

DECEMBER 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

March 2016

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Norfolk, Virginia 23511

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

March 2016

A handwritten signature in black ink that reads "Brian Caldwell".

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Brian Caldwell  
Contract Task Order Manager

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List of Acronyms and Abbreviations

DOT	Department of Transportation
IDW	Investigation Derived Waste
Katahdin	Katahdin Analytical Services
NG	Northrop Grumman
NWIRP	Naval Weapons Industrial Reserve Plant
NYSDEC	New York State Department of Environmental Conservation
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

## 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 December 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 47 Navy-owned monitoring wells by Resolution Consultants on behalf of the Navy, and by ARCADIS on behalf of the Navy at the direction of Northrop Grumman (NG) as part of an agreement between the Navy and NG. 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, to evaluate the southernmost extend of the OU2 plume, and to evaluate outpost wells intended to provide early warning of plume migration to public water supply wells. The locations of monitoring wells sampled as part of this effort are shown in Figure 2. Well construction information and sampling responsibility are listed in Table 1.

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## 2.0 FIELD PROGRAM

Field tasks were conducted in December 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 December 2015 quarterly sampling round consisted of a total of 47 wells (Table 1). Of these, 30 groundwater wells were sampled by Resolution Consultants and 17 were sampled by ARCADIS, NG consultant. ARCADIS sampled the following wells after initial sampling by Resolution Consultants in September 2015 (Resolution Consultants, 2016): RE117D1, RE117D2, RE118D1, RE119D1 and BPOW5-7. These wells were transitioned to ARCADIS for the December 2015 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 for ARCADIS-sampled wells are provided in Table 4, Table 5, Table 6; data validation packages are included in Appendix C.

Additional Navy-owned wells were sampled by ARCADIS in the fourth quarter of 2015 as part of separate and ongoing OU2 monitoring programs, as summarized in the sampling schedule in Appendix D. ARCADIS will document these activities and results in their 2015 Annual Groundwater

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Monitoring Report, scheduled for submission to New York State Department of Environmental Conservation in the spring of 2016.

## 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 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, 5 and 6. Data validation packages for wells sampled by ARCADIS are included in Appendix C.

Additional Navy-owned wells sampled by ARCADIS in the fourth quarter of 2015 as part of separate and ongoing OU2 monitoring programs are summarized in Appendix D.

#### 4.0 REFERENCES

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

Resolution Consultants, 2016. September *2015 Groundwater Sampling Data Summary Report, Bethpage, NY*. February.

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
RE107D1	530	505	525	515	5	VPB141	Resolution
RE107D2	585	560	580	570	5	VPB141	Resolution
RE107D3	670	645	665	655	5	VPB141	Resolution
RE108D1	555	530	550	540	5	VPB142	Resolution
RE108D2	655	630	650	640	5	VPB142	Resolution
RE114D1	560	535	555	545	5	VPB148	Resolution
RE114D2	635	610	630	620	5	VPB148	Resolution
RE114D3	725	700	720	710	5	VPB148	Resolution
RE117D1	760	730	755	742.5	5	VPB151	ARCADIS
RE117D2	810	780	805	792.5	5	VPB151	ARCADIS
RE118D1	795	765	790	777.5	5	VPB152	ARCADIS
RE119D1	745	715	740	727.5	5	VPB153	ARCADIS
RE120D1	655	630	650	640	5	VPB154	Resolution
RE120D2	713	690	710	700	3	VPB154	Resolution
RE120D3	765	740	760	750	5	VPB154	Resolution
RE121D1	575	550	570	560	5	VPB155	Resolution
RE121D2	755	730	750	740	5	VPB155	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



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
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
BPOW5-7	555	525	550	537.5	5	VPB153	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)	RE103D1	RE103D2	RE103D3	RE122D1
		12/14/2015	12/14/2015	12/14/2015	12/15/2015
Sample Date		RE103D1-GW-121415	RE103D2-GW-121415	RE103D3-GW-121415	RE122D1-GW-121515
Sample ID					
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	12	3.2	2.5	4.7
1,1,2-TRICHLOROETHANE	1	0.62 J	< 0.50 U	< 0.50 U	< 0.50 U
1,1-DICHLOROETHANE	5	1.1	< 0.50 U	< 0.50 U	< 0.50 U
1,1-DICHLOROETHENE	5	7.6	0.77 J	0.62 J	0.63 J
1,2,4-TRICHLOROENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DIBROMO-3-CHLOROPROPANE	0.04	< 0.75 U	< 0.75 U	< 0.75 U	< 0.75 U
1,2-DIBROMOETHANE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DICHLOROBENZENE	3	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DICHLOROETHANE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DICHLOROETHENE, TOTAL	5	3.2	1.1 J	1.0 J	1.9 J
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	12	1.2	0.81	8.7
2-BUTANONE	50	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 U
2-HEXANONE	50	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 U
4-METHYL-2-PENTANONE	NL	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 U
ACETONE	50	< 2.5 U	< 2.5 U	< 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 UJ	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ
CARBON DISULFIDE	60	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CARBON TETRACHLORIDE	5	< 0.50 U	< 0.50 U	0.24 J	< 0.50 U
CHLOROBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
CHLOROFORM	7	< 0.50 U	< 0.50 U	0.79 J	< 0.50 U
CHLOROMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
CIS-1,2-DICHLOROETHENE	5	3.2	1.1	1.0	1.9
CIS-1,3-DICHLOROPROPENE	0.4	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CYCLOHEXANE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
DIBROMOCHLOROMETHANE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
DICHLORODIFLUOROMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
ETHYLBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
ISOPROPYLBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
M- AND P-XYLENE	NL	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
METHYL ACETATE	NL	< 0.75 U	< 0.75 U	< 0.75 U	< 0.75 U
METHYL CYCLOHEXANE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
METHYL TERT-BUTYL ETHER	10	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
METHYLENE CHLORIDE	5	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 U
O-XYLENE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
STYRENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TETRACHLOROETHENE	5	3.0 J	0.72 J	< 0.50 UJ	1.5 J
TOLUENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TRANS-1,2-DICHLOROETHENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TRANS-1,3-DICHLOROPROPENE	0.4	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TRICHLOROETHENE	5	930	620	510	600
TRICHLOROFLUOROMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
VINYL CHLORIDE	2	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
XYLENES, TOTAL	5	< 1.5 U	< 1.5 U	< 1.5 U	< 1.5 U

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

Location	NYSDEC Groundwater Guidance or Standard Value (Note 1)	RE122D2	RE122D3	RE104D1	RE104D2
		12/15/2015	12/15/2015	12/15/2015	12/15/2015
Sample Date		RE122D2-GW-121515	RE122D3-GW-121515	RE104D1-GW-121515	RE104D2-GW-121515
Sample ID					
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	21	< 0.50 U	4.6	< 0.50 U
1,1,2-TRICHLOROETHANE	1	3.1	< 0.50 U	< 0.50 U	< 0.50 U
1,1-DICHLOROETHANE	5	1.5	< 0.50 U	< 0.50 U	< 0.50 U
1,1-DICHLOROETHENE	5	8.9	< 0.50 U	0.80 J	< 0.50 U
1,2,4-TRICHLOROENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DIBROMO-3-CHLOROPROPANE	0.04	< 0.75 U	< 0.75 U	< 0.75 U	< 0.75 U
1,2-DIBROMOETHANE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DICHLOROBENZENE	3	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DICHLOROETHANE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DICHLOROETHENE, TOTAL	5	5.7	< 1.0 U	1.1 J	2.7
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	11	< 0.17 U	6.9	0.22 J
2-BUTANONE	50	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 U
2-HEXANONE	50	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 U
4-METHYL-2-PENTANONE	NL	< 2.5 U	< 2.5 UJ	< 2.5 U	< 2.5 U
ACETONE	50	< 2.5 U	< 2.5 U	< 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 UJ	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ
CARBON DISULFIDE	60	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CARBON TETRACHLORIDE	5	1.9	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
CHLOROFORM	7	2.6	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
CIS-1,2-DICHLOROETHENE	5	5.7	< 0.50 U	1.1	2.7
CIS-1,3-DICHLOROPROPENE	0.4	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CYCLOHEXANE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
DIBROMOCHLOROMETHANE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
DICHLORODIFLUOROMETHANE	5	< 1.0 U	< 1.0 U	0.68 J	< 1.0 U
ETHYLBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
ISOPROPYLBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
M- AND P-XYLENE	NL	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
METHYL ACETATE	NL	< 0.75 U	< 0.75 U	< 0.75 U	< 0.75 U
METHYL CYCLOHEXANE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
METHYL TERT-BUTYL ETHER	10	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
METHYLENE CHLORIDE	5	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 U
O-XYLENE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
STYRENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TETRACHLOROETHENE	5	2.3 J	< 0.50 UJ	1.9 J	< 0.50 UJ
TOLUENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TRANS-1,2-DICHLOROETHENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TRANS-1,3-DICHLOROPROPENE	0.4	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TRICHLOROETHENE	5	4700	2.5	110	6.8
TRICHLOROFLUOROMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
VINYL CHLORIDE	2	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
XYLENES, TOTAL	5	< 1.5 U	< 1.5 U	< 1.5 U	< 1.5 U

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

Location	NYSDEC Groundwater Guidance or Standard Value (Note 1)	RE104D2	RE104D3	RE114D1	RE114D2
Sample Date		12/15/2015	12/15/2015	12/21/2015	12/16/2015
Sample ID		DUPLICATE1-GW-121515	RE104D3-GW-121515	RE114D1-GW-122115	RE114D2-GW-121615
Sample type code		FD	N	N	N
<b>VOC 8260C (ug/L)</b>					
1,1,1-TRICHLOROETHANE	5	< 0.50 U	< 0.50 U	0.64 J	< 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	20 J	14
1,1,2-TRICHLOROETHANE	1	< 0.50 U	< 0.50 U	1.6 J	< 0.50 U
1,1-DICHLOROETHANE	5	< 0.50 U	< 0.50 U	1.5 J	< 0.50 U
1,1-DICHLOROETHENE	5	< 0.50 U	< 0.50 U	4.0 J	< 0.50 U
1,2,4-TRICHLOROENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DIBROMO-3-CHLOROPROPANE	0.04	< 0.75 U	< 0.75 U	< 0.75 U	< 0.75 U
1,2-DIBROMOETHANE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DICHLOROBENZENE	3	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DICHLOROETHANE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DICHLOROETHENE, TOTAL	5	2.7	< 1.0 U	5.1 J	0.82 J
1,2-DICHLOROPROPANE	1	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,3-DICHLOROBENZENE	3	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,4-DICHLOROBENZENE	3	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,4-DIOXANE (Method 8270D_SIM)	NL	0.28	< 0.17 U	5.5	2.5
2-BUTANONE	50	< 2.5 U	< 2.5 U	< 2.5 UJ	< 2.5 U
2-HEXANONE	50	< 2.5 U	< 2.5 U	< 2.5 UJ	< 2.5 U
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 UJ	< 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 UJ	< 1.0 UJ	< 1.0 U	< 1.0 UJ
CARBON DISULFIDE	60	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CARBON TETRACHLORIDE	5	< 0.50 U	< 0.50 U	2.5 J	< 0.50 U
CHLOROBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROETHANE	5	< 1.0 U	< 1.0 U	< 1.0 UJ	< 1.0 U
CHLOROFORM	7	< 0.50 U	< 0.50 U	2.9 J	0.40 J
CHLOROMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
CIS-1,2-DICHLOROETHENE	5	2.7	< 0.50 U	5.1 J	0.82 J
CIS-1,3-DICHLOROPROPENE	0.4	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CYCLOHEXANE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
DIBROMOCHLOROMETHANE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
DICHLORODIFLUOROMETHANE	5	< 1.0 U	< 1.0 U	1.0 J	< 1.0 U
ETHYLBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
ISOPROPYLBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
M- AND P-XYLENE	NL	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
METHYL ACETATE	NL	< 0.75 U	< 0.75 U	< 0.75 U	< 0.75 U
METHYL CYCLOHEXANE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
METHYL TERT-BUTYL ETHER	10	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
METHYLENE CHLORIDE	5	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 U
O-XYLENE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
STYRENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TETRACHLOROETHENE	5	< 0.50 UJ	< 0.50 UJ	< 0.50 U	< 0.50 UJ
TOLUENE	5	< 0.50 U	< 0.50 U	0.30 J	< 0.50 U
TRANS-1,2-DICHLOROETHENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TRANS-1,3-DICHLOROPROPENE	0.4	< 0.50 U	< 0.50 U	< 0.50 UJ	< 0.50 U
TRICHLOROETHENE	5	6.8	< 0.50 U	370	70
TRICHLOROFUOROMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
VINYL CHLORIDE	2	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
XYLENES, TOTAL	5	< 1.5 U	< 1.5 U	< 1.5 U	< 1.5 U

