampler(s) Name(s): <b>PATRICIA PREZORSKI, BRENDAN TONG</b>		<b>Project Information</b> Project Name: <b>AGMNYM62235 // OU2 Outpost Wells Northrop Grumman OU2 Hydro</b> Street: _____ Billing Information (if different from Report to): Company Name: <b>Arcadis, U.S., Inc. Attn: Accts Payable</b> Street Address: <b>630 Plaza Drive, Suite 600</b> City: <b>Highlands Ranch, CO 80129</b> Attention: <b>Soma Das</b>		FED-EX Tracking # <b>#4</b> Accutest Quote # _____ Bottle Order Control # _____ Accutest Job # <b>JC3705</b>											
Requested Analysis (see TEST CODE sheet) <b>V5242NG FULL LIST</b> <b>B8270SIM14DIOX</b>		Matrix Codes DW - Drinking Water GW - Ground Water WW - Water SW - Surface Water SO - Soil SL - Sludge SED - Sediment OI - Oil LIQ - Other Liquid AIR - Air SOL - Other Solid WP - Wipe FB - Field Blank EB - Equipment Blank RB - Rinse Blank TB - Trip Blank		LAB USE ONLY <b>V30</b>											
Accutest Sample # <b>1</b> <b>2</b> <b>3</b>	Field ID / Point of Collection <b>FB091158T1</b> <b>TB091158T1</b> <b>BPOW 6-6</b>	MECH/DI Val #  	Date <b>9/11/15</b> <b>9/11/15</b> <b>9/11/15</b>	Time <b>0840</b> <b>0830</b> <b>1230</b>	Sampled by <b>GP</b> <b>GP</b> <b>GP</b>	Matrix <b>FB</b> <b>TB</b> <b>GW</b>	# of bottles <b>2</b> <b>2</b> <b>3</b>	FCI <b>2</b> <b>2</b> <b>3</b>	NIENH  	FNCO3  	HPSO4  	NONE  	DI Value  	MECH  	ENCORE  
Turnaround Time (Business days) <input type="checkbox"/> Std. 15 Business Days <input checked="" type="checkbox"/> Std. 10 Business Days (by Contract only) <input type="checkbox"/> 10 Day RUSH <input type="checkbox"/> 5 Day RUSH <input type="checkbox"/> 3 Day EMERGENCY <input type="checkbox"/> 2 Day EMERGENCY <input type="checkbox"/> 1 Day EMERGENCY Emergency & Rush T/A data available VIA Lablink		Approved By (Accutest PM) / Date: <b>INITIAL ASSESSMENT IB OB</b> <b>LABEL VERIFICATION JK</b>		Data Deliverable Information <input type="checkbox"/> Commercial "A" (Level 1) <input type="checkbox"/> Commercial "B" (Level 2) <input type="checkbox"/> FULLT1 (Level 3+4) <input type="checkbox"/> NJ Reduced <input type="checkbox"/> Commercial "C" <input type="checkbox"/> NYASP Category A <input type="checkbox"/> NYASP Category B <input type="checkbox"/> State Forms <input type="checkbox"/> EDD Format <input checked="" type="checkbox"/> Other <b>COMMC+</b> Commercial "A" = Results Only Commercial "B" = Results + QC Summary NJ Reduced = Results + QC Summary + Partial Raw data		Comments / Special Instructions <b>OU2 Hydro Analyte List (V5242NG -FULL LIST) plus 1,4-Dioxane (B8270SIM14DIOX)</b> Please use " " for MS/MSD QA/QC Sample.									
Relinquished by Sampler: <b>1 Patricia Prezorski</b> <b>3 Brendan Tong</b> <b>5</b>		Date Time: <b>9/11/15 1525</b> <b>9/11/15 15:25</b>		Received By: <b>Patrizia Prezorski</b> <b>Patrizia Prezorski</b>		Relinquished By: <b>Patrizia Prezorski</b> <b>Patrizia Prezorski</b>		Date Time: <b>9-11-2015</b> <b>9-11-2015</b>		Received By: <b>Patrizia Prezorski</b> <b>Patrizia Prezorski</b>					
Relinquished by:  		Date Time:  		Received By:  		Relinquished By:  		Date Time:  		Received By:  					
Custody Seal # _____ <input type="checkbox"/> Intact <input type="checkbox"/> Not intact		Preserved where applicable <input type="checkbox"/>		Cooler Temp. <b>42°C</b>											

5.1  
5

<b>FED-EX Tracking #</b> #4 <b>Accutest Quote #</b>		<b>Bottle Order Control #</b> <b>Accutest Job #</b> JC3705	
<b>Client / Reporting Information</b> Company Name: <b>Arcadis</b> Street Address: <b>2 Huntington Quad, Suite 1S10</b> City, State, Zip: <b>Melville NY 11747</b> Project Contact: <b>Soma Das, soma.das@arcadis-us.com</b> Phone #: <b>631-391-5247</b> Fax #: <b>631-249-7610</b> Sample(s) Name(s): <b>PATRICIA PRZORSKI, BRENDAN TONG</b> Phone #: <b>516 287-6247</b>		<b>Project Information</b> Project Name: <b>AGMNYM62235 // OU2 Outpost Wells</b> Street: <b>Northrop Grumman OU2 Hydro</b> Billing Information (if different from Report to): Company Name: <b>Arcadis, U.S., Inc. Attn: Accts Payable</b> Street Address: <b>630 Plaza Drive, Suite 600</b> City, State, Zip: <b>Highlands Ranch, CO 80129</b> Client Purchase Order #: <b>NY001496.1514.NAVI3</b> Work Authorization #: <b>NY001496_2015</b> Project Manager: <b>Carlo San Giovanni</b> Attention: <b>Soma Das</b>	
<b>Requested Analysis (see TEST CODE sheet)</b> V5242NG FULL LIST B8270SIM14DIOX		<b>Matrix Codes</b> DW - Drinking Water GW - Ground Water WW - Water SW - Surface Water SO - Soil SL - Sludge SED - Sediment OI - Oil LIQ - Other Liquid AIR - Air SOL - Other Solid WP - Waste FB - Field Blank EB - Equipment Blank RB - Rinse Blank TB - Trip Blank	
<b>Turnaround Time (Business days)</b> <input type="checkbox"/> Std. 15 Business Days <input checked="" type="checkbox"/> Std. 10 Business Days (by Contract only) <input type="checkbox"/> 10 Day RUSH <input type="checkbox"/> 5 Day RUSH <input type="checkbox"/> 3 Day EMERGENCY <input type="checkbox"/> 2 Day EMERGENCY <input type="checkbox"/> 1 Day EMERGENCY Emergency & Rush TIA date available via Lablink		<b>Data Deliverable Information</b> <input type="checkbox"/> Commercial "A" (Level 1) <input type="checkbox"/> NYASP Category A <input type="checkbox"/> Commercial "B" (Level 2) <input type="checkbox"/> NYASP Category B <input type="checkbox"/> FULLT1 (Level 3+4) <input type="checkbox"/> State Forms <input type="checkbox"/> NJ Reduced <input type="checkbox"/> EDD Format <input type="checkbox"/> Commercial "C" <input checked="" type="checkbox"/> Other <b>COMMC+</b> Commercial "A" = Results Only Commercial "B" = Results + QC Summary NJ Reduced = Results + QC Summary + Partial Raw data	
<b>Comments / Special Instructions</b> OU2 Hydro Analyte List (V5242NG-FULL LIST) plus 1,4-Dioxane (B8270SIM14DIOX) Please use * <i>BP</i> * for MS/MSD QA/QC Sample.			
<b>Sample Custody must be documented below each time samples change possession, including courier delivery.</b>			
Relinquished by Sampler: <b>1 Patricia Przorski</b>	Date/Time: <b>9/11/15 1500</b>	Received By: <b>FX</b>	Relinquished By: <b>FX</b>
Relinquished by Sampler: <b>3 Brendan Tong</b>	Date/Time:	Received By:	Relinquished By:
Relinquished by:	Date/Time:	Received By:	Relinquished By:
Emergency & Rush TIA date available via Lablink		On Ice <input type="checkbox"/> Cooler Temp. <b>16°C</b>	

5.1  
5

17A



## Report of Analysis

<b>Client Sample ID:</b> FB091115BT1		<b>Date Sampled:</b> 09/11/15
<b>Lab Sample ID:</b> JC3705-1		<b>Date Received:</b> 09/11/15
<b>Matrix:</b> AQ - Field Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B99369.D	1	09/15/15	MD	n/a	n/a	V1B4711
Run #2							

Run #	Purge Volume
Run #1	5.0 ml
Run #2	

## Special VOA List

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

ND = Not detected MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

## Report of Analysis

<b>Client Sample ID:</b> FB091115BT1		<b>Date Sampled:</b> 09/11/15
<b>Lab Sample ID:</b> JC3705-1		<b>Date Received:</b> 09/11/15
<b>Matrix:</b> AQ - Field Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	103%		78-114%
460-00-4	4-Bromofluorobenzene	89%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

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ND = Not detected      MDL = Method Detection Limit      J = Indicates an estimated value  
 RL = Reporting Limit      B = Indicates analyte found in associated method blank  
 E = Indicates value exceeds calibration range      N = Indicates presumptive evidence of a compound

4.1  
4

## Report of Analysis

<b>Client Sample ID:</b> FB091115BT1	<b>Date Sampled:</b> 09/11/15
<b>Lab Sample ID:</b> JC3705-1	<b>Date Received:</b> 09/11/15
<b>Matrix:</b> AQ - Field Blank Water	<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8270D BY SIM SW846 3510C	
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY	

Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 <sup>a</sup>	I97753.D	1	09/16/15	AMA	09/14/15	M:OP44621	M:MSI3656
Run #2							

Run #	Initial Volume	Final Volume
Run #1	850 ml	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
123-91-1	1,4-Dioxane	ND	0.24	0.089	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	71%		26-121%		
321-60-8	2-Fluorobiphenyl	65%		28-107%		
1718-51-0	Terphenyl-d14	81%		29-129%		

(a) Analysis performed at Accutest Laboratories, Marlborough, MA.