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

Location	NYSDEC Groundwater Guidance or Standard Value (Note 1)	RE114D3	RE105D1	RE105D2	TT101D
		12/16/2015	12/17/2015	12/17/2015	12/17/2015
Sample Date		RE114D3-GW-121615	RE105D1-GW-121715	RE105D2-GW-121715	TT101D-GW-121715
Sample ID					
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	13	8.7	26	16
1,1,2-TRICHLOROETHANE	1	< 0.50 U	< 0.50 U	1.3	< 0.50 U
1,1-DICHLOROETHANE	5	< 0.50 U	< 0.50 U	1.9	0.84 J
1,1-DICHLOROETHENE	5	1.1	1.3	7.0	3.4
1,2,4-TRICHLOROENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DIBROMO-3-CHLOROPROPANE	0.04	< 0.75 U	< 0.75 U	< 0.75 U	< 0.75 U
1,2-DIBROMOETHANE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DICHLOROBENZENE	3	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DICHLOROETHANE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DICHLOROETHENE, TOTAL	5	0.67 J	1.7 J	4.0	3.1
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	2.1	10	5.8	8.4
2-BUTANONE	50	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 U
2-HEXANONE	50	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 U
4-METHYL-2-PENTANONE	NL	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ
ACETONE	50	< 2.5 U	< 2.5 U	< 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 UJ	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ
CARBON DISULFIDE	60	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CARBON TETRACHLORIDE	5	< 0.50 U	< 0.50 U	3.0	< 0.50 U
CHLOROBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
CHLOROFORM	7	< 0.50 U	0.38 J	2.0	0.55 J
CHLOROMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
CIS-1,2-DICHLOROETHENE	5	0.67 J	1.7	4.0	3.1
CIS-1,3-DICHLOROPROPENE	0.4	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CYCLOHEXANE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
DIBROMOCHLOROMETHANE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
DICHLORODIFLUOROMETHANE	5	< 1.0 U	0.58 J	0.45 J	2.2
ETHYLBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
ISOPROPYLBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
M- AND P-XYLENE	NL	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
METHYL ACETATE	NL	< 0.75 U	< 0.75 U	< 0.75 U	< 0.75 U
METHYL CYCLOHEXANE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
METHYL TERT-BUTYL ETHER	10	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
METHYLENE CHLORIDE	5	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 U
O-XYLENE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
STYRENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TETRACHLOROETHENE	5	< 0.50 UJ	< 0.50 UJ	1.9 J	< 0.50 UJ
TOLUENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TRANS-1,2-DICHLOROETHENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TRANS-1,3-DICHLOROPROPENE	0.4	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TRICHLOROETHENE	5	43	120	1800	74
TRICHLOROFUOROMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
VINYL CHLORIDE	2	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
XYLENES, TOTAL	5	< 1.5 U	< 1.5 U	< 1.5 U	< 1.5 U

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

Location	NYSDEC Groundwater Guidance or Standard Value (Note 1)	TT101D1	TT101D2	TT101D2	RE107D1
Sample Date		12/17/2015	12/21/2015	12/21/2015	12/18/2015
Sample ID		TT101D1-GW-121715	TT101D2-GW-122115	DUPLICATE-GW-122115	RE107D1-GW-121815
Sample type code		N	N	FD	N
<b>VOC 8260C (ug/L)</b>					
1,1,1-TRICHLOROETHANE	5	< 0.50 U	0.34 J	< 0.50 U	< 0.50 U
1,1,2,2-TETRACHLOROETHANE	5	< 0.50 U	< 0.50 UJ	< 0.50 U	< 0.50 U
1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	5	16	19	24	0.95 J
1,1,2-TRICHLOROETHANE	1	0.48 J	0.50 J	0.65 J	< 0.50 U
1,1-DICHLOROETHANE	5	< 0.50 U	< 0.50 UJ	0.81 J	< 0.50 U
1,1-DICHLOROETHENE	5	4.6	3.6 J	5.0 J	< 0.50 U
1,2,4-TRICHLOROENZENE	5	< 0.50 U	< 0.50 UJ	< 0.50 U	< 0.50 U
1,2-DIBROMO-3-CHLOROPROPANE	0.04	< 0.75 U	< 0.75 UJ	< 0.75 U	< 0.75 U
1,2-DIBROMOETHANE	NL	< 0.50 U	< 0.50 UJ	< 0.50 U	< 0.50 U
1,2-DICHLOROBENZENE	3	< 0.50 U	< 0.50 UJ	< 0.50 U	< 0.50 U
1,2-DICHLOROETHANE	5	< 0.50 U	< 0.50 UJ	< 0.50 U	< 0.50 U
1,2-DICHLOROETHENE, TOTAL	5	1.9 J	1.7 J	2.0	0.21 J
1,2-DICHLOROPROPANE	1	< 0.50 U	< 0.50 UJ	< 0.50 U	< 0.50 U
1,3-DICHLOROBENZENE	3	< 0.50 U	< 0.50 UJ	< 0.50 U	< 0.50 U
1,4-DICHLOROBENZENE	3	< 0.50 U	< 0.50 UJ	< 0.50 U	< 0.50 U
1,4-DIOXANE (Method 8270D_SIM)	NL	11	1.7	2.2	6.9
2-BUTANONE	50	< 2.5 U	< 2.5 U	< 2.5 UJ	< 2.5 U
2-HEXANONE	50	< 2.5 U	< 2.5 U	< 2.5 UJ	< 2.5 U
4-METHYL-2-PENTANONE	NL	< 2.5 UJ	< 2.5 UJ	< 2.5 U	< 2.5 U
ACETONE	50	< 2.5 U	< 2.5 U	< 2.5 UJ	< 2.5 UJ
BENZENE	1	< 0.50 U	< 0.50 UJ	< 0.50 U	< 0.50 U
BROMODICHLOROMETHANE	50	< 0.50 U	< 0.50 UJ	< 0.50 U	< 0.50 U
BROMOFORM	50	< 0.50 U	< 0.50 UJ	< 0.50 U	< 0.50 U
BROMOMETHANE	5	< 1.0 UJ	< 1.0 UJ	< 1.0 U	< 1.0 U
CARBON DISULFIDE	60	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CARBON TETRACHLORIDE	5	1.9	1.3 J	1.4	< 0.50 U
CHLOROBENZENE	5	< 0.50 U	< 0.50 UJ	< 0.50 U	< 0.50 U
CHLOROETHANE	5	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ	< 1.0 UJ
CHLOROFORM	7	1.0	0.90 J	0.92 J	< 0.50 U
CHLOROMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 UJ
CIS-1,2-DICHLOROETHENE	5	1.9	1.7 J	2.0	0.21 J
CIS-1,3-DICHLOROPROPENE	0.4	< 0.50 U	< 0.50 UJ	< 0.50 U	< 0.50 U
CYCLOHEXANE	NL	< 0.50 U	< 0.50 UJ	< 0.50 U	< 0.50 U
DIBROMOCHLOROMETHANE	5	< 0.50 U	< 0.50 UJ	< 0.50 U	< 0.50 U
DICHLORODIFLUOROMETHANE	5	1.8 J	< 1.0 U	< 1.0 U	< 1.0 U
ETHYLBENZENE	5	< 0.50 U	< 0.50 UJ	< 0.50 U	< 0.50 U
ISOPROPYLBENZENE	5	< 0.50 U	< 0.50 UJ	< 0.50 U	< 0.50 U
M- AND P-XYLENE	NL	< 1.0 U	< 1.0 UJ	< 1.0 U	< 1.0 U
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 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 UJ	< 2.5 U	< 2.5 U
O-XYLENE	NL	< 0.50 U	< 0.50 UJ	< 0.50 U	< 0.50 U
STYRENE	5	< 0.50 U	< 0.50 UJ	< 0.50 U	< 0.50 U
TETRACHLOROETHENE	5	< 0.50 UJ	< 0.50 UJ	0.94 J	1.6 J
TOLUENE	5	< 0.50 U	< 0.50 UJ	< 0.50 U	< 0.50 U
TRANS-1,2-DICHLOROETHENE	5	< 0.50 U	< 0.50 UJ	< 0.50 U	< 0.50 U
TRANS-1,3-DICHLOROPROPENE	0.4	< 0.50 U	< 0.50 UJ	< 0.50 UJ	< 0.50 UJ
TRICHLOROETHENE	5	200	510	590	17
TRICHLOROFUOROMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
VINYL CHLORIDE	2	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
XYLENES, TOTAL	5	< 1.5 U	< 1.5 UJ	< 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)	RE107D2	RE107D3	RE123D1	RE123D2
Sample Date		12/18/2015	12/29/2015	12/21/2015	12/21/2015
Sample ID		RE107D2-GW-121815	RE107D3-GW-122915	RE123D1-GW-122115	RE123D2-GW-122115
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	15	4.9	< 0.50 U	< 0.50 U
1,1,2-TRICHLOROETHANE	1	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,1-DICHLOROETHANE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,1-DICHLOROETHENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,2,4-TRICHLOROBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DIBROMO-3-CHLOROPROPANE	0.04	< 0.75 U	< 0.75 U	< 0.75 U	< 0.75 U
1,2-DIBROMOETHANE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DICHLOROBENZENE	3	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DICHLOROETHANE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DICHLOROETHENE, TOTAL	5	2.7	< 1.0 U	< 1.0 U	< 1.0 U
1,2-DICHLOROPROPANE	1	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,3-DICHLOROBENZENE	3	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,4-DICHLOROBENZENE	3	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,4-DIOXANE (Method 8270D_SIM)	NL	9.3	< 0.17 U	5.0	0.70
2-BUTANONE	50	< 2.5 UJ	< 2.5 U	< 2.5 U	< 2.5 U
2-HEXANONE	50	< 2.5 UJ	< 2.5 U	< 2.5 U	< 2.5 U
4-METHYL-2-PENTANONE	NL	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 U
ACETONE	50	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ
BENZENE	1	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
BROMODICHLOROMETHANE	50	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
BROMOFORM	50	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
BROMOMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
CARBON DISULFIDE	60	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CARBON TETRACHLORIDE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROETHANE	5	< 1.0 UJ	< 1.0 U	< 1.0 UJ	< 1.0 UJ
CHLOROFORM	7	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CHLOROMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 UJ	< 1.0 UJ
CIS-1,2-DICHLOROETHENE	5	2.7	< 0.50 U	< 0.50 U	< 0.50 U
CIS-1,3-DICHLOROPROPENE	0.4	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CYCLOHEXANE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
DIBROMOCHLOROMETHANE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
DICHLORODIFLUOROMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
ETHYLBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
ISOPROPYLBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
M- AND P-XYLENE	NL	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
METHYL ACETATE	NL	< 0.75 U	< 0.75 U	< 0.75 UJ	< 0.75 UJ
METHYL CYCLOHEXANE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
METHYL TERT-BUTYL ETHER	10	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
METHYLENE CHLORIDE	5	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 U
O-XYLENE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
STYRENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TETRACHLOROETHENE	5	6.4	< 0.50 UJ	< 0.50 UJ	0.59 J
TOLUENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TRANS-1,2-DICHLOROETHENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TRANS-1,3-DICHLOROPROPENE	0.4	< 0.50 UJ	< 0.50 U	< 0.50 UJ	< 0.50 UJ
TRICHLOROETHENE	5	140	0.36 J	6.1	1.5
TRICHLOROFLUOROMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
VINYL CHLORIDE	2	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
XYLENES, TOTAL	5	< 1.5 U	< 1.5 U	< 1.5 U	< 1.5 U

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

Location	NYSDEC Groundwater Guidance or Standard Value (Note 1)	RE123D3	RE120D1	RE120D2	RE120D3
Sample Date		12/21/2015	12/18/2015	12/29/2015	12/29/2015
Sample ID		RE123D3-GW-122115	RE120D1-GW-121815	RE120D2-GW-122915	RE120D3-GW-122915
Sample type code		N	N	N	N
<b>VOC 8260C (ug/L)</b>					
1,1,1-TRICHLOROETHANE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,1,2,2-TETRACHLOROETHANE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	5	< 0.50 U	42	25	3.1
1,1,2-TRICHLOROETHANE	1	< 0.50 U	1.4	0.64 J	< 0.50 U
1,1-DICHLOROETHANE	5	< 0.50 U	3.2	1.1	< 0.50 U
1,1-DICHLOROETHENE	5	< 0.50 U	23	5.7	< 0.50 U
1,2,4-TRICHLOROBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DIBROMO-3-CHLOROPROPANE	0.04	< 0.75 U	< 0.75 U	< 0.75 U	< 0.75 U
1,2-DIBROMOETHANE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DICHLOROBENZENE	3	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DICHLOROETHANE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,2-DICHLOROETHENE, TOTAL	5	< 1.0 U	4.0	3.4	< 1.0 U
1,2-DICHLOROPROPANE	1	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,3-DICHLOROBENZENE	3	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,4-DICHLOROBENZENE	3	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
1,4-DIOXANE (Method 8270D_SIM)	NL	< 0.17 U	12	8.8	0.28
2-BUTANONE	50	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 U
2-HEXANONE	50	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 U
4-METHYL-2-PENTANONE	NL	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 U
ACETONE	50	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ	< 2.5 UJ
BENZENE	1	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
BROMODICHLOROMETHANE	50	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
BROMOFORM	50	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
BROMOMETHANE	5	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
CARBON DISULFIDE	60	0.56 J	< 0.50 U	< 0.50 U	< 0.50 U
CARBON TETRACHLORIDE	5	< 0.50 U	0.79 J	0.69 J	< 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 U	< 1.0 U
CHLOROFORM	7	< 0.50 U	0.99 J	0.77 J	< 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	4.0	3.4	< 0.50 U
CIS-1,3-DICHLOROPROPENE	0.4	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
CYCLOHEXANE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
DIBROMOCHLOROMETHANE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
DICHLORODIFLUOROMETHANE	5	< 1.0 U	0.38 J	< 1.0 U	< 1.0 U
ETHYLBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
ISOPROPYLBENZENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
M- AND P-XYLENE	NL	< 1.0 U	< 1.0 U	< 1.0 U	< 1.0 U
METHYL ACETATE	NL	< 0.75 UJ	< 0.75 UJ	< 0.75 U	< 0.75 U
METHYL CYCLOHEXANE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
METHYL TERT-BUTYL ETHER	10	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
METHYLENE CHLORIDE	5	< 2.5 U	< 2.5 U	< 2.5 U	< 2.5 U
O-XYLENE	NL	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
STYRENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TETRACHLOROETHENE	5	< 0.50 UJ	2.1 J	3.7 J	< 0.50 UJ
TOLUENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TRANS-1,2-DICHLOROETHENE	5	< 0.50 U	< 0.50 U	< 0.50 U	< 0.50 U
TRANS-1,3-DICHLOROPROPENE	0.4	< 0.50 UJ	< 0.50 UJ	< 0.50 U	< 0.50 U
TRICHLOROETHENE	5	< 0.50 U	1300	680	29
TRICHLOROFLUOROMETHANE	5	< 1.0 U	0.39 J	0.26 J	< 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)	RE108D1	RE108D2	RE121D1	RE121D2
Sample Date		12/22/2015	12/22/2015	12/21/2015	12/21/2015
Sample ID		RE108D1-GW-122215	RE108D2-GW-122215	RE121D1-GW-122115	RE121D2-GW-122115
Sample type code		N	N	N	N
<b>VOC 8260C (ug/L)</b>					
1,1,1-TRICHLOROETHANE	5	< 0.50 U	1.4 J	0.38 J	0.48 J
1,1,2,2-TETRACHLOROETHANE	5	< 0.50 U	< 2.5 U	< 0.50 U	< 0.50 U
1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	5	1.4	6.2	8.3	17 J
1,1,2-TRICHLOROETHANE	1	< 0.50 U	< 2.5 U	< 0.50 U	0.64 J
1,1-DICHLOROETHANE	5	< 0.50 U	5.1	< 0.50 U	0.51 J
1,1-DICHLOROETHENE	5	0.44 J	9.0	2.1	3.1 J
1,2,4-TRICHLOROBENZENE	5	< 0.50 U	< 2.5 U	< 0.50 U	< 0.50 U
1,2-DIBROMO-3-CHLOROPROPANE	0.04	< 0.75 U	< 3.8 U	< 0.75 U	< 0.75 U
1,2-DIBROMOETHANE	NL	< 0.50 U	< 2.5 U	< 0.50 U	< 0.50 U
1,2-DICHLOROBENZENE	3	< 0.50 U	< 2.5 U	< 0.50 U	< 0.50 U
1,2-DICHLOROETHANE	5	< 0.50 U	< 2.5 U	0.38 J	< 0.50 U
1,2-DICHLOROETHENE, TOTAL	5	0.61 J	9.0 J	0.96 J	2.1 J
1,2-DICHLOROPROPANE	1	< 0.50 U	< 2.5 U	< 0.50 U	< 0.50 U
1,3-DICHLOROBENZENE	3	< 0.50 U	< 2.5 U	< 0.50 U	< 0.50 U
1,4-DICHLOROBENZENE	3	< 0.50 U	< 2.5 U	< 0.50 U	< 0.50 U
1,4-DIOXANE (Method 8270D_SIM)	NL	6.7	8.8	6.8	4.9
2-BUTANONE	50	< 2.5 UJ	< 12 UJ	< 2.5 UJ	< 2.5 UJ
2-HEXANONE	50	< 2.5 UJ	< 12 UJ	< 2.5 UJ	< 2.5 UJ
4-METHYL-2-PENTANONE	NL	< 2.5 U	< 12 U	< 2.5 U	< 2.5 U
ACETONE	50	< 2.5 UJ	< 12 UJ	< 2.5 UJ	< 2.5 UJ
BENZENE	1	< 0.50 U	< 2.5 U	< 0.50 U	< 0.50 U
BROMODICHLOROMETHANE	50	< 0.50 U	< 2.5 U	< 0.50 U	< 0.50 U
BROMOFORM	50	< 0.50 U	< 2.5 U	< 0.50 U	< 0.50 U
BROMOMETHANE	5	< 1.0 U	< 5.0 U	< 1.0 U	< 1.0 U
CARBON DISULFIDE	60	< 0.50 U	< 2.5 U	< 0.50 U	< 0.50 U
CARBON TETRACHLORIDE	5	< 0.50 U	1.8 J	0.34 J	3.1 J
CHLOROBENZENE	5	< 0.50 U	< 2.5 U	< 0.50 U	< 0.50 U
CHLOROETHANE	5	< 1.0 UJ	< 5.0 UJ	< 1.0 UJ	< 1.0 UJ
CHLOROFORM	7	< 0.50 U	4.4 J	0.47 J	1.7 J
CHLOROMETHANE	5	< 1.0 U	< 5.0 U	< 1.0 U	< 1.0 U
CIS-1,2-DICHLOROETHENE	5	0.61 J	9.0	0.96 J	2.1 J
CIS-1,3-DICHLOROPROPENE	0.4	< 0.50 U	< 2.5 U	< 0.50 U	< 0.50 U
CYCLOHEXANE	NL	< 0.50 U	< 2.5 U	< 0.50 U	< 0.50 U
DIBROMOCHLOROMETHANE	5	< 0.50 U	< 2.5 U	< 0.50 U	< 0.50 U
DICHLORODIFLUOROMETHANE	5	< 1.0 U	< 5.0 U	2.2	0.85 J
ETHYLBENZENE	5	< 0.50 U	< 2.5 U	< 0.50 U	< 0.50 U
ISOPROPYLBENZENE	5	< 0.50 U	< 2.5 U	< 0.50 U	< 0.50 U
M- AND P-XYLENE	NL	< 1.0 U	< 5.0 U	< 1.0 U	< 1.0 U
METHYL ACETATE	NL	< 0.75 U	< 3.8 U	< 0.75 U	< 0.75 U
METHYL CYCLOHEXANE	NL	< 0.50 U	< 2.5 U	< 0.50 U	< 0.50 U
METHYL TERT-BUTYL ETHER	10	< 0.50 U	< 2.5 U	< 0.50 U	< 0.50 U
METHYLENE CHLORIDE	5	< 2.5 U	< 12 U	< 2.5 U	< 2.5 U
O-XYLENE	NL	< 0.50 U	< 2.5 U	< 0.50 U	< 0.50 U
STYRENE	5	< 0.50 U	< 2.5 U	< 0.50 U	< 0.50 U
TETRACHLOROETHENE	5	1.2	< 2.5 U	< 0.50 U	< 0.50 U
TOLUENE	5	< 0.50 U	< 2.5 U	< 0.50 U	< 0.50 U
TRANS-1,2-DICHLOROETHENE	5	< 0.50 U	< 2.5 U	< 0.50 U	< 0.50 U
TRANS-1,3-DICHLOROPROPENE	0.4	< 0.50 UJ	< 2.5 UJ	< 0.50 UJ	< 0.50 UJ
TRICHLOROETHENE	5	110	2900 J	29	480
TRICHLOROFLUOROMETHANE	5	< 1.0 U	< 5.0 U	< 1.0 U	< 1.0 U
VINYL CHLORIDE	2	< 1.0 U	< 5.0 U	< 1.0 U	< 1.0 U
XYLENES, TOTAL	5	< 1.5 U	< 7.5 U	< 1.5 U	< 1.5 U