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ND = Not detected	MDL = Method Detection Limit	J = Indicates an estimated value
RL = Reporting Limit		B = Indicates analyte found in associated method blank
E = Indicates value exceeds calibration range		N = Indicates presumptive evidence of a compound

4.1  
4

### Report of Analysis

<b>Client Sample ID:</b> TB091115BT1	<b>Date Sampled:</b> 09/11/15
<b>Lab Sample ID:</b> JC3705-2	<b>Date Received:</b> 09/11/15
<b>Matrix:</b> AQ - Trip Blank Water	<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1	
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY	

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B99370.D	1	09/15/15	MD	n/a	n/a	V1B4711
Run #2							

Run #1	Purge Volume
Run #1	5.0 ml
Run #2	

**Special VOA List**

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

ND = Not detected      MDL = Method Detection Limit      J = Indicates an estimated value  
 RL = Reporting Limit      B = Indicates analyte found in associated method blank  
 E = Indicates value exceeds calibration range      N = Indicates presumptive evidence of a compound

4.2  
 4

## Report of Analysis

<b>Client Sample ID:</b> TB091115BT1		<b>Date Sampled:</b> 09/11/15
<b>Lab Sample ID:</b> JC3705-2		<b>Date Received:</b> 09/11/15
<b>Matrix:</b> AQ - Trip Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	103%		78-114%
460-00-4	4-Bromofluorobenzene	89%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

ND = Not detected      MDL = Method Detection Limit  
 RL = Reporting Limit  
 E = Indicates value exceeds calibration range

J = Indicates an estimated value  
 B = Indicates analyte found in associated method blank  
 N = Indicates presumptive evidence of a compound

4.2  
4

## Report of Analysis

<b>Client Sample ID:</b> BPOW6-6		<b>Date Sampled:</b> 09/11/15
<b>Lab Sample ID:</b> JC3705-3		<b>Date Received:</b> 09/11/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B99371.D	1	09/15/15	MD	n/a	n/a	V1B4711
Run #2							

Run #	Purge Volume
Run #1	5.0 ml
Run #2	

## Special VOA List

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

ND = Not detected MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

## Report of Analysis

<b>Client Sample ID:</b> BPOW6-6		<b>Date Sampled:</b> 09/11/15
<b>Lab Sample ID:</b> JC3705-3		<b>Date Received:</b> 09/11/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> EPA 524.2 REV 4.1		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**Special VOA List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethylene	ND	0.50	0.024	ug/l	
75-01-4	Vinyl chloride	ND	0.50	0.032	ug/l	
	m,p-Xylene	ND	0.50	0.13	ug/l	
95-47-6	o-Xylene	ND	0.50	0.029	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2199-69-1	1,2-Dichlorobenzene-d4	102%		78-114%
460-00-4	4-Bromofluorobenzene	87%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

ND = Not detected      MDL = Method Detection Limit  
 RL = Reporting Limit  
 E = Indicates value exceeds calibration range

J = Indicates an estimated value  
 B = Indicates analyte found in associated method blank  
 N = Indicates presumptive evidence of a compound

4.3  
4

## Report of Analysis

<b>Client Sample ID:</b> BPOW6-6	<b>Date Sampled:</b> 09/11/15
<b>Lab Sample ID:</b> JC3705-3	<b>Date Received:</b> 09/11/15
<b>Matrix:</b> AQ - Ground Water	<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8270D BY SIM SW846 3510C	
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY	

Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 <sup>a</sup>	I97754.D	1	09/16/15	AMA	09/14/15	M:OP44621	M:MSI3656
Run #2							

Run #	Initial Volume	Final Volume
Run #1	900 ml	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
123-91-1	1,4-Dioxane	ND	0.22	0.084	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	64%		26-121%		
321-60-8	2-Fluorobiphenyl	60%		28-107%		
1718-51-0	Terphenyl-d14	72%		29-129%		

(a) Analysis performed at Accutest Laboratories, Marlborough, MA.

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ND = Not detected	MDL = Method Detection Limit	J = Indicates an estimated value
RL = Reporting Limit		B = Indicates analyte found in associated method blank
E = Indicates value exceeds calibration range		N = Indicates presumptive evidence of a compound

4.3  
4



**Appendix D**

***ARCADIS Report: Results of Third Quarter  
2015 Groundwater Monitoring (ARCADIS, 2015)***

Mr. Henry Wilkie  
Project Manager  
New York State Department of Environmental Conservation  
Remedial Bureau A  
625 Broadway  
Albany, New York 12233-7015

Mr. Steven Scharf, P.E.  
Project Manager  
New York State Department of Environmental Conservation  
Remedial Bureau A  
625 Broadway  
Albany, New York 12233-7015

Subject:

Results of Third Quarter 2015 Groundwater Monitoring,  
Operable Unit 2, Northrop Grumman Systems Corporation and Naval Weapons  
Industrial Reserve Plant (NWIRP) Sites, Bethpage, New York.  
(NYSDEC Site #s 1-30-003A and B)

Dear Mr. Scharf:

On behalf of Northrop Grumman Systems Corporation (Northrop Grumman), ARCADIS is providing the NYSDEC with the validated results of Operable Unit 2 (OU2) groundwater monitoring, performed in accordance with the approved Groundwater Monitoring Plan (ARCADIS of New York, Inc. 2012) and the Public Water Supply Contingency Plan (PWSCP) (ARCADIS G&M, Inc. 2003). Table 1 summarizes OU2 remedial system performance operational data and water balance. Tables 2 and 3 provide the validated analytical results of monitoring for this period. Table 4, 5 and 6 provide the validated analytical results for vapor samples collected from the system for this period. Figure 1 shows the site plan with well locations.

Please contact us if you have any questions or comments.

Sincerely,

Arcadis of New York, Inc.



David E. Stern  
Senior Hydrogeologist

Enclosures

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ENVIRONMENT

Date:

November 30, 2015

Contact:

David Stern

Phone:

631.391.5284

Email:

[david.stern@arcadis.com](mailto:david.stern@arcadis.com)

Our ref:

NY001496. 314I.NAVI4

Mr. Henry Wilkie  
Mr. Steven Scharf, P.E.  
November 30, 2015

Copies:

Ed Hannon – Northrop Grumman  
Fred Weber – Northrop Grumman  
Walter Parish – NYSDEC Region 1  
Steven Karpinski – New York State Department of Health  
Michael Alarcon – Nassau County Department of Health  
Joseph DeFranco – Nassau County Department of Health  
Lora Fly – NAVFAC Midlant Environmental  
David Brayack – TetraTech NUS, Inc.  
Roger Smith – Glenn Springs Holdings, Inc.  
Kevin Lumpe – Steel Equities  
Thomas Taccone – USEPA  
Robert Alvey – USEPA  
Carol Stein-USEPA  
Matthew Russo – Town of Oyster Bay  
Stan Carey – Massapequa Water District  
Matthew Snyder – New York American Water  
Charles Prucha – South Farmingdale Water District  
John Reinhardt – Town of Hempstead Water District  
Michael Boufis – Bethpage Water District  
Lois Lovisolo – Bethpage Public Library (Public Repository)  
File

# TABLES



**Table 1**  
**Operational Summary for the On-Site Portion of the Operable Unit 2 Groundwater Remedy**  
**Third Quarter 2015**  
**Northrop Grumman Systems Corporation, Operable Unit 2**  
**Bethpage, New York**