**Notes:**

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

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

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

Yellow highlighted values exceed Groundwater Standards or guidance value

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

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

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

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

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

TABLE 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	12/17/2015	15.26	4.63	0.078	0.47	225.0	0.10	34.01	850
TT101D1	12/17/2015	15.23	5.03	0.082	0.87	231.8	0.10	35.01	850
TT101D2	12/21/2015	15.30	5.09	0.034	7.41	157.4	0	35.38	700
RE103D1	12/14/2015	15.19	5.25	0.137	4.24	102.6	0.14	40.87	500
RE103D2	12/14/2015	14.87	5.38	0.039	7.30	194.9	0.21	40.15	490
RE103D3	12/14/2015	15.03	5.58	0.041	5.31	161.4	0.66	41.22	500
RE104D1	12/15/2015	14.23	5.09	0.053	4.67	254	0.39	37.02	500
RE104D2	12/15/2015	14.38	4.81	0.019	7.20	188.1	2.80	39.38	500
RE104D3	12/15/2015	14.28	4.49	0.016	5.78	120.1	2.7	39.9	600
RE105D1	12/17/2015	14.43	4.65	0.085	2.72	129.3	0.17	38.50	675
RE105D2	12/17/2015	14.60	4.58	0.056	5.28	205.7	0.10	36.78	500
RE107D1	12/18/2015	13.91	5.52	0.083	4.02	152.9	414.00	41.80	525
RE107D2	12/18/2015	14.56	5.69	0.077	3.49	238.1	49.5	42.68	700
RE107D3	12/29/2015	14.03	5.65	0.033	7.42	20.7	48.4	42.21	250
RE108D1	12/22/2015	14.56	5.10	0.096	7.63	282.1	0.55	40.24	600
RE108D2	12/22/2015	14.32	4.70	0.082	5.42	38.1	0.20	40.72	550
RE114D1	12/21/2015	13.85	5.82	0.073	2.62	148.7	21.00	31.64	500
RE114D2	12/16/2015	14.28	5.99	0.070	0.66	100.9	34.1	31.84	500
RE114D3	12/16/2015	14.35	5.45	0.033	5.78	245.2	1.42	32.26	500
RE120D1	12/18/2015	14.43	4.79	0.094	1.81	104.9	0.42	36.61	525
RE120D2	12/29/2015	14.87	5.22	0.080	5.91	42.6	1.45	36.70	500
RE120D3	12/29/2015	14.75	4.88	0.026	4.72	98.1	1.2	37.14	500
RE121D1	12/21/2015	15.97	5.50	0.071	0.57	154.1	4.21	34.40	863
RE121D2	12/21/2015	15.15	5.16	0.071	2.77	-54.5	8.85	34.41	450
RE122D1	12/15/2015	14.54	5.73	0.079	2.92	89.8	1.02	42.48	600
RE122D2	12/15/2015	14.80	5.21	0.071	5.13	213.7	0.68	43.00	600
RE122D3	12/15/2015	14.60	4.67	0.020	3.32	180.4	14.3	42.82	500
RE123D1	12/21/2015	15.59	5.25	0.087	9.32	172.8	3.08	47.63	700
RE123D2	12/21/2015	14.38	5.54	0.022	8.37	202.9	4.52	48.99	550
RE123D3	12/21/2015	14.77	5.72	0.066	0.66	-82.8	9.25	48.95	600

°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 through BPOW5-7, Fourth Quarter 2015**  
**Operable Unit 2 (Groundwater),**  
**Bethpage, New York**

Well:	BPOW5-1	BPOW5-2	BPOW5-3	BPOW5-4	BPOW5-5	BPOW5-6	BPOW5-7
Sample ID:	BPOW5-1	BPOW5-2	BPOW5-3	BPOW5-4	BPOW5-5	BPOW5-6	BPOW5-7
Date:	11/12/2015	11/12/2015	12/3/2015	11/16/2015	11/13/2015	11/13/2015	11/20/2015
CONSTITUENT							
Units (ug/L)							
<b><u>Volatile Organic Compounds (VOCs) <sup>(1)</sup></u></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 B	< 5.0	< 5.0	< 5.0 B	< 5.0 B	< 5.0	< 5.0 B
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>0.15 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 B	< 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	<b>0.49 J</b>
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>(2)</sup></b>	<b>0</b>	<b>0</b>	<b>0</b>	<b>0</b>	<b>0</b>	<b>0.15</b>	<b>0.49</b>
<b>1,4-Dioxane <sup>(3)</sup></b>	<b>&lt;0.22</b>	<b>&lt; 0.23</b>	<b>0.39</b>	<b>0.28</b>	<b>0.42</b>	<b>&lt; 0.21</b>	<b>&lt; 0.21</b>

See last page for Notes and Abbreviations

**Table 4.**  
**Concentrations of Volatile Organic Compounds**  
**and 1,4-Dioxane in Monitoring Wells BPOW 5-1 through BPOW 5-7, Fourth Quarter 2015**  
**Operable Unit 2 (Groundwater),**  
**Bethpage, New York**

**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).

<b>Bold</b>	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	Constituent detected in associated blank sample
<0.50	Constituent not detected above its laboratory detection limit

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

Well:	BPOW6-1	BPOW6-2	BPOW6-3	BPOW6-4	BPOW6-4	BPOW6-5	BPOW6-6	
Sample ID:	BPOW6-1	BPOW6-2	BPOW6-3	BPOW6-4	REP120115PP1	BPOW6-5	BPOW6-6	
Date:	11/30/2015	11/30/2015	12/1/2015	12/1/2015	12/1/2015	12/2/2015	12/2/2015	
CONSTITUENT								
Units (ug/L)								
<b><u>Volatile Organic Compounds (VOCs) <sup>(1)</sup></u></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 B	< 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>0.89</b>	<b>0.40 J</b>	
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>(2)</sup></b>	<b>0</b>	<b>0</b>	<b>0</b>	<b>0</b>	<b>0</b>	<b>0.89</b>	<b>0.4</b>	
<b>1,4-Dioxane <sup>(3)</sup></b>	<b>&lt; 0.22</b>	<b>&lt; 0.22</b>	<b>&lt; 0.23</b>	<b>&lt; 0.22</b>	<b>&lt; 0.21</b>	<b>&lt; 0.21</b>	<b>&lt; 0.22</b>	

See last page for Notes and Abbreviations.

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

**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).

<b>Bold</b>	Constituent detected
TCL	Target Compound List
VOC	Volatile Organic Compound
USEPA	United States Environmental Protection Agency
REP	Blind duplicate sample
SIM	Selected Ion Monitoring
µg/L	Micrograms per liter
J	Constituent value is estimated
<0.50	Constituent not detected above its laboratory detection limit

**Table 6.**  
**Concentrations of Volatile Organic Compounds**  
**and 1,4-Dioxane in Monitoring Wells RE117D1,**  
**RE117D2, RE118D1 and RE119D1, Fourth Quarter 2015**  
**Operable Unit 2 (Groundwater),**  
**Bethpage, New York**

Well:	RE117D1	RE117D2	RE118D1	RE119D1
Sample ID:	RE117D1	RE117D2	RE118D1	RE119D1
Date:	11/18/2015	11/18/2015	11/23/2015	11/20/2015
CONSTITUENT				
Units (ug/L)				
<b><u>Volatile Organic Compounds (VOCs) <sup>(1)</sup></u></b>				
1,1,1-Trichloroethane	< 1.0	< 1.0 J	< 1.0	< 1.0
1,1,2,2-Tetrachloroethane	< 1.0	< 1.0 J	< 1.0	< 1.0
1,1,2-trichloro-1,2,2-trifluoroethane	< 5.0	< 5.0	< 5.0	< 5.0
1,1,2-Trichloroethane	< 1.0	< 1.0 J	< 1.0	< 1.0
1,1-Dichloroethane	< 1.0	< 1.0 J	< 1.0	< 1.0
1,1-Dichloroethene	< 1.0	< 1.0 J	< 1.0	< 1.0
1,2-Dichloroethane	< 1.0	< 1.0 J	< 1.0	< 1.0
1,2-Dichloropropane	< 1.0	< 1.0 J	< 1.0	< 1.0
2-Butanone (MEK)	< 10	< 10 J	< 10	< 10
2-Hexanone	< 5.0	< 5.0 J	< 5.0	< 5.0
4-methyl-2-pentanone (MIK)	< 5.0	< 5.0 J	< 5.0	< 5.0
Acetone	< 10	< 10 J	< 10	< 10
Benzene	< 0.50	< 0.50 J	< 0.50	< 0.50
Bromodichloromethane	< 1.0	< 1.0 J	< 1.0	< 1.0
Bromoform	< 1.0	< 1.0 J	< 1.0	< 1.0
Bromomethane	< 2.0	< 2.0 J	< 2.0	< 2.0
Carbon Disulfide	< 2.0	< 2.0 J	< 2.0	< 2.0
Carbon tetrachloride	< 1.0	< 1.0 J	< 1.0	< 1.0
Chlorobenzene	< 1.0	< 1.0 J	< 1.0	< 1.0
Chloroethane	< 1.0	< 1.0 J	< 1.0	< 1.0
Chloroform	< 1.0	< 1.0 J	< 1.0	< 1.0
Chloromethane	< 1.0	< 1.0 J	< 1.0	< 1.0
cis-1,2-dichloroethene	< 1.0	< 1.0 J	< 1.0	< 1.0
cis-1,3-dichloropropene	< 1.0	< 1.0 J	< 1.0	< 1.0
Dibromochloromethane	< 1.0	< 1.0 J	< 1.0	< 1.0
Ethylbenzene	< 1.0	< 1.0 J	< 1.0	< 1.0
Methylene Chloride	< 2.0	< 2.0 J	< 2.0	< 2.0
Styrene	< 1.0	< 1.0 J	< 1.0	< 1.0
Tetrachloroethene	< 1.0	< 1.0 J	< 1.0	< 1.0
Toluene	<b>0.75 J</b>	<b>0.98 J</b>	<b>0.57 J</b>	<b>0.72</b>
trans-1,2-dichloroethene	< 1.0	< 1.0 J	< 1.0	< 1.0
trans-1,3-dichloropropene	< 1.0	< 1.0 J	< 1.0	< 1.0
Trichloroethylene	<b>9.4</b>	< 1.0 J	< 1.0	< 1.0
Vinyl Chloride	< 1.0	< 1.0 J	< 1.0	< 1.0
Xylene-o	< 1.0	< 1.0 J	< 1.0	< 1.0
Xylenes - m,p	< 1.0	< 1.0 J	< 1.0	< 1.0
<b>Total VOCs <sup>(2)</sup></b>	<b>10.15</b>	<b>0.98</b>	<b>0.57</b>	<b>0.72</b>
<b>1,4-Dioxane <sup>(3)</sup></b>	< 0.21	< 0.20	< 0.22	< 0.22

See last page for Notes and Abbreviations.