	Quarterly Flow Rates (gpm)		Quarterly Flow Volumes (MG)			Quarterly VOC Concentrations (µg/L)		VOC Mass Removed (lbs) <sup>(7)</sup>
	Design <sup>(2)</sup>	Average <sup>(3,4,17)</sup>	Design <sup>(2)</sup>	Actual <sup>(3,4)</sup>	% of Design	TCE <sup>(5)</sup>	TVOC <sup>(5,6)</sup>	Quarterly
<b><u>Influent Groundwater</u></b>								
Well 1 <sup>(11,13)</sup>	800	803	104.8	101.0	96%	852	900	760
Well 3R <sup>(12,13)</sup>	700	916	91.7	118.9	130%	557	620	603
Well 17 <sup>(14,16)</sup>	1,000	998	131.0	126.9	97%	178	220	228
Well 18 <sup>(16)</sup>	600	616	78.6	79.1	101%	61	83	54
Well 19 <sup>(15,16)</sup>	700	750	91.7	92.4	101%	153	180	136
<b>Total</b>	<b>3,800</b>	<b>4,083</b>	<b>498</b>	<b>518</b>	<b>104%</b>	<b>--</b>	<b>--</b>	<b>1,781</b>
<b><u>Effluent Groundwater</u> <sup>(8)</sup></b>								
Calpine	100 - 400	497	--	65.9	--	--	--	--
OXY Biosparge <sup>(10)</sup>	2 - 42	3.8	--	0.5	--	--	--	--
West Recharge Basins	1,112 - 1,455	852	--	111.6	--	--	0.90	--
South Recharge Basins	2,231	2,597	292.4	340.3	116%	--	1.18	--
<b>Total Effluent Groundwater</b>	<b>--</b>	<b>3,950</b>	<b>--</b>	<b>518</b>	<b>--</b>	<b>--</b>	<b>--</b>	<b>--</b>
<b><u>Additional Flow to South Recharge Basins</u></b>								
Storm Water Runoff Contributing to South Recharge Basins Flow Volume <sup>(18)</sup>				20.7				
<b>Total Flow Volume to South Recharge Basins</b> <sup>(19)</sup>			<b>292.4</b>	<b>361.0</b>	<b>123%</b>			
<b><u>Treatment Efficiencies</u> <sup>(9)</sup></b>								
Tower 96 System:	99.9%							
Tower 102 System:	>99.9%							

See notes on next page

**Table 1**  
**Operational Summary for the On-Site Portion of the Operable Unit 2 Groundwater Remedy**  
**Third Quarter 2015**  
**Northrop Grumman Systems Corporation, Operable Unit 2**  
**Bethpage, New York**

**Notes:**

- (1) Quarterly reporting period: July 06, 2015 through October 05, 2015
- (2) "Design" flow rates were determined for the five remedial wells and for the South Recharge Basins based on computer modeling (ARCADIS G&M, Inc. 2003c, modified in April 2005). Flow rates for Calpine, OXY Biosparge and West Recharge Basins are typical flow rates and are provided for reader information. "Design" flow volumes represent the volume of water that should be pumped/discharged during the reporting period and is calculated by multiplying the design rate by the reporting period duration.
- (3) "Average" flow rates for the remedial wells represent the average actual pumping rates when the pumps are operational and do not take into account the time that a well is not operational. During this reporting period, the remedial wells operated for the following percentage of the time: Well 1 (96%), Well 3R (99%), Well 17 (97%), Well 18 (98%), and Well 19 (94%). "Actual" volumes are determined via totalizing flow meters.
- (4) "Average" flow rates for the system discharges represent the average flow rate during the entire reporting period and are determined by dividing the total flow during the reporting period by the reporting period duration. The Calpine and South Recharge Basins flow volumes are determined via totalizing flow meters. The West Recharge Basin flow is calculated by subtracting the cumulative flow to the other discharges from the total influent flow. Actual flow to the recharge basins are greater than shown because storm water combines with the plant effluent prior to discharge to the recharge basins.
- (5) The TCE and TVOC concentrations for the remedial wells are from the quarterly sampling event performed during this reporting period on September 9, 2015 (Table 2).
- (6) The TVOC concentration for the two sets of recharge basins are their respective average monthly SPDES concentration for the current quarter.
- (7) TVOC mass removed for the reporting period is calculated by multiplying the TVOC concentration from the quarterly sampling event and the quantity of water pumped during the reporting period.
- (8) There are four discharges for the effluent groundwater: South Recharge Basins, West Recharge Basins, Calpine and OXY Biosparge system. Treated water is continuously discharged to the south and west recharge basins, and is available "on-demand" to both the Calpine Power Plant (Calpine) for use as make-up water, and the biosparge remediation system operated by Occidental Chemical (OXY Biosparge).
- (9) Treatment System Efficiencies are calculated by dividing the difference between the influent and effluent TVOC concentrations by the influent concentration.
- (10) The flow rate and volume for OXY Biosparge (Occidental Chemical) were estimated based on the average pumping rate calculated from data from April 2007 through March 2012.
- (11) Well 1 was shutdown from September 21 through 24, 2015 due to replacement of the existing drop pipe with stainless steel pipe.
- (12) A Well 3R pilot study was started on July 14, 2014 in an effort to increase the VOC mass removal through an increased pumping rate to approximately 1,000 gpm. The TVOC concentration and mass removal have increased since the initiation of the pilot study. On April 3, 2015 Well 3R flow rate was decreased to approximately 900 gpm to eliminate the increasing trend of TVOC concentration at Tower 96 effluent. Well 3R was brought online December 2013 to replace Well 3 due to decreasing specific capacity at Well 3 indicative of imminent well failure.
- (13) Well 1 and Well 3R were shutdown on August 26, 2015 due to cleaning of the distribution tray at the Tower 96 air stripper.
- (14) Well 17 was shut down due to pressure gauge issues and replacement, and installation of additional ball valve on the sampling port line.
- (15) Well 19 was shut down from August 21 through 25, 2015 due to flow meter calibration and repair.
- (16) Wells 17, 18 and 19 were shut down various times throughout the reporting period due to Tower 102 treatment system alarm testing, routine maintenance and well communication alarms.
- (17) Total pumpage/recharge rates are accurate to +/- 15% due to limitations in metering. Flow meter calibration is scheduled.
- (18) Storm Water Runoff Volume is calculated by multiplying the adjusted tributary area and NOAA precipitation data for the reporting period. The adjusted tributary area is tributary area that is adjusted by the runoff coefficient to exclude the infiltration volume from the total rainfall volume. The tributary area, runoff coefficient, and adjusted tributary area are from Dvirka and Bartilucci Consulting Engineers' Storm Water Permit Evaluation Report (January, 28, 2010). The NOAA precipitation data are calculated as a sum of NOAA daily precipitation data for the reporting period. NOAA precipitation data are retrieved from Station GHCND:USW00054787 - FARMINGDALE REPUBLIC AIRPORT, NY US.
- (19) Total Flow Volume to South Recharge Basins is estimated as a sum of flow volumes contributed from the Effluent Groundwater to South Recharge Basins and from Storm Water Runoff to South Recharge Basins.

**Abbreviations:**

--	Not Available or Not Applicable	TCE	Trichloroethene
TVOC	Total Volatile Organic Compounds	lbs	pounds
µg/L	micrograms per liter	MG	Million Gallons
VOC	Volatile Organic Compounds	NOAA	National Oceanic and Atmospheric Administration
gpm	gallons per minute	SPDES	State Pollutant Discharge Elimination System

**Table 2**  
**Water Sample Analytical Results for Groundwater**  
**Remedial Wells and Treatment Systems, Third Quarter 2015**  
**Northrop Grumman Systems Corporation, Operable Unit 2**  
**Bethpage, New York**

Constituent (Units in µg/L)	Location: Sample ID: Date:	WELL 1 WELL 1 9/9/2015	WELL 3R WELL 3R 9/9/2015	96 EFFLUENT T96 EFFLUENT (GW) 9/9/2015
<u>Volatile Organic Compounds (VOCs)<sup>(1)</sup></u>				
1,1,1-Trichloroethane		<5.0	<5.0	<1.0
1,1,2,2-Tetrachloroethane		<5.0	<5.0	<1.0
1,1,2-Trichloroethane		<5.0	<5.0	<1.0
1,1-Dichloroethane		<5.0	<b>1.3 J</b>	<1.0
1,1-Dichloroethene		<b>2.8 J</b>	<b>3.9 J</b>	<1.0
1,2-Dichloroethane		<5.0	<5.0	<1.0
1,2-Dichloropropane		<b>5.0</b>	<5.0	<1.0
2-Butanone (MEK)		<50	<50	<10
2-Hexanone (MBK)		<25	<25	<5.0
4-methyl-2-pentanone (MIK)		<25	<25	<5.0
Acetone		<50	<50	<10
Benzene		<2.5	<2.5	<0.50
Bromodichloromethane		<5.0	<5.0	<1.0
Bromoform		<5.0	<5.0	<1.0
Bromomethane		<10	<10	<2.0
Carbon Disulfide		<10	<10	<2.0
Carbon tetrachloride		<5.0	<5.0	<1.0
Chlorobenzene		<5.0	<5.0	<1.0
Chloroethane		<5.0	<5.0	<1.0
Chloroform		<5.0	<5.0	<1.0
Chloromethane		<5.0	<5.0	<1.0
cis-1,2-dichloroethene		<b>4.6 J</b>	<b>6.4</b>	<1.0
cis-1,3-dichloropropene		<5.0	<5.0	<1.0
Dibromochloromethane		<5.0	<5.0	<1.0
Ethylbenzene		<5.0	<5.0	<1.0
Methylene Chloride		<10	<10	<2.0
Styrene		<5.0	<5.0	<1.0
Tetrachloroethene		<b>35.4</b>	<b>36.6</b>	<1.0
Toluene		<5.0	<5.0	<1.0
trans-1,2-dichloroethene		<5.0	<5.0	<1.0
trans-1,3-dichloropropene		<5.0	<5.0	<1.0
Trichloroethylene		<b>852</b>	<b>557</b>	<b>0.66 J</b>
Trichlorotrifluoroethane (Freon 113)		<b>4.3 J</b>	<b>4.8 J</b>	<5.0
Vinyl Chloride		<5.0	<b>12.7</b>	<1.0
Xylene-o		<5.0	<5.0	<1.0
Xylenes-m,p		<5.0	<5.0	<1.0
<b>Total VOCs<sup>(2)</sup></b>		<b>900</b>	<b>620</b>	<b>0.66</b>
1,4-Dioxane <sup>(3)</sup>		<b>3.4</b>	<b>8.5</b>	<b>6.8</b>

See notes on last page.

**Table 2**  
**Water Sample Analytical Results for Groundwater**  
**Remedial Wells and Treatment Systems, Third Quarter 2015**  
**Northrop Grumman Systems Corporation, Operable Unit 2**  
**Bethpage, New York**

Constituent (Units in µg/L)	Location: Sample ID: Date:	WELL 17 WELL 17 9/9/2015	WELL 17-REP REP-090915-EE-1 9/9/2015	WELL 18 WELL 18 9/9/2015	WELL 19 WELL 19 9/9/2015	102 EFFLUENT T102 EFFLUENT (GW) 9/9/2015
<u>Volatile Organic Compounds (VOCs)<sup>(1)</sup></u>						
1,1,1-Trichloroethane		<b>0.51 J</b>	<b>0.54 J</b>	<b>0.87 J</b>	<b>0.43 J</b>	<1.0
1,1,2,2-Tetrachloroethane		<1.0	<1.0	<1.0	<1.0	<1.0 J
1,1,2-Trichloroethane		<1.0	<1.0	<1.0	<1.0	<1.0
1,1-Dichloroethane		<b>1.2</b>	<b>1.1</b>	<b>1.4</b>	<b>0.77 J</b>	<1.0
1,1-Dichloroethene		<b>1.9</b>	<b>2.0</b>	<1.0	<b>1.3</b>	<1.0
1,2-Dichloroethane		<1.0	<1.0	<1.0	<b>0.46 J</b>	<1.0
1,2-Dichloropropane		<1.0	<1.0	<1.0	<1.0	<1.0
2-Butanone (MEK)		<10	<10	<10	<10	<10 J
2-Hexanone (MBK)		<5.0	<5.0	<5.0	<5.0	<5.0 J
4-methyl-2-pentanone (MIK)		<5.0	<5.0	<5.0	<5.0	<5.0 J
Acetone		<10	<10	<10	<10	<10 J
Benzene		<0.50	<0.50	<0.50	<0.50	<0.50
Bromodichloromethane		<1.0	<1.0	<1.0	<1.0	<1.0
Bromoform		<1.0	<1.0	<1.0	<1.0	<1.0
Bromomethane		<2.0	<2.0	<2.0	<2.0	<2.0
Carbon Disulfide		<2.0	<2.0	<2.0	<2.0	<2.0
Carbon tetrachloride		<1.0	<1.0	<1.0	<1.0	<1.0
Chlorobenzene		<1.0	<1.0	<1.0	<1.0	<1.0
Chloroethane		<1.0	<1.0	<1.0	<1.0	<1.0
Chloroform		<b>0.43 J</b>	<b>0.47 J</b>	<1.0	<b>0.44 J</b>	<1.0
Chloromethane		<1.0	<1.0	<1.0	<1.0	<1.0
cis-1,2-dichloroethene		<b>3.8</b>	<b>3.7</b>	<b>2.0</b>	<b>18.3</b>	<1.0
cis-1,3-dichloropropene		<1.0	<1.0	<1.0	<1.0	<1.0
Dibromochloromethane		<1.0	<1.0	<1.0	<1.0	<1.0
Ethylbenzene		<1.0	<1.0	<1.0	<1.0	<1.0
Methylene Chloride		<2.0	<2.0	<2.0	<2.0	<2.0
Styrene		<1.0	<1.0	<1.0	<1.0	<1.0
Tetrachloroethene		<b>30</b>	<b>30.5</b>	<b>15.2</b>	<b>6.7</b>	<1.0
Toluene		<1.0	<1.0	<1.0	<1.0	<1.0
trans-1,2-dichloroethene		<1.0	<1.0	<1.0	<1.0	<1.0
trans-1,3-dichloropropene		<1.0	<1.0	<1.0	<1.0	<1.0
Trichloroethylene		<b>178</b>	<b>177</b>	<b>61.3</b>	<b>153</b>	<1.0
Trichlorotrifluoroethane (Freon 113)		<b>2.4 J</b>	<b>2.5 J</b>	<b>2.3 J</b>	<b>0.97 J</b>	<5.0
Vinyl Chloride		<1.0	<1.0	<1.0	<1.0	<1.0
Xylene-o		<1.0	<1.0	<1.0	<1.0	<1.0
Xylenes-m,p		<1.0	<1.0	<1.0	<1.0	<1.0
<b>Total VOCs<sup>(2)</sup></b>		<b>220</b>	<b>220</b>	<b>83</b>	<b>180</b>	0
1,4-Dioxane <sup>(3)</sup>		<b>2.8</b>	<b>3.0</b>	<b>2.2</b>	<b>1.9</b>	<b>2.8</b>

See notes on last page.



**Table 2**  
**Water Sample Analytical Results for Groundwater**  
**Remedial Wells and Treatment Systems, Third Quarter 2015**  
**Northrop Grumman Systems Corporation, Operable Unit 2**  
**Bethpage, New York**

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**Notes and Abbreviations:**

- (1) VOC samples analyzed using Method 8260C.
- (2) Total VOC results rounded to two significant figures.
- (3) 1,4-Dioxane samples analyzed using Method 8270D SIM.

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

<b>Bold</b>	Constituent detected
VOCs	Volatile Organic Compounds
µg/L	Micrograms per liter
J	Constituent value is estimated
REP	Field replicate
<5.0	Compound not detected above its laboratory quantification limit.
OU2	Operable Unit 2

**Table 3**  
**Concentrations of Site Related Volatile Organic Compounds**  
**and 1,4-Dioxane in Outpost Wells<sup>(1)</sup>**  
**Northrop Grumman Systems Corporation, Operable Unit 2**  
**Bethpage, New York**

Constituent (Units in µg/L)	Well: Sample ID: Date:	BPOW 1-1 BPOW 1-1 8/11/2015	BPOW 1-2 BPOW 1-2 8/11/2015	BPOW 1-3 BPOW 1-3 8/13/2015	BPOW 1-4 BPOW 1-4 8/18/2015	BPOW 1-5 BPOW 1-5 8/18/2015	BPOW 1-6 BPOW 1-6 8/17/2015	BPOW 2-1 BPOW 2-1 8/10/2015
<u>Volatile Organic Compounds (VOCs) <sup>(3)</sup></u>								
1,1,1-Trichloroethane		<b>0.67</b>	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50
1,1,2,2-Tetrachloroethane		<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50
1,1,2-Trichloroethane		<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50
1,1-Dichloroethane		<b>0.21J</b>	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50
1,1-Dichloroethene		<b>0.6</b>	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50
1,2-Dichloroethane		<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50
Carbon Tetrachloride		<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50
Chlorobenzene		<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50
Chloroform		<b>0.15J</b>	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50
cis-1,2-Dichloroethene		<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50
Trichlorotrifluoroethane (Freon 113)		<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0
Tetrachloroethene		<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50
trans-1,2-Dichloroethene		<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50
Trichloroethene		<b>1.1</b>	<b>0.30J</b>	<0.50	<0.50	<0.50	<0.50	<0.50
<b>Total Site-Related VOCs <sup>(4)</sup> :</b>		<b>2.7</b>	<b>0.3</b>	0	0	0	0	0
<hr/>								
1,4-Dioxane <sup>(5)</sup>		<0.20	<0.20	<0.21	<0.21	<0.20	<0.22	<b>0.09J</b>

See notes on last page.