**Table 6.  
Concentrations of Volatile Organic Compounds  
and 1,4-Dioxane in Monitoring Wells RE-117D1,  
RE-117D2, RE-118D1 and RE-119D1, Fourth Quarter  
2015 Operable Unit 2 (Groundwater),  
Bethpage, New York**

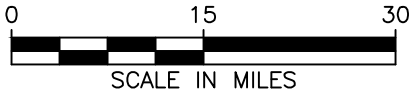
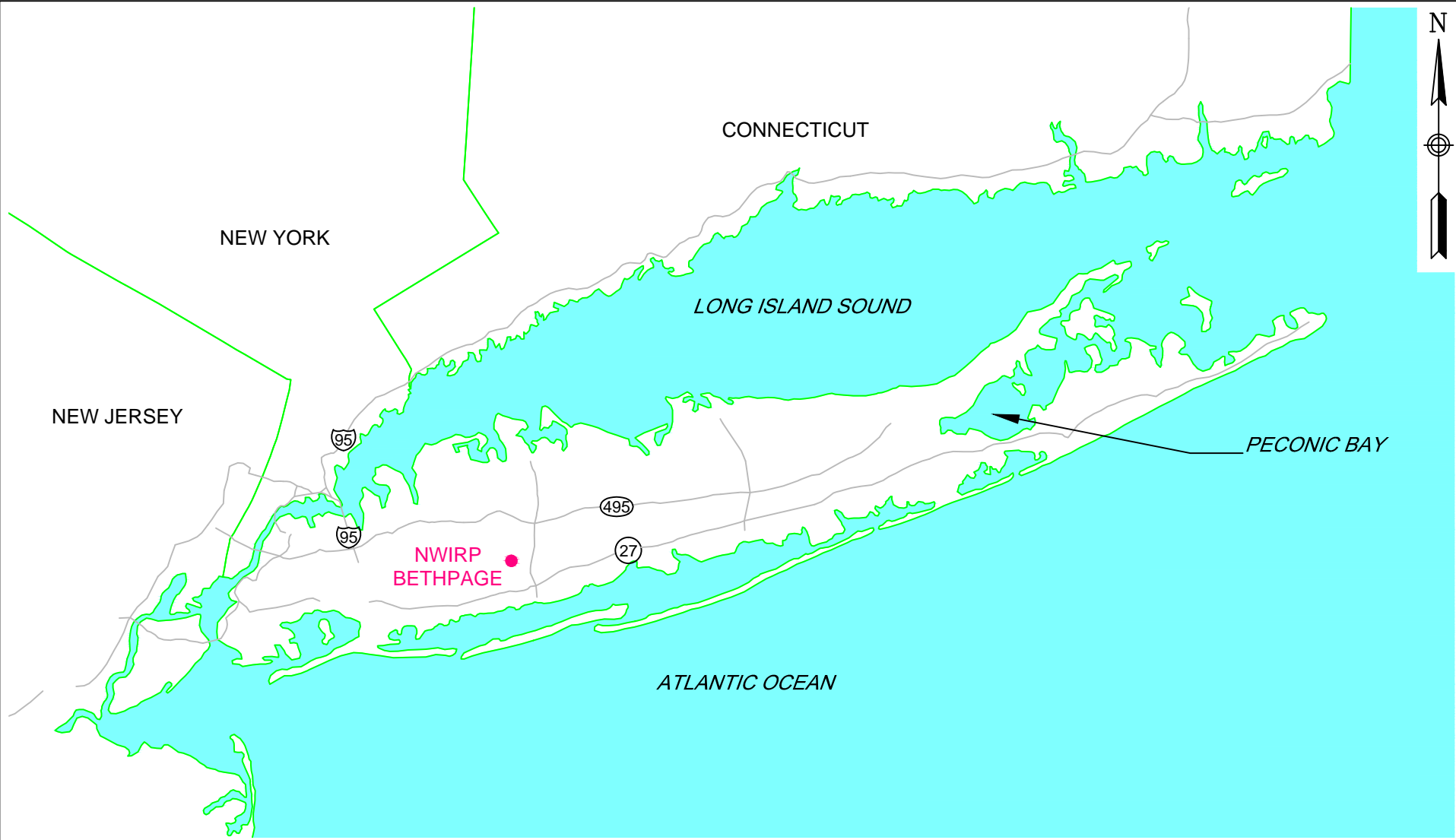
**Notes and Abbreviations:**

- (1) Samples were analyzed for the TCL VOCs using Method 8260C.
- (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).

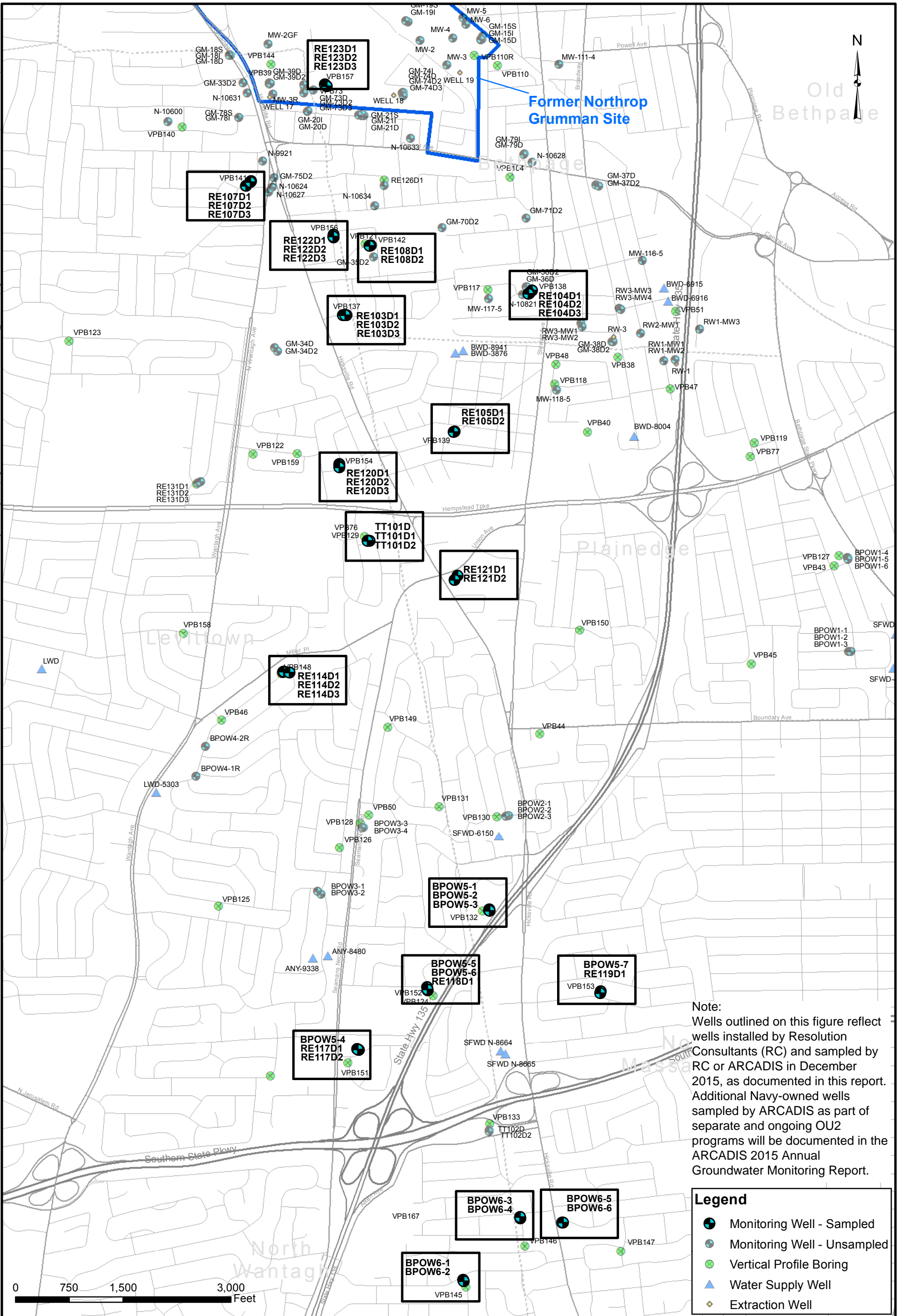
<b>Bold</b>	Constituent detected
TCL	Target Compound List
VOC	Volatile Organic Compound
USEPA	United States Environmental Protection Agency
SIM	Selected Ion Monitoring
µg/L	Micrograms per liter
J	Constituent value is estimated
<0.50	Constituent not detected above its laboratory detection limit

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



Note:  
Wells outlined on this figure reflect wells installed by Resolution Consultants (RC) and sampled by RC or ARCADIS in December 2015, as documented in this report. Additional Navy-owned wells sampled by ARCADIS as part of separate and ongoing OU2 programs will be documented in the ARCADIS 2015 Annual Groundwater Monitoring Report.

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



LOCATION MAP  
DECEMBER 2015 GROUNDWATER SAMPLING  
NAVAL WEAPONS INDUSTRIAL RESERVE PLANT  
BETHPAGE, NEW YORK

CONTRACT NUMBER N62470-11-D8013	CTO NUMBER WE 15
APPROVED BY EV	DATE 2/9/2016
APPROVED BY	DATE
FIGURE NO. 2	REV 0

## Appendices

Appendix A

Groundwater Sampling Forms – Resolution Consultants



RESOLUTION  
CONSULTANTS

Well ID: RE103D1

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 12/14 / 15 Time: Start 1135 am/pm  
 Project No: 60266526 Finish 1330 am/pm  
 Site Location: Bethpage  
 Weather Conds: Cloudy 60° Collector(s): Caterle Foster

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

a. Total Well Length 640 c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material 4-inch PVC  
 b. Water Table Depth 41.48 d. Calculated System Volume (see back) 9.8

### 2. WELL PURGE DATA

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

#### b. Acceptance Criteria defined (see workplan)

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

c. Field Testing Equipment used: \_\_\_\_\_ Make \_\_\_\_\_ Model \_\_\_\_\_ Serial Number \_\_\_\_\_

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
11:40	-	15.31	7.39	0.106	4.35	120.5	-	500	41.63	
11:50	-	15.27	6.02	0.105	1.62	101.8	1.41	500	41.65	Clear
12:00	-	15.26	5.76	0.106	1.47	96.5	1.00	500	41.66	
12:05	-	15.27	5.56	0.108	3.17	97.0	-	500	41.60	Clear
12:15	5 Gal	15.30	5.47	0.108	3.44	97.9	0.14	500	41.40	
12:25	-	15.28	5.37	0.109	3.63	98.0	-	500	41.09	Clear

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>RE103D1-GW-12/14/15</u>	40-mL vials	3	HCl	VOCs	<u>1300</u>
	1-L amber	2	none	1,4-Dioxane	

Comments \_\_\_\_\_

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







Well ID: RE10302

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 12/14/15 Time: Start 1045 am/pm  
 Project No: 60266526 Finish 1330 am/pm  
 Site Location: Avoca  
 Weather Conds: Cloudy 55° Collector(s): \_\_\_\_\_

### 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 41.41 d. Calculated System Volume (see back) 13.1

### 2. WELL PURGE DATA

a. Purge Method: Geotech bladder pump with drop tube assembly  
 b. Acceptance Criteria defined (see workplan)  
 - Temperature ± 3% - Turbidity ± 10% - D.O. ± 10% (values >0.5 mg/L)  
 - pH ± 0.1 unit - ORP ± 10mV  
 - Sp. Cond. ± 3% - Drawdown < 0.3' Remove a minimum 1 screen volume 13.1

c. Field Testing Equipment used:

Make	Model	Serial Number

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
1115										OK
1130		14.91	5.52	0.042	5.96	147.2		300		
1135		14.95	5.49	0.040	5.36	151.9		475		
1140		14.95	5.45	0.047	5.46	156.8		475	41.10	
1145		14.94	5.40	0.040	6.87	167.2	1.43	490	41.07	
1150		14.93	5.39	0.039	7.24	170.9	0.97	490	41.05	

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>RE103102-GW 12/14/15</u>	40-mL vials	3	HCl	VOCs	1315
	1-L amber	2	none	1,4-Dioxane	

Comments \_\_\_\_\_

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





RESOLUTION  
CONSULTANTS

Well ID: RE10303

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 12/14/15 Time: Start 11:30 am/pm  
 Project No: 60266526 Finish 1330 am/pm  
 Site Location: Bethpage, NY  
 Weather Conds: Foggy Collector(s): F.B.M./P.K./R.P.

### 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 40.95 d. Calculated System Volume (see back) 98

### 2. WELL PURGE DATA

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

#### b. Acceptance Criteria defined (see workplan)

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

c. Field Testing Equipment used: \_\_\_\_\_ Make \_\_\_\_\_ Model \_\_\_\_\_ Serial Number \_\_\_\_\_

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
11:30	-	15.08	6.82	0.043	7.02	112.4	-	-	-	Clear.
11:50	-	15.08	6.52	0.042	5.87	119.8	2.02	475	41.30	Clear
12:00	-	15.07	6.08	0.041	5.46	136.5	-	500	41.34	Clear
12:05	-	15.07	6.07	0.041	5.49	142.6	0.86	-	41.35	Clear
12:15		15.11	5.60	0.041	5.25	149.9	0.98		41.35	
12:25		15.06	5.63	0.041	5.62	158.8	0.57		41.28	

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
<u>RE10303-GW-121415</u>	<u>40-mL vials</u>	<u>3</u>	<u>HCl</u>	<u>VOCs</u>	<u>1245</u>
	<u>1-L amber</u>	<u>2</u>	<u>none</u>	<u>1,4-Dioxane</u>	

Comments: ms/mss collected here

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





Well ID: RE10401

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 12/15/15 Time: Start 745 am/pm  
 Project No: 60266526 Finish 1100 am/pm  
 Site Location: Hilltop  
 Weather Conds: Sunny 66° Collector(s): \_\_\_\_\_

- 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 36.49 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%
  - pH ± 0.1 unit
  - Sp. Cond. ± 3%
  - Turbidity ± 10%
  - ORP ± 10mV
  - Drawdown < 0.3'
  - D.O. ± 10% (values >0.5 mg/L)
- Remove a minimum 1 screen volume
- c. Field Testing Equipment used:
- | Make       | Model          | Serial Number     |
|------------|----------------|-------------------|
| <u>YSI</u> | <u>556 MTS</u> | <u>05G1942 AC</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
9:00	-	14.62	6.22	0.065	11.84	170.1	-	450	36.60	Clear
9:10	-	13.95	5.10	0.054	6.01	230.3	-	450	-	Clear
9:15	-	13.95	5.06	0.056	6.07	247.7	0.059	500	36.71	Clear
9:20	-	14.01	5.07	0.055	5.76	263.9	-	500	36.71	Clear
9:25	-	14.03	5.07	0.054	5.58	269.7	1.09	500	36.75	Clear
9:40	-	14.08	5.08	0.053	5.19	281.6	-	500	36.81	Clear

- d. Acceptance criteria pass/fail
- |                                     | Yes                      | No                       | N/A                      |
|-------------------------------------|--------------------------|--------------------------|--------------------------|
| Has required volume been removed    | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Has required turbidity been reached | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Have parameters stabilized          | <input 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>RE10401-GW-12/5X</u>	40-mL vials	3	HCl	VOCs	<u>1100</u>
	1-L amber	2	none	1,4-Dioxane	

Comments \_\_\_\_\_

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







Well ID: RE10402

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 12/15/15 Time: Start 745 am/pm  
 Project No: 60266526 Finish 1100 am/pm  
 Site Location: Hilltop  
 Weather Conds: sunny 60° Collector(s): \_\_\_\_\_

**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 39.97 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% - Turbidity ± 10% - D.O. ± 10% (values >0.5 mg/L)  
 - 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>450231X</u>
<u>Hanna</u>	<u>98703</u>	<u>480211X</u>

Time (24hr)	Volume Removed (Liters) <sup>1</sup>	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>845</u>		<u>14.75</u>	<u>6.97</u>	<u>0.045</u>	<u>12.03</u>	<u>112.0</u>	<u>2.56</u>	<u>500</u>	<u>40.01</u>	<u>OH</u>
<u>890</u>	<u>2.5</u>	<u>14.38</u>	<u>4.55</u>	<u>0.021</u>	<u>12.71</u>	<u>131.9</u>	<u>1.56</u>	<u>500 ml</u>	<u>40.0</u>	<u>clear</u>
<u>905</u>							<u>2.22</u>			<u>HOLD</u>
<u>915</u>		<u>14.24</u>	<u>5.18</u>	<u>0.020</u>	<u>7.90</u>	<u>150.0</u>		<u>475</u>		<u>OH</u>
<u>920</u>	<u>5901</u>	<u>14.22</u>	<u>5.14</u>	<u>0.019</u>	<u>7.84</u>	<u>146.4</u>		<u>475</u>	<u>3.9.98</u>	
<u>925</u>		<u>14.28</u>	<u>5.01</u>	<u>0.019</u>	<u>7.49</u>	<u>148.5</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
<u>RE10402-GW-121515</u>	<u>40-mL vials</u>	<u>3</u>	<u>HCl</u>	<u>VOCs</u>	<u>1030</u>
	<u>1-L amber</u>	<u>2</u>	<u>none</u>	<u>1,4-Dioxane</u>	
<u>DUPLICATED - GW - 121515</u>					<u>1045</u>

Comments: 905, blowing air, pull pump & reset bladder

Signature: Paul Knecht Date: 12/15/15







RESOLUTION CONSULTANTS

Well ID: RE104D3

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 12/15/15 Time: Start 745 am/pm  
 Project No: 60266526 Finish 1100 am/pm  
 Site Location: Hilltop  
 Weather Conds: slurry 60° Collector(s): ER

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

a. Total Well Length 785 c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material 4-inch PVC  
 b. Water Table Depth 39.67 d. Calculated System Volume (see back) \_\_\_\_\_ 13.1 gal to remove

### 2. WELL PURGE DATA

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

#### b. Acceptance Criteria defined (see workplan)

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

#### c. Field Testing Equipment used:

Make	Model	Serial Number
<u>YSI</u>		<u>07H100384</u>
<u>Hanna Turbidity meter</u>		

Time (24hr)	Volume Removed (Liters) gal	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>0855</u>		<u>13.86</u>	<u>4.73</u>	<u>0.022</u>	<u>6.11</u>	<u>150.3</u>	<u>7.29</u>	<u>350</u>	<u>39.71</u>	<u>clear</u>
<u>0905</u>	<u>1.25</u>	<u>14.15</u>	<u>4.53</u>	<u>0.019</u>	<u>5.25</u>	<u>114.7</u>	<u>9.81</u>	<u>550</u>	<u>39.70</u>	<u>clear</u>
<u>0915</u>	<u>3.5</u>	<u>14.15</u>	<u>4.39</u>	<u>0.017</u>	<u>5.76</u>	<u>110.2</u>	<u>9.81</u>	<u>550</u>	<u>39.70</u>	<u>clear</u>
<u>0925</u>	<u>4.5</u>	<u>14.17</u>	<u>4.39</u>	<u>0.017</u>	<u>5.90</u>	<u>111.9</u>	<u>9.80</u>	<u>550</u>	<u>39.7</u>	<u>clear</u>
<u>0935</u>	<u>5.5</u>	<u>14.20</u>	<u>4.38</u>	<u>0.016</u>	<u>5.55</u>	<u>113.3</u>	<u>7.23</u>	<u>550</u>	<u>39.7</u>	<u>clear</u>
<u>0945</u>	<u>6.5</u>	<u>14.20</u>	<u>4.44</u>	<u>0.017</u>	<u>5.87</u>	<u>114.7</u>	<u>7.20</u>	<u>550</u>	<u>39.7</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>RE104D3-GW-121515</u>	<u>40-mL vials</u>	<u>3</u>	<u>HCl</u>	<u>VOCs</u>	<u>1040</u>
	<u>1-L amber</u>	<u>2</u>	<u>none</u>	<u>1,4-Dioxane</u>	<u>1040</u>

Comments \_\_\_\_\_

Signature Rich Page

Date 12-15-14





RESOLUTION CONSULTANTS

Well ID: RE10501

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 12/17/15 Time: Start 8:45 am/pm  
 Project No: 60266526 Finish 1:30 am/pm  
 Site Location: Lincoln  
 Weather Conds: overcast, drizzle 50° Collector(s): \_\_\_\_\_

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

a. Total Well Length 555 c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material 4-inch PVC  
 b. Water Table Depth 38.14 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%
  - pH ± 0.1 unit
  - Sp. Cond. ± 3%
  - Turbidity ± 10%
  - ORP ± 10mV
  - Drawdown < 0.3'
  - D.O. ± 10% (values >0.5 mg/L)
- Remove a minimum 1 screen volume

#### c. Field Testing Equipment used:

Make	Model	Serial Number
YSI	550 DDS	78291
Solinst	WLM	1169413K
Hanna	H198703	U80211X
MPIO	GED	U51644X

Pump on → 9:15

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
9:30		14.39	4.78	0.084	5.50	118.7	0.18	700	38.37	Clear / no odor
09:40	5 gallons	14.41	4.65	0.086	3.00	119.7	0.46	650	38.41	
09:50		14.42	4.66	0.084	2.90	119.4	0.39	650	38.44	
10:00		14.42	4.65	0.085	2.86	121.4	0.36	675	38.46	
10:05	10 gallons	14.43	4.66	0.084	2.82	123.2	0.23	675	38.47	
10:10		14.44	4.65	0.084	2.80	124.8	0.28	675	38.48	

#### 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"/> |

(continued on back)

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>RE10501-GW-121715</u>	<u>40-mL vials</u>	<u>3</u>	<u>HCl</u>	<u>VOCs</u>	<u>10:30</u>
<u>RE10501-GW-121715</u>	<u>1-L amber</u>	<u>2</u>	<u>none</u>	<u>1,4-Dioxane</u>	<u>10:30</u>

Comments \_\_\_\_\_

Signature \_\_\_\_\_

Date \_\_\_\_\_





RESOLUTION  
CONSULTANTS

Well ID: RE10502

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 12/17/15 Time: Start 845 am/pm  
 Project No: 60266526 Finish 1130 am/pm  
 Site Location: Lincoln  
 Weather Conds: overcast, drizzle 50° Collector(s): \_\_\_\_\_

### 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 38.80 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%
  - pH ± 0.1 unit
  - Sp. Cond. ± 3%
  - Turbidity ± 10%
  - ORP ± 10mV
  - Drawdown < 0.3'
  - D.O. ± 10% (values >0.5 mg/L)
- Remove a minimum 1 screen volume

#### c. Field Testing Equipment used:

Make	Model	Serial Number
<u>VST</u>	<u>556</u>	<u>50231</u>
<u>Hanna</u>	<u>99703</u>	<u>80211</u>

Time (24hr)	Volume		pH	Spec.			Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
	Removed (Liters)	Temp. (°C)		Cond. (mS/cm)	DO (mg/L)	ORP (mV)				
<u>930</u>										
<u>940</u>		<u>14.57</u>	<u>5.43</u>	<u>0.052</u>	<u>597</u>	<u>149.4</u>	<u>0.25</u>	<u>475</u>	<u>36.95</u>	
<u>945</u>		<u>14.61</u>	<u>5.25</u>	<u>0.052</u>	<u>4.53</u>	<u>159.7</u>				
<u>950</u>		<u>14.61</u>	<u>5.06</u>	<u>0.052</u>	<u>3.84</u>	<u>164.9</u>				
<u>955</u>		<u>14.61</u>	<u>4.95</u>	<u>0.052</u>	<u>3.56</u>	<u>169.4</u>		<u>500</u>	<u>36.82</u>	
<u>1000</u>		<u>14.60</u>	<u>4.85</u>	<u>0.053</u>	<u>4.64</u>	<u>179.4</u>				

d. Acceptance criteria pass/fail

Has required volume been removed	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	(continued on back)
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>RE10502-G10-0217K</u>	<u>40-mL vials</u>	<u>3</u>	<u>HCl</u>	<u>VOCs</u>	<u>1110</u>
	<u>1-L amber</u>	<u>2</u>	<u>none</u>	<u>1,4-Dioxane</u>	<u>1110</u>

Comments \_\_\_\_\_

Signature Paul Kallath Date 12/17/15







RESOLUTION  
CONSULTANTS

Well ID: RE10201

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 12/19/15 Time: Start 7:45 am/pm  
 Project No: 60266526 Finish 11:00 am/pm  
 Site Location: Crescent  
 Weather Conds: cloudy 50° Collector(s): \_\_\_\_\_

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

a. Total Well Length 530 c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material  
 b. Water Table Depth 42.20 d. Calculated System Volume (see back) 13.1 4-inch PVC