**Table 3**  
**Concentrations of Site Related Volatile Organic Compounds**  
**and 1,4-Dioxane in Outpost Wells<sup>(1)</sup>**  
**Northrop Grumman Systems Corporation, Operable Unit 2**  
**Bethpage, New York**

Constituent (Units in µg/L)	Well: Sample ID: Date:	BPOW 2-2 BPOW 2-2 8/14/2015	BPOW 2-3 BPOW 2-3 8/14/2015	BPOW 3-1 BPOW 3-1 8/12/2015	BPOW 3-2 BPOW 3-2 9/17/2015	BPOW 3-3 BPOW 3-3 8/25/2015	BPOW 3-4 BPOW 3-4 8/25/2015	BPOW 3-4 REP082515BT1 8/25/2015
<u>Volatile Organic Compounds (VOCs) <sup>(3)</sup></u>								
1,1,1-Trichloroethane		<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50
1,1,2,2-Tetrachloroethane		<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50
1,1,2-Trichloroethane		<0.50	<0.50	<0.50	<0.50	<0.50	<b>0.79J</b>	<b>0.74J</b>
1,1-Dichloroethane		<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50
1,1-Dichloroethene		<0.50	<0.50	<0.50	<0.50	<0.50	<b>0.61</b>	<b>0.64</b>
1,2-Dichloroethane		<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50
Carbon Tetrachloride		<0.50	<0.50	<0.50	<0.50	<0.50	<b>1</b>	<b>1.1</b>
Chlorobenzene		<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50
Chloroform		<0.50	<0.50	<0.50	<0.50	<0.50	<b>1.3</b>	<b>1.4</b>
cis-1,2-Dichloroethene		<0.50	<0.50	<0.50	<0.50	<0.50	<b>1.1</b>	<b>1.1</b>
Trichlorotrifluoroethane (Freon 113)		<1.0	<1.0	<1.0	<1.0	<1.0	<b>0.75J</b>	<b>0.72J</b>
Tetrachloroethene		<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50
trans-1,2-Dichloroethene		<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50
Trichloroethene		<0.50	<0.50	<0.50	<0.50	<0.50	<b>60.9D</b>	<b>61D</b>
<b>Total Site-Related VOCs <sup>(4)</sup> :</b>		<b>0</b>	<b>0</b>	<b>0</b>	<b>0</b>	<b>0</b>	<b>66</b>	<b>67</b>
1,4-Dioxane <sup>(5)</sup>		<0.21	<0.21	<b>0.28</b>	<b>1.1</b>	<b>0.6</b>	<b>0.77J</b>	<b>0.32J</b>

See notes on last page.

**Table 3**  
**Concentrations of Site Related Volatile Organic Compounds**  
**and 1,4-Dioxane in Outpost Wells<sup>(1)</sup>**  
**Northrop Grumman Systems Corporation, Operable Unit 2**  
**Bethpage, New York**

Constituent (Units in µg/L)	Well: Sample ID: Date:	BPOW 4-1R <sup>(2)</sup> BPOW 4-1R 8/26/2015	BPOW 4-2R <sup>(2)</sup> BPOW 4-2R 8/26/2015
<u>Volatile Organic Compounds (VOCs) <sup>(3)</sup></u>			
1,1,1-Trichloroethane		<0.50	<b>0.063J</b>
1,1,2,2-Tetrachloroethane		<0.50	<0.50
1,1,2-Trichloroethane		<0.50	<0.50
1,1-Dichloroethane		<0.50	<0.50
1,1-Dichloroethene		<b>0.38J</b>	<b>0.52</b>
1,2-Dichloroethane		<b>0.13J</b>	<b>0.072J</b>
Carbon Tetrachloride		<b>0.21J</b>	<b>0.19J</b>
Chlorobenzene		<0.50	<0.50
Chloroform		<b>0.15J</b>	<b>0.090J</b>
cis-1,2-Dichloroethene		<b>0.086J</b>	<b>0.18J</b>
Trichlorotrifluoroethane (Freon 113)		<b>14.7</b>	<b>12.9</b>
Tetrachloroethene		<0.50	<0.50
trans-1,2-Dichloroethene		<0.50	<0.50
Trichloroethene		<b>0.92J</b>	<b>1.6</b>
<b>Total Site-Related VOCs <sup>(4)</sup> :</b>		<b>17</b>	<b>16</b>
<hr/>			
1,4-Dioxane <sup>(5)</sup>		<b>0.58</b>	<b>0.41</b>

See notes on last page.

**Table 3**  
**Concentrations of Site Related Volatile Organic Compounds**  
**and 1,4-Dioxane in Outpost Wells<sup>(1)</sup>**  
**Northrop Grumman Systems Corporation, Operable Unit 2**  
**Bethpage, New York**

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**Notes and Abbreviations:**

- (1) These outpost wells have been recently repurposed for use as plume monitoring wells per the June 2015 Groundwater Monitoring Plan Addendum (ARCADIS of New York, Inc., 2015) as conditionally approved by the NYSDEC (August 25, 2015). Therefore, TVOC trigger levels that may have been previously established are no longer shown.
- (2) The NAVY abandoned original Wells BPOW4-1 and BPOW4-2 and installed replacement Wells BPOW4-1R and BPOW4-2R between August, 2014 and October, 2014.
- (3) Samples analyzed for site related VOCs per the PWSCP (ARCADIS G&M, Inc. 2003) using USEPA Method 524.2
- (4) Site-related VOCs were established for the wells identified above in the Public Water Supply Contingency Plan (PWSCP) (ARCADIS G&M, Inc. 2003). Total Site-Related VOCs rounded to two significant figures.
- (5) Samples analyzed for 1,4-Dioxane by USEPA Method 8270D SIM.

**Bold** Bold value indicates constituent detected.  
 J Constituent value is estimated  
 D Concentration is based on a diluted sample analysis  
 REP Replicate Sample  
 TVOCs Total Volatile Organic Compounds  
 USEPA United States Environmental Protection Agency  
 VOC Volatile Organic Compounds  
 µg/L micrograms per liter  
 <0.5 Compound not detected above its laboratory quantification limit.

**Table 4**  
**Vapor Sample Analytical Results for Treatment Systems**  
**Third Quarter 2015**  
**Northrop Grumman Systems Corporation, Operable Unit 2**  
**Bethpage, New York**

Constituent□ (Units in µg/m3)	Sample ID: Date:	T96 Influent 9/9/2015	T96 Midtrain 9/9/2015	T96 Effluent 9/9/2015	T102 Influent 9/9/2015	T102 Effluent 9/9/2015
1,1,1-Trichloroethane		<65	<b>9.3</b>	<0.55	<b>10</b>	<0.55
1,1,2,2-Tetrachloroethane		<82	<5.5	<0.69	<5.5	<0.69
1,1,2-Trichloroethane		<65	<4.4	<0.55	<4.4	<0.55
1,1-Dichloroethane	49.0 J		<b>39</b>	<b>34</b>	<b>23</b>	<b>8.1</b>
1,1-Dichloroethylene		<b>148</b>	<b>142</b>	<b>60.7</b>	<b>52.7</b>	<b>35</b>
1,2-Dichloroethane	<97		<6.5	<0.81	<6.5	<0.81
1,2-Dichloropropane	84.1 J		<b>18</b>	<0.92	<7.4	<0.92
Benzene	<77		<5.1	<0.64	<5.1	<b>0.51</b>
Bromodichloromethane	<80		< 5.4	<0.67	< 5.4	<0.67
Bromoform	<49		< 3.3	<0.41	< 3.3	<0.41
Bromomethane	<93		< 6.2	<0.78	< 6.2	<0.78
Carbon disulfide	<75		< 5.0	<0.62	< 5.0	<0.62
Carbon tetrachloride	<30		<2.0	<0.25	<2.0	<0.25
Chlorobenzene	<110		<7.4	<0.92	<7.4	<0.92
Chloroethane	<63		<b>11</b>	<b>13</b>	<4.2	<0.53
Chloroform	<120		6.8 J	<b>2.5</b>	<b>8.8</b>	<b>1.6</b>
Chloromethane	<50		<3.3	<b>1.6</b>	<3.3	<b>1.3</b>
cis-1,3-Dichloropropene	<110		<7.3	<0.91	<7.3	<0.91
Dibromochloromethane	<100		<6.8	<0.85	<6.8	<0.85
Ethylbenzene	<100		<6.9	<0.87	<6.9	0.69 J
Methylene chloride	<b>154</b>		<5.6	<b>1.7</b>	<5.6	<b>1.5</b>
Styrene	<100		<6.8	<0.85	<6.8	<0.85
Tetrachloroethylene	<b>1,460</b>		<b>178</b>	<b>0.37</b>	<b>411</b>	<b>3.7</b>
Toluene	<90		<6.0	<0.75	< 6.0	<b>21</b>
trans-1,3-Dichloropropene	<110		<7.3	<0.91	<7.3	<0.91
Trichloroethylene	<b>29,900</b>		<b>8,010</b>	<b>3.8</b>	<b>3,710</b>	<b>12</b>
Trichlorotrifluoroethane (Freon 113)	<b>189</b>		<b>109</b>	<b>2.6</b>	<b>52</b>	<b>7.7</b>
Vinyl chloride	<b>268</b>		<b>284</b>	<b>28.4</b>	<0.82	<0.10
Xylene-o	<100		<6.9	<0.87	<6.9	0.56 J
Xylenes - m,p	<100		<6.9	<0.87	<6.9	<b>1.9</b>
<b>Total VOCs <sup>(2)</sup></b>		<b>32,104</b>	<b>8,617</b>	<b>54</b>	<b>4,182</b>	<b>96</b>

See notes on last page.

**Table 4**  
**Vapor Sample Analytical Results for Treatment Systems**  
**Third Quarter 2015**  
**Northrop Grumman Systems Corporation, Operable Unit 2**  
**Bethpage, New York**

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**Notes and Abbreviations:**

- (1) Vapor samples collected by ARCADIS on the dates shown and submitted to a NYSDOH ELAP certified laboratory for VOC analyses per Modified USEPA Method TO-15.
- (2) "Total VOCs" represents the sum of individual concentrations of compounds detected rounded to the nearest whole number.

**Acronyms\Key:**

**700** Bold data indicates that the analyte was detected at or above its reporting limit.

16 Data that is not bold indicates analyte detected but below its reporting limit; the value is estimated.

ELAP Environmental Laboratory Approval Program

NYSDOH New York State Department of Health

USEPA United States Environmental Protection Agency.