### 2. WELL PURGE DATA

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

c. Field Testing Equipment used:

Make	Model	Serial Number
<u>Ysi</u>	<u>556 mFS</u>	<u>050231X</u>
<u>HANNA</u>	<u>98703</u>	<u>080211</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
0834										
0835	0.5	13.26	6.14	0.088	7.70	204.	887	400		cloudy
0845	1	13.80	5.47	0.080	5.70	124.4	61.2		42.28	cloudy
850	1.5	13.84	5.53	0.080	5.55	29.7	55.1	325	42.28	
855	2.5	13.80	5.48	0.079	5.17	-5.5	38.5			
900	3	13.81	5.50	0.079	4.93	37.8	43.8			

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 type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>
Have parameters stabilized	<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>

If no or N/A - Explain below.  
turbidity greater than 10 NTU, purged for 2 hours

### 3. SAMPLE COLLECTION:

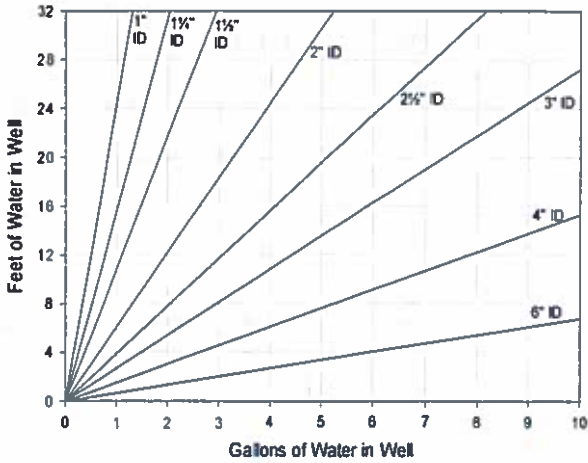
Method: Geotech bladder pump with drop tube assembly

Sample ID	Container Type	No. of Containers	Preservation	Analysis Req.	Time
<u>RE10201-GW-121815</u>	40-mL vials	3	HCl	VOCs	
	1-L amber	2	none	1,4-Dioxane	<u>1040</u>

Comments \_\_\_\_\_

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

One screen volume  
(4-inch well)

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

Well ID: RE107D1

(continued from front)

Time (24 hr)	Volume Removed (Liters)	Temp (°C)	pH	Specific Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
905		13.76	5.53	0.080	4.73	138.9			42.33	
910		13.69	5.48	0.080	4.57	140.5				
915		13.99	5.44	0.081	4.37	144.5				
920	5 gal									
925		13.97	5.47	0.082	4.30	147.5	149	950	42.12	cloudy
930	6	13.95	5.46	0.082	4.35	149.5	209			
935	7	14.03	5.47	0.082	4.0	150.8	252		41.88	
940	9	14.03	5.52	0.083	3.92	150.6	295	500		cloudy
945	9.5	14.08	5.53	0.083	3.87	151.7	326		41.75	
950	10.0	14.01	5.52	0.083	3.83	152.8	351			
955	10.2	13.98	5.53	0.083	3.82	151.7	390	500	41.78	cloudy
1000	11	14.06	5.51	0.083	3.84	148.9				
1005	12	13.99	5.58	0.083	3.91	149.5	433		41.84	cloudy
1010	12.5	13.94	5.49	0.083	3.92	151.0	384	525		cloudy
1015	13	13.97	5.51	0.083	3.90	150.2	387			
1020	14	13.92	5.51	0.083	3.94	149.2	449	525	41.85	
1025	14.5	13.86	5.53	0.083	4.07	149.8	414			cloudy
1030	15	13.89	5.52	0.083	4.07	149.0				
1035	16.5	13.91	5.52	0.083	4.05	152.9		525	41.80	
1240										sample





RESOLUTION  
CONSULTANTS

Well ID: RE10702

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 12/18/15 Time: Start 745 am/pm  
 Project No: 60266526 Finish 1100 am/pm  
 Site Location: Crescent  
 Weather Conds: cloudy 50° Collector(s): \_\_\_\_\_

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

a. Total Well Length 585 c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material 4-inch PVC  
 b. Water Table Depth 42.59 d. Calculated System Volume (see back) 13.1

### 2. WELL PURGE DATA

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

b. Acceptance Criteria defined (see workplan)  
 - Temperature  $\pm 3\%$  - Turbidity  $\pm 10\%$  - D.O.  $\pm 10\%$  (values  $>0.5$  mg/L)  
 - pH  $\pm 0.1$  unit - ORP  $\pm 10$ mV  
 - Sp. Cond.  $\pm 3\%$  - 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>473042x</u>
<u>Hanna</u>	<u>99703</u>	<u>80211</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>815</u>										<u>ON</u>
<u>830</u>	<u>2.5</u>	<u>14.58</u>	<u>5.63</u>	<u>0.072</u>	<u>5.97</u>	<u>241.2</u>	<u>8.85</u>	<u>700</u>	<u>42.6</u>	<u>cloudy</u>
<u>840</u>	<u>5</u>	<u>14.59</u>	<u>5.65</u>	<u>0.075</u>	<u>4.93</u>	<u>242.6</u>	<u>21.5</u>	<u>700</u>	<u>42.6</u>	<u>cloudy</u>
<u>845</u>	<u>6</u>	<u>14.59</u>	<u>5.66</u>	<u>0.075</u>	<u>4.72</u>	<u>241.2</u>	<u>21.2</u>	<u>700</u>	<u>42.6</u>	<u>cloudy</u>
<u>850</u>	<u>7.5</u>	<u>14.60</u>	<u>5.52</u>	<u>0.075</u>	<u>4.59</u>	<u>240.6</u>	<u>19.5</u>	<u>700</u>		<u>cloudy</u>
<u>855</u>	<u>9</u>	<u>14.55</u>	<u>5.67</u>	<u>0.076</u>	<u>4.35</u>	<u>240</u>	<u>20.3</u>	<u>700</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>RE10702-GW-121815</u>	<u>40-mL vials</u>	<u>3</u>	<u>HCl</u>	<u>VOCs</u>	<u>930</u>
	<u>1-L amber</u>	<u>2</u>	<u>none</u>	<u>1,4-Dioxane</u>	

Comments \_\_\_\_\_

Signature Paul Kavetha Date 12-18-15





Well ID: RE10703

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 12/29/15 Time: Start 1345 am/pm  
 Project No: 60266526 Finish 1600 am/pm  
 Site Location: \_\_\_\_\_  
 Weather Conds: \_\_\_\_\_ Collector(s): \_\_\_\_\_

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

a. Total Well Length 725 c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material 4-inch PVC  
 b. Water Table Depth 43.6 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% - Turbidity ± 10% - D.O. ± 10% (values >0.5 mg/L)  
 - 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

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
14:10	5g	13.47	6.15	0.092	5.41	24.9	142	500	43.6	muddy
14:15		13.47	6.13	0.092	5.64	25.0		500	43.6	muddy
14:20		13.77	6.07	0.078	4.74	16.1	630	500	43.2	"
14:25		13.92	5.93	0.082	5.48	12.2	288	500	43.21	"
14:30		13.86	5.74	0.038	5.73	15.0	210	500	43.21	"
14:35		13.84	5.69	0.031	5.75	15.6	90.1	300	43.20	"

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>RE10703-GW-122915</u>	40-mL vials	3	HCl	VOCs	
	1-L amber	2	none	1,4-Dioxane	

Comments \_\_\_\_\_

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





RESOLUTION  
CONSULTANTS

Well ID: RE108D1

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 12/22/15 Time: Start 845 am/pm  
 Project No: 60266526 Finish 1030 am/pm  
 Site Location: Corona + Cicl  
 Weather Conds: Overcast, sprinkles - 50°F Collector(s): SC

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

a. Total Well Length 555 c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material 4-inch PVC  
 b. Water Table Depth 40.18 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  $\pm 3\%$  - Turbidity  $\pm 10\%$  - D.O.  $\pm 10\%$  (values  $>0.5$  mg/L)  
 - pH  $\pm 0.1$  unit - ORP  $\pm 10$ mV  
 - Sp. Cond.  $\pm 3\%$  - Drawdown  $< 0.3'$  Remove a minimum 1 screen volume

c. Field Testing Equipment used:

Make	Model	Serial Number
Hanna	1+298703	15021X
YSI	556 MFS	073042x

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
0850										Pump on
0900		14.62	5.33	0.078	5.64	259.1	2.29	600	40.24	clear
905		14.64	5.17	0.096	7.54	259.1	1.59	600	40.24	clear
910		14.63	5.12	0.096	7.73	268.4	0.99	600	40.24	clear.
920		14.60	5.11	0.096	7.82	277.8	0.81	600	40.24	"
930		14.58	5.11	0.096	7.91	274.1	0.79	600	40.24	"

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
RE108D1-GW-12222015	40-mL vials	3	HCl	VOCs	100S
RE108D1-GW-12222015	1-L amber	2	none	1,4-Dioxane	

Comments \_\_\_\_\_

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





RESOLUTION CONSULTANTS

Well ID: RE108D2

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 12/22/15 Time: Start 820 am/pm  
 Project No: 60266526 Finish 1030 am/pm  
 Site Location: Corona & Cail  
 Weather Conds: Rain Collector(s): RR JL

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

a. Total Well Length \_\_\_\_\_ c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material 4-inch PVC  
 b. Water Table Depth 40.71 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  $\pm 3\%$  - Turbidity  $\pm 10\%$  - D.O.  $\pm 10\%$  (values  $>0.5$  mg/L)  
 - pH  $\pm 0.1$  unit - ORP  $\pm 10$ mV  
 - Sp. Cond.  $\pm 3\%$  - Drawdown  $< 0.3'$  Remove a minimum 1 screen volume

c. Field Testing Equipment used:

Make	Model	Serial Number
<u>Hanna</u>	<u>HI 98703</u>	<u>V80211X</u>
<u>YSI</u>	<u>556 MPS</u>	<u>54577</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
850		14.45	4.91	0.083	5.15	79.5	0.81	550	40.40	clear
0900		14.44	4.86	0.083	4.12	58.1		550	40.40	
0905		14.45	4.86	0.083	4.10	58.0	0.37	550	40.72	
0910		14.42	4.79	0.081	4.92	50.6	0.33	550		
0920		14.36	4.75	0.081	5.41	44.5	0.25	550	40.72	clear
095		14.35	4.74	0.081	6.36	43.1		550	40.72	clear

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>RE108D2-GW-122215</u>	<u>40-mL vials</u>	<u>3</u>	<u>HCl</u>	<u>VOCs</u>	<u>1020</u>
	<u>1-L amber</u>	<u>2</u>	<u>none</u>	<u>1,4-Dioxane</u>	

Comments \_\_\_\_\_

Signature \_\_\_\_\_

Date \_\_\_\_\_







Well ID: RE11401

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 12/21/15 Time: Start 1100 am/pm  
 Project No: 60266526 Finish 1330 am/pm  
 Site Location: Elm Dr. N  
 Weather Conds: partly sunny 50° Collector(s): \_\_\_\_\_

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

a. Total Well Length 560 c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material 4-inch PVC  
 b. Water Table Depth 31.68 d. Calculated System Volume (see back) 13.1

**2. WELL PURGE DATA**

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

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

c. Field Testing Equipment used: Make VSI Model 556 Serial Number 05H 1965/450231X

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>1130</u>		<u>1</u>								<u>ON</u>
<u>1140</u>		<u>13.95</u>	<u>6.73</u>	<u>0.066</u>	<u>13.36</u>	<u>142.7</u>		<u>350</u>	<u>31.60</u>	<u>cloudy</u>
<u>1145</u>		<u>13.44</u>	<u>6.13</u>	<u>0.068</u>	<u>9.00</u>	<u>142.7</u>			<u>31.60</u>	
<u>1150</u>		<u>13.53</u>	<u>5.94</u>	<u>0.070</u>	<u>6.37</u>	<u>145.5</u>		<del><u>350</u></del>		
<u>1155</u>		<u>13.56</u>	<u>5.92</u>	<u>0.070</u>	<u>6.36</u>	<u>143.4</u>				
<u>1200</u>		<u>13.57</u>	<u>5.90</u>	<u>0.071</u>	<u>4.33</u>	<u>140.7</u>		<u>500</u>	<u>31.62</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
<u>RE10401-GW-12215</u>	<u>40-mL vials</u>	<u>3</u>	<u>HCl</u>	<u>VOCs</u>	<u>1320</u>
	<u>1-L amber</u>	<u>2</u>	<u>none</u>	<u>1,4-Dioxane</u>	

Comments: of tubing  
Bottom depth is way off, at least 40ft of bonded is out of the well.  
thing point is 10ft at the bottom  
\* Tubing is marked "TOP" \* Need to check tubing length and trim to the correct length.