VOC Volatile organic compound.

$\mu\text{g}/\text{m}^3$  Micrograms per cubic meter.

**Table 5A**  
**Summary of SCREEN3 Model Input and Outputs**  
**Tower 96 Treatment System, Third Quarter 2015**  
**Northrop Grumman Systems Corporation, Operable Unit 2**  
**Bethpage, New York**

Parameters	Date Sampled: 11/13/2014	3/16/2015	5/11/2015	9/9/2015
<b>SCREEN3 Model Input</b>				
Source Type	Point	Point	Point	Point
Emission Rate (g/s)	1	1	1	1
Stack Height (ft)	55	55	55	55
Stack Height (m)	16.8	16.8	16.8	16.8
Stack Inside Diameter (m)	0.508	0.508	0.508	0.508
Air Flow Rate (scfm @stack temp) <sup>(1)</sup>	4,606	4,627	4,688	4,581
Air Flow Rate (acfm) <sup>(2), (3)</sup>	4,780	4,810	4,936	4,840
Stack Gas Exit Temperature (K) <sup>(2)</sup>	305	306	310	311
Ambient Air Temperature (K) <sup>(4)</sup>	286	277	287	293
Receptor Height (m) <sup>(5)</sup>	1.5	1.5	1.5	1.5
Urban/Rural	Urban	Urban	Urban	Urban
Building Height (m)	6.7	6.7	6.7	6.7
Min Horizontal Bldg Dim (m)	9.8	9.8	9.8	9.8
Max Horizontal Bldg Dim (m)	12.8	12.8	12.8	12.8
Consider Bldg Downwash?	Yes	Yes	Yes	Yes
Simple/Complex Terrain Above Stack	Simple	Simple	Simple	Simple
Simple/Complex Terrain Above Stack Base	Simple	Simple	Simple	Simple
Meteorology	Full	Full	Full	Full
Automated Distances Array	Yes	Yes	Yes	Yes
Terrain Height Above Stack Base	0	0	0	0
<b>SCREEN3 Model Output</b>				
1-HR Max Concentration at Receptor Height ( $\mu\text{g}/\text{m}^3$ ) <sup>(6)</sup>	198	197	195	199
Annualization Factor <sup>(7)</sup>	0.08	0.08	0.08	0.08
Average Annual Concentration at Receptor Height ( $\mu\text{g}/\text{m}^3$ ) <sup>(8)</sup>	15.8	15.8	15.6	15.9
Distance To Max Concentration (m) <sup>(9)</sup>	110	110	110	109

See notes on last page.



**Table 5A**  
**Summary of SCREEN3 Model Input and Outputs**  
**Tower 96 Treatment System, Third Quarter 2015**  
**Northrop Grumman Systems Corporation, Operable Unit 2**  
**Bethpage, New York**

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**Notes:**

- (1) The stack air flow rate at the stack temperature (in scfm) was calculated by multiplying the stack air flow rate in acfm by the ratio of the standard temperature to the actual stack gas exit temperature in degrees Kelvin.
- (2) The stack air flow rate (in acfm) and temperature were measured using inline instrumentation. Values were measured at the blower effluent location.
- (3) The stack air flow rate is taken from the actual stack air flow rate on the day of sampling.
- (4) The ambient temperature was recorded from weather.newsday.com website for Islip, New York. The mean actual temperature from the website was used in the model calculation
- (5) The receptor height corresponds to the average inhalation level.
- (6) SCREEN3 calculated constituent concentration at listed conditions at the specified inhalation level.
- (7) A USEPA time averaging conversion factor of 1/0.08 was used to convert the 1-hour maximum concentration output to an annual average.
- (8) Average annual constituent concentration at the receptor height was calculated by multiplying the one hour maximum concentration by the annualization factor.
- (9) SCREEN3 calculated distance to the 1-hour maximum concentration.

**Acronyms\Key:**

$\mu\text{g}/\text{m}^3$	Micrograms per cubic meter.
acfm	Actual cubic feet per minute.
ft	Feet.
g/s	Grams per second.
K	Kelvin.
m	Meters.
scfm	Standard cubic feet per minute.
USEPA	United States Environmental Protection Agency.

**Table 5B**  
**Summary of Maximum Allowable Stack Concentration Calculations**  
**Tower 96 Treatment System, Third Quarter 2015**  
**Northrop Grumman Systems Corporation, Operable Unit 2**  
**Bethpage, New York**

Compound	SGC <sup>(1)</sup> (µg/m <sup>3</sup> )	Actual Effluent Concentrations <sup>(2)</sup> (µg/m <sup>3</sup> )			
		11/13/2014	3/16/2015	5/11/2015	9/9/2015
1,1-Dichloroethane	95,000 <sup>(3)</sup>	0	8.9	0.57	34
1,1-Dichloroethene	188,000 <sup>(3)</sup>	3.7	138	2.3	60.7
Benzene	1,300	0.38	0	0	0
Chloroethane	619,000 <sup>(3)</sup>	19	20	12	13
Chloroform	150	0	0.63	0	2.5
Chloromethane	22,000	1.2	3.7	3.5	1.6
Methylene Chloride	14,000	18	1.5	1.7	1.7
Tetrachloroethylene	1,000	1.4	14	2.7	0.37
Toluene	37,000	0.49	0	0	0
Trichloroethylene	14,000	2	203	32	3.8
Trichlorotrifluoroethane (Freon 113)	960,000	0	4.5	0	2.6
Vinyl chloride	180,000	99.9	281	1.0	28.4

Compound	AGC <sup>(4)</sup> (µg/m <sup>3</sup> )	Annual MASC <sup>(5)</sup> (µg/m <sup>3</sup> )			
		11/13/2014	3/16/2015	5/11/2015	9/9/2015
1,1-Dichloroethane	0.63	1.77E+04	1.76E+04	1.73E+04	1.73E+04
1,1-Dichloroethene	200	5.61E+06	5.58E+06	5.50E+06	5.51E+06
Benzene	0.13	3.65E+03	3.62E+03	3.58E+03	3.58E+03
Chloroethane	10,000	2.81E+08	2.79E+08	2.75E+08	2.75E+08
Chloroform	14.700	4.12E+05	4.10E+05	4.05E+05	4.05E+05
Chloromethane	90	2.53E+06	2.51E+06	2.48E+06	2.48E+06
Methylene Chloride	60	1.68E+06	1.67E+06	1.65E+06	1.65E+06
Tetrachloroethylene	4.0	1.12E+05	1.12E+05	1.10E+05	1.10E+05
Toluene	5,000	1.40E+08	1.39E+08	1.38E+08	1.38E+08
Trichloroethylene	0.2	5.61E+03	5.58E+03	5.50E+03	5.51E+03
Trichlorotrifluoroethane (Freon 113)	180,000	5.05E+09	5.02E+09	4.95E+09	4.96E+09
Vinyl chloride	0.068	1.91E+03	1.90E+03	1.87E+03	1.87E+03

See notes on last page.

**Table 5B**  
**Summary of Maximum Allowable Stack Concentration Calculations**  
**Tower 96 Treatment System, Third Quarter 2015**  
**Northrop Grumman Systems Corporation, Operable Unit 2**  
**Bethpage, New York**

Compound	Percent of Annual MASC <sup>(6)</sup>				Cumulative % MASC <sup>(7)</sup>
	11/13/2014	3/16/2015	5/11/2015	9/9/2015	
1,1-Dichloroethane	0.0%	0.05%	0.0%	0.20%	0.08%
1,1-Dichloroethene	0.0%	0.0%	0.0%	0.0%	0.0%
Benzene	0.01%	0.0%	0.0%	0.0%	0.0%
Chloroethane	0.0%	0.0%	0.0%	0.0%	0.0%
Chloroform	0.0%	0.0%	0.0%	0.0%	0.0%
Chloromethane	0.0%	0.0%	0.0%	0.0%	0.0%
Methylene Chloride	0.0%	0.0%	0.0%	0.0%	0.0%
Tetrachloroethylene	0.0%	0.01%	0.0%	0.0%	0.0%
Toluene	0.0%	0.0%	0.0%	0.0%	0.0%
Trichloroethylene	0.04%	3.64%	0.58%	0.07%	1.35%
Trichlorotrifluoroethane (Freon 113)	0.0%	0.0%	0.0%	0.0%	0.0%
Vinyl chloride	5.24%	14.82%	0.05%	1.52%	6.45%

**Notes:**

- (1) Refers to the compound-specific SGC per the NYSDEC DAR-1 AGC/SGC tables revised February 28, 2014.
- (2) Only VOCs that were detected in the effluent vapor sample (T96 EFF) over the past year of system operation are included in this table.
- (3) An SGC was not provided in the DAR-1 AGC/SGC Tables, dated February 28, 2014. An interim SGC was developed based on guidelines provided in Section IV.A.2.b.1 of the NYSDEC DAR-1 Guidelines for the Control of Toxic Ambient Air Contaminants, 1991 edition.
- (4) AGC refers to the compound-specific annual guideline concentration per the NYSDEC DAR-1 AGC/SGC tables, revised February 28, 2014. NYSDEC DAR-1 AGCs were scaled using the results of a site-specific annual USEPA SCREEN 3 model to calculate the annual MASC per monitoring event.
- (5) Annual MASC was calculated by dividing the product of the AGC of a compound and the ratio of the SCREEN3 gas emission rate and the SCREEN 3 average annual concentration at receptor height by the air flow rate at the stack temperature and multiplying by the appropriate conversion factors.
- (6) Percent of MASC was calculated by dividing the actual effluent concentration by the MASC for the past four quarters of operation. Percentages have been rounded to two digits.
- (7) Cumulative percent of the MASC was calculated using a time-weighted average of the percent MASC per event. Percentages have been rounded to two digits.

**Acronyms\Key:**

µg/m <sup>3</sup>	Micrograms per cubic meter.
AGC	Annual guideline concentration.
SGC	Short-term guideline concentration
DAR-1	Division of Air Resources-1
MASC	Maximum allowable stack concentration.
NL	Compound concentration is not listed
NYSDEC	New York State Department of Environmental Conservation

**Table 6A**  
**Summary of SCREEN3 Model Input and Outputs**  
**Tower 102 Treatment System, Third Quarter 2015**  
**Northrop Grumman Systems Corporation, Operable Unit 2**  
**Bethpage, New York**

Parameters	Date Sampled:	11/13/14	3/16/15	5/11/15	9/9/15
<b>SCREEN3 Model Input</b>					
Source Type		Point	Point	Point	Point
Emission Rate (g/s)		1	1	1	1
Stack Height (ft)		69.52	69.52	69.52	69.52
Stack Height (m)		21.19	21.19	21.19	21.19
Stack Inside Diameter (m)		0.61	0.61	0.61	0.61
Air Flow Rate (scfm @ stack temp) <sup>(1)</sup>		8,123	8,231	8,068	7,930
Air Flow Rate (acfm) <sup>(2), (3)</sup>		8,320	8,420	8,220	8,080
Stack Gas Exit Temperature (K) <sup>(2)</sup>		301	301	300	300
Ambient Air Temperature (K) <sup>(4)</sup>		286	277	287	293
Receptor Height (m) <sup>(5)</sup>		1.5	1.5	1.5	1.5
Urban/Rural		Urban	Urban	Urban	Urban
Building Height (m)		7.62	7.62	7.62	7.62
Min Horizontal Bldg Dim (m)		12.5	12.5	12.5	12.5
Max Horizontal Bldg Dim (m)		15.54	15.54	15.54	15.54
Consider Bldg Downwash?		Yes	Yes	Yes	Yes
Simple/Complex Terrain Above Stack		Simple	Simple	Simple	Simple
Simple/Complex Terrain Above Stack Base		Simple	Simple	Simple	Simple
Meteorology		Full	Full	Full	Full
Automated Distances Array		Yes	Yes	Yes	Yes
Terrain Height Above Stack Base		0	0	0	0
<b>SCREEN3 Model Output</b>					
1-HR Max Concentration at Receptor Height ( $\mu\text{g}/\text{m}^3$ ) <sup>(6)</sup>		108.0	106.6	108.7	110.6
Annualization Factor <sup>(7)</sup>		0.08	0.08	0.08	0.08
Average Annual Concentration at Receptor Height ( $\mu\text{g}/\text{m}^3$ ) <sup>(8)</sup>		8.6	8.5	8.7	8.8
Distance To Max Concentration (m) <sup>(9)</sup>		146	146	145	144

See notes on last page.

**Table 6A**  
**Summary of SCREEN3 Model Input and Outputs**  
**Tower 102 Treatment System, Third Quarter 2015**  
**Northrop Grumman Systems Corporation, Operable Unit 2**  
**Bethpage, New York**

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**Notes:**

- (1) The stack air flow rate at the stack temperature (in scfm) was calculated by multiplying the stack air flow rate in acfm by the ratio of the standard temperature to the actual stack gas exit temperature in degrees Kelvin.
- (2) The stack air flow rate (in acfm) and temperature were measured using inline instrumentation. Values were measured at the blower effluent location.
- (3) The stack air flow rate is taken from the actual stack air flow rate on the day of sampling.
- (4) The ambient temperature was recorded from weather.newsday.com website for Islip, New York. The mean actual temperature from the website was used in the model calculation
- (5) The receptor height corresponds to the average inhalation level.
- (6) SCREEN3 calculated constituent concentration at listed conditions at the specified inhalation level.
- (7) A USEPA time averaging conversion factor of 1/0.08 was used to convert the 1-hour maximum concentration output to an annual average.
- (8) Average annual constituent concentration at the receptor height was calculated by multiplying the one hour maximum concentration by the annualization
- (9) SCREEN3 calculated distance to the 1-hour maximum concentration.

**Acronyms\Key:**

$\mu\text{g}/\text{m}^3$	Micrograms per cubic meter.
acfm	Actual cubic feet per minute.
ft	Feet.
g/s	Grams per second.
K	Kelvin.
m	Meters.
scfm	Standard cubic feet per minute.
USEPA	United States Environmental Protection Agency.

**Table 6B**  
**Summary of Maximum Allowable Stack Concentration Calculations**  
**Tower 102 Treatment System, Third Quarter 2015**  
**Northrop Grumman Systems Corporation, Operable Unit 2**  
**Bethpage, New York**

Compound	SGC <sup>(1)</sup> ( $\mu\text{g}/\text{m}^3$ )	Actual Effluent Concentrations <sup>(2)</sup> ( $\mu\text{g}/\text{m}^3$ )			
		11/13/2014	3/16/2015	5/11/2015	9/9/2015
1,1,1-Trichloroethane	9,000	0	0	0.55	0
1,1-Dichloroethane	95,000 <sup>(3)</sup>	4.5	3.2	5.7	8.1
1,1-Dichloroethene	188,000 <sup>(3)</sup>	15	11	21	35
Benzene	1,300	0	0	0	0.51
Carbon Disulfide	6,200	0	1.6	0	0
Chloroform	150	1.1	0.88	1.6	1.6
Chloromethane	22,000	1.1	1.2	0.99	1.3
Ethylbenzene	54,000	0	0	0	0.69
Methylene Chloride	14,000	13	16	1.3	1.5
Tetrachloroethene	1,000	7.5	5.2	1.6	3.7
Toluene	37,000	0	0	0	21
Trichloroethene	14,000	22	20	34	12
Trichlorotrifluoroethane (Freon 113)	960,000	4.4	3.1	7.1	7.7
Vinyl Chloride	180,000	0.41	0.31	0	0
Xylene-m,p	4,300	0	0	0	1.9
Xylene-o	4,300	0	0	0	0.56

See notes on last page.

**Table 6B**  
**Summary of Maximum Allowable Stack Concentration Calculations**  
**Tower 102 Treatment System, Third Quarter 2015**  
**Northrop Grumman Systems Corporation, Operable Unit 2**  
**Bethpage, New York**

Compound	AGC <sup>(4)</sup> (µg/m <sup>3</sup> )	Annual MASC <sup>(5)</sup> (µg/m <sup>3</sup> )			
		11/13/2014	3/16/2015	5/11/2015	9/9/2015
1,1,1-Trichloroethane	5,000	1.48E+08	1.48E+08	1.48E+08	1.49E+08
1,1-Dichloroethane	0.63	1.87E+04	1.87E+04	1.87E+04	1.88E+04
1,1-Dichloroethene	200	5.92E+06	5.92E+06	5.93E+06	5.96E+06
Benzene	0.13	3.85E+03	3.85E+03	3.85E+03	3.87E+03
Carbon Disulfide	700	2.07E+07	2.07E+07	2.07E+07	2.09E+07
Chloroform	14.700	4.35E+05	4.35E+05	4.36E+05	4.38E+05
Chloromethane	90	2.67E+06	2.66E+06	2.67E+06	2.68E+06
Ethylbenzene	1,000	2.96E+07	2.96E+07	2.96E+07	2.98E+07
Methylene Chloride	60.0	1.78E+06	1.78E+06	1.78E+06	1.79E+06
Tetrachloroethene	4	1.18E+05	1.18E+05	1.19E+05	1.19E+05
Toluene	5,000	1.48E+08	1.48E+08	1.48E+08	1.49E+08
Trichloroethene	0.2	5.92E+03	5.92E+03	5.93E+03	5.96E+03
Trichlorotrifluoroethane (Freon 113)	180,000	5.33E+09	5.33E+09	5.33E+09	5.36E+09
Vinyl Chloride	0.068	2.01E+03	2.01E+03	2.01E+03	2.03E+03
Xylene-m,p	100	2.96E+06	2.96E+06	2.96E+06	2.98E+06
Xylene-o	100	2.96E+06	2.96E+06	2.96E+06	2.98E+06

See notes on last page.

**Table 6B**  
**Summary of Maximum Allowable Stack Concentration Calculations**  
**Tower 102 Treatment System, Third Quarter 2015**  
**Northrop Grumman Systems Corporation, Operable Unit 2**  
**Bethpage, New York**

Compound	Percent of Annual MASC <sup>(6)</sup>				Cumulative % MASC <sup>(7)</sup>
	11/13/14	03/16/15	5/11/2015	9/9/2015	
1,1,1-Trichloroethane	0.0%	0.0%	0.0%	0.0%	0.0%
1,1-Dichloroethane	0.02%	0.02%	0.03%	0.04%	0.03%
1,1-Dichloroethene	0.0%	0.0%	0.0%	0.0%	0.0%
Benzene	0.0%	0.0%	0.0%	0.01%	0.0%
Carbon Disulfide	0.0%	0.0%	0.0%	0.0%	0.0%
Chloroform	0.0%	0.0%	0.0%	0.0%	0.0%
Chloromethane	0.0%	0.0%	0.0%	0.0%	0.0%
Ethylbenzene	0.0%	0.0%	0.0%	0.0%	0.0%
Methylene Chloride	0.0%	0.0%	0.0%	0.0%	0.0%
Tetrachloroethene	0.01%	0.0%	0.0%	0.0%	0.0%
Toluene	0.0%	0.0%	0.0%	0.0%	0.0%
Trichloroethene	0.37%	0.34%	0.57%	0.20%	0.33%
Trichlorotrifluoroethane (Freon 113)	0.0%	0.0%	0.0%	0.0%	0.0%
Vinyl Chloride	0.02%	0.02%	0.0%	0.0%	0.01%
Xylene-m,p	0.0%	0.0%	0.0%	0.0%	0.0%
Xylene-o	0.0%	0.0%	0.0%	0.0%	0.0%

See notes on last page.



**Table 6B**  
**Summary of Maximum Allowable Stack Concentration Calculations**  
**Tower 102 Treatment System, Third Quarter 2015**  
**Northrop Grumman Systems Corporation, Operable Unit 2**  
**Bethpage, New York**

**Notes:**

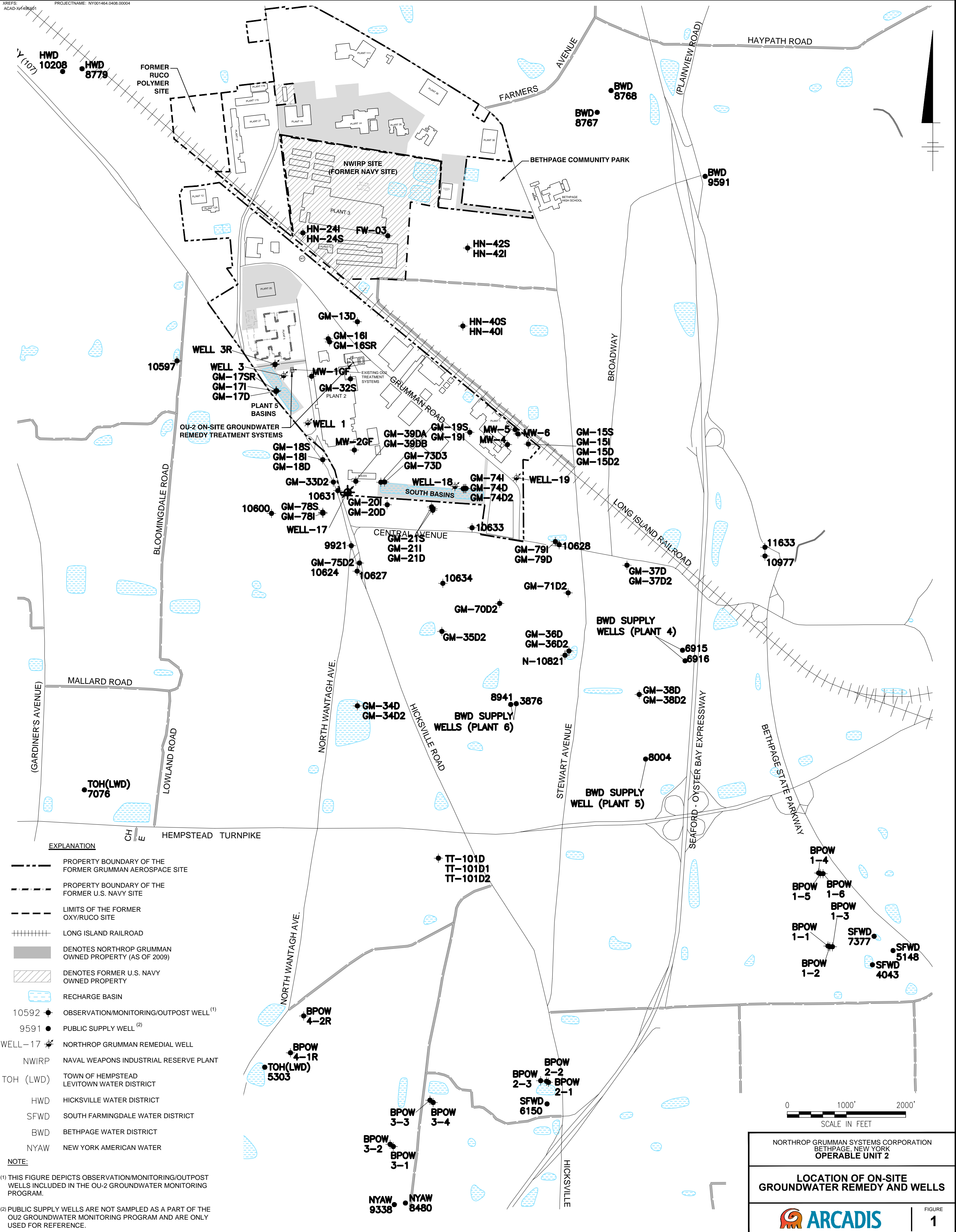
- (1) Refers to the compound-specific SGC per the NYSDEC DAR-1 AGC/SGC tables revised February 28, 2014.
- (2) Only VOCs that were detected in the effluent vapor sample (T102 EFF) over the past year of system operation are included in this table.
- (3) An SGC was not provided in the DAR-1 AGC/SGC Tables, dated February 28, 2014. An interim SGC was developed based on guidelines provided in Section IV.A.2.b.1 of the NYSDEC DAR-1 Guidelines for the Control of Toxic Ambient Air Contaminants, 1991 edition.
- (4) AGC refers to the compound-specific annual guideline concentration per the NYSDEC DAR-1 AGC/SGC tables, revised February 28, 2014. NYSDEC DAR-1 AGCs were scaled using the results of a site-specific annual USEPA SCREEN 3 model to calculate the annual MASC per monitoring event.
- (5) Annual MASC was calculated by dividing the product of the AGC of a compound and the ratio of the SCREEN3 gas emission rate and the SCREEN 3 average annual concentration at receptor height by the air flow rate at the stack temperature and multiplying by the appropriate conversion factors.
- (6) Percent of MASC was calculated by dividing the actual effluent concentration by the MASC for the past four quarters of operation. Percentages have been rounded to two digits.
- (7) Cumulative percent of the MASC was calculated using a time-weighted average of the percent MASC per event. Percentages have been rounded to two digits.

**Acronyms\Key:**

--	Compound not reported, unable to compute MASC
AGC	Annual guideline concentration.
NL	Compound concentration not listed
DAR-1	Division of Air Resources-1
MASC	Maximum allowable stack concentration.
$\mu\text{g}/\text{m}^3$	Micrograms per cubic meter.
NYSDEC	New York State Department of Environmental Conservation

# FIGURES





**EXPLANATION**

- PROPERTY BOUNDARY OF THE FORMER GRUMMAN AEROSPACE SITE
- PROPERTY BOUNDARY OF THE FORMER U.S. NAVY SITE
- LIMITS OF THE FORMER OXY/RUCO SITE
- +++++ LONG ISLAND RAILROAD
- DENOTES NORTHROP GRUMMAN OWNED PROPERTY (AS OF 2009)
- ▨ DENOTES FORMER U.S. NAVY OWNED PROPERTY
- RECHARGE BASIN
- 10592 ● OBSERVATION/MONITORING/OUTPOST WELL (1)
- 9591 ● PUBLIC SUPPLY WELL (2)
- WELL-17 ● NORTHROP GRUMMAN REMEDIAL WELL
- NWIRP NAVAL WEAPONS INDUSTRIAL RESERVE PLANT
- TOH (LWD) TOWN OF HEMPSTEAD LEVITOWN WATER DISTRICT
- HWD HICKSVILLE WATER DISTRICT
- SFWD SOUTH FARMINGDALE WATER DISTRICT
- BWD BETHPAGE WATER DISTRICT
- NYAW NEW YORK AMERICAN WATER

**NOTE:**  
 (1) THIS FIGURE DEPICTS OBSERVATION/MONITORING/OUTPOST WELLS INCLUDED IN THE OU-2 GROUNDWATER MONITORING PROGRAM.  
 (2) PUBLIC SUPPLY WELLS ARE NOT SAMPLED AS A PART OF THE OU2 GROUNDWATER MONITORING PROGRAM AND ARE ONLY USED FOR REFERENCE.

0 1000' 2000'  
SCALE IN FEET

NORTHROP GRUMMAN SYSTEMS CORPORATION  
BETHPAGE, NEW YORK  
OPERABLE UNIT 2

**LOCATION OF ON-SITE  
GROUNDWATER REMEDY AND WELLS**

**ARCADIS**

FIGURE  
**1**