Signature: Paul K... Date: 12/21/15





RESOLUTION  
CONSULTANTS

Well ID: RE11402

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 12/16/15 Time: Start 1300 am/pm  
 Project No: 60266526 Finish 1630 am/pm  
 Site Location: Elm Place  
 Weather Conds: Sunny 60° Collector(s): \_\_\_\_\_

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

a. Total Well Length 635 c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material  
4-inch PVC  
 b. Water Table Depth \_\_\_\_\_ 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  $\pm 3\%$
  - pH  $\pm 0.1$  unit
  - Sp. Cond.  $\pm 3\%$
  - Turbidity  $\pm 10\%$
  - ORP  $\pm 10$ mV
  - Drawdown  $< 0.3'$
  - D.O.  $\pm 10\%$  (values  $> 0.5$  mg/L)
- Remove a minimum 1 screen volume

c. Field Testing Equipment used: \_\_\_\_\_ Make \_\_\_\_\_ Model \_\_\_\_\_ Serial Number \_\_\_\_\_

Time (24hr)	Volume		Temp. (°C)	pH	Spec.			Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
	Removed (Liters)				Cond. (mS/cm)	DO (mg/L)	ORP (mV)				
1345											OK
1355											blow out
1435								400			reset bladder
1445	3 gal		14.30	5.85	0.069	3.29	114.6	44	500	32.03	
1450											
1455			14.13	6.08	0.072	1.94	107.3		475	42.6	

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
<u>RE11402-GW-121615</u>	<u>40-mL vials</u>	<u>3</u>	<u>HCl</u>	<u>VOCs</u>	<u>1615</u>
	<u>1-L amber</u>	<u>2</u>	<u>none</u>	<u>1,4-Dioxane</u>	

Comments: 1355 p. 11 pump reset bladder

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





Well ID: RE11403

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 12/16/15 Time: Start 1300 am/pm  
 Project No: 60266526 Finish 1830 am/pm  
 Site Location: \_\_\_\_\_  
 Weather Conds: Sunny 60° Collector(s): \_\_\_\_\_

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

a. Total Well Length 725 c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material 4-inch PVC  
 b. Water Table Depth 32.35 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%
- pH ± 0.1 unit
- Sp. Cond. ± 3%
- Turbidity ± 10%
- ORP ± 10mV
- Drawdown < 0.3'
- D.O. ± 10% (values >0.5 mg/L)
- Remove a minimum 1 screen volume

c. Field Testing Equipment used:

Make	Model	Serial Number
<u>YSI</u>	<u>556 MPS</u>	<u>05G1942 AC</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>14:20</u>	<u>-</u>	<u>14.41</u>	<u>10.00</u>	<u>0.073</u>	<u>9.65</u>	<u>176.9</u>	<u>-</u>	<u>500</u>	<u>32.27</u>	
<u>14:40</u>	<u>-</u>	<u>14.39</u>	<u>5.93</u>	<u>0.042</u>	<u>5.08</u>	<u>218.0</u>	<u>29.3</u>	<u>500</u>	<u>-</u>	<u>Clear/odor</u>
<u>14:45</u>	<u>-</u>	<u>14.40</u>	<u>5.63</u>	<u>0.035</u>	<u>5.61</u>	<u>224.3</u>	<u>-</u>	<u>500</u>	<u>32.25</u>	
<u>14:50</u>	<u>-</u>	<u>14.37</u>	<u>5.63</u>	<u>0.035</u>	<u>5.62</u>	<u>224.9</u>	<u>-</u>	<u>500</u>	<u>32.26</u>	
<u>14:55</u>	<u>48 Gal</u>	<u>14.37</u>	<u>5.60</u>	<u>0.035</u>	<u>5.68</u>	<u>228.5</u>	<u>2.75</u>	<u>500</u>	<u>32.24</u>	
<u>15:05</u>	<u>-</u>	<u>14.39</u>	<u>5.52</u>	<u>0.034</u>	<u>5.72</u>	<u>233.4</u>	<u>1.82</u>	<u>500</u>	<u>32.27</u>	

d. Acceptance criteria pass/fail (continued on back)

	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 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>RE11403-GW-121615</u>	<u>40-mL vials</u>	<u>3</u>	<u>HCl</u>	<u>VOCs</u>	<u>1800</u>
<u>RE11403-GW-121615</u>	<u>1-L amber</u>	<u>2</u>	<u>none</u>	<u>1,4-Dioxane</u>	

Comments \_\_\_\_\_

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





RESOLUTION  
CONSULTANTS

Well ID: RE12001

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 12/18/15 Time: Start 1320 am/pm  
 Project No: 60266526 Finish 1525 am/pm  
 Site Location: Shelley  
 Weather Conds: cloudy 45° Collector(s): F. V. Avdow

## 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 37.45 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  $\pm 3\%$  - Turbidity  $\pm 10\%$  - D.O.  $\pm 10\%$  (values  $>0.5$  mg/L)  
 - pH  $\pm 0.1$  unit - ORP  $\pm 10$ mV  
 - Sp. Cond.  $\pm 3\%$  - 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>4577</u>
<u>HANNA</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										
1330	1.5	13.74	5.54	0.096	10.12	172	3.13	275	37	clear
1335	1	13.95	5.54	0.095	4.04	153	1.95			
1340	1.5	14.06	5.44	0.095	2.75	145	4.35	300	36.92	Clear
1345	2	14.14	5.3	0.094	2.30	139	5.51			
1350	2.5	14.20	5.23	0.093	2.33	134	2.07	450		

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>RE10701-GW-121815</u>	<u>40-mL vials</u>	<u>3</u>	<u>HCl</u>	<u>VOCs</u>	<u>1525</u>
	<u>1-L amber</u>	<u>2</u>	<u>none</u>	<u>1,4-Dioxane</u>	<u>1525</u>

Comments: Bubbles in flow thru cell affect DO, Turbidity in well

Signature: [Signature] Date: 12/18/15









RESOLUTION  
CONSULTANTS

Well ID: RE 12002

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 12/29/15 Time: Start \_\_\_\_\_ am/pm  
 Project No: 60266526 Finish \_\_\_\_\_ am/pm  
 Site Location: \_\_\_\_\_  
 Weather Conds: hazy Collector(s): JL/TP

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

a. Total Well Length \_\_\_\_\_ c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material  
 b. Water Table Depth 36.6 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% - Turbidity ± 10% - D.O. ± 10% (values >0.5 mg/L)  
 - 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

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
10:25		17.44			5.52	42.4			36.6	Clear
11:30	1094	14.84	5.22	0.080	5.93	42.6		500	36.71	Clear
11:35	-	14.87	5.22	0.080	5.71	42.5		500	36.71	"
11:40	-	14.92	5.22	0.080	5.73	42.2	1.48	500	36.71	"
11:45	-	14.88	5.22	0.080	5.88	42.3	1.47	500	36.71	"
11:50	-	14.87	5.22	0.080	5.91	42.6	1.45	500	36.75	"

d. Acceptance criteria pass/fail

	Yes	No	N/A
Has required volume been removed	<input type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Has required turbidity been reached	<input type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Have parameters stabilized	<input 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>RE12002-GW-122915</u>	40-mL vials	3	HCl	VOCs	
	1-L amber	2	none	1,4-Dioxane	

Comments \_\_\_\_\_

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





RESOLUTION  
CONSULTANTS

Well ID: RE120D3

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 12/29/15 Time: Start \_\_\_\_\_ am/pm  
 Project No: 60266526 Finish \_\_\_\_\_ am/pm  
 Site Location: \_\_\_\_\_  
 Weather Conds: 40s Rainy Collector(s): JC / JP

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

a. Total Well Length \_\_\_\_\_ c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material  
 4-inch PVC  
 b. Water Table Depth 36.9 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  $\pm 3\%$
  - pH  $\pm 0.1$  unit
  - Sp. Cond.  $\pm 3\%$
  - Turbidity  $\pm 10\%$
  - ORP  $\pm 10mV$
  - Drawdown  $< 0.3'$
  - D.O.  $\pm 10\%$  (values  $> 0.5$  mg/L)
- Remove a minimum 1 screen volume

c. Field Testing Equipment used:	Make	Model	Serial Number

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>10 gal</u>								<u>38.9</u>	
<u>12:00</u>	<u>10 gal</u>	<u>14.75</u>	<u>4.86</u>	<u>0.021</u>	<u>11.7</u>	<u>56.1</u>	<u>1.21</u>	<u>500</u>	<u>37.14</u>	<u>clear</u>
<u>12:05</u>	<u>-</u>	<u>14.75</u>	<u>4.88</u>	<u>0.026</u>	<u>4.72</u>	<u>98.1</u>	<u>1.20</u>	<u>500</u>	<u>37.14</u>	<u>-</u>

d. Acceptance criteria pass/fail

Has required volume been removed	Yes <input type="checkbox"/>	No <input type="checkbox"/>	N/A <input type="checkbox"/>
Has required turbidity been reached	<input type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Have parameters stabilized	<input 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-122915</u>	<u>40-mL vials</u>	<u>3</u>	<u>HCl</u>	<u>VOCs</u>	
	<u>1-L amber</u>	<u>2</u>	<u>none</u>	<u>1,4-Dioxane</u>	

Comments \_\_\_\_\_

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





RESOLUTION  
CONSULTANTS

Well ID: RE 121 01

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 12/21/15 Time: Start 830 am/pm  
 Project No: 60266526 Finish 13 am/pm  
 Site Location: \_\_\_\_\_  
 Weather Conds: \_\_\_\_\_ Collector(s): \_\_\_\_\_

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

a. Total Well Length \_\_\_\_\_ c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material  
4-inch PVC  
 b. Water Table Depth \_\_\_\_\_ 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%
  - pH ± 0.1 unit
  - Sp. Cond. ± 3%
  - Turbidity ± 10%
  - ORP ± 10mV
  - Drawdown < 0.3'
  - D.O. ± 10% (values >0.5 mg/L)
- Remove a minimum 1 screen volume

c. Field Testing Equipment used:	Make	Model	Serial Number

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
855		14.92	10.02	0.136	2.40	1651			34.40	
900		14.90	6.31	0.080	1.48	146.3	824			
905		14.88	6.25	0.085	1.38	146.6			34.72	
910		14.91	6.08	0.075	1.30	146.0	455			
915		14.97	5.58	0.072	1.10	144.9			35.40	
920		14.95	5.55	0.071	0.95	150.9	10.9			

d. Acceptance criteria pass/fail

	Yes	No	N/A	(continued on back)
Has required volume been removed	<input type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Has required turbidity been reached	<input type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Have parameters stabilized	<input 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 vials	3	HCl	VOCs	
	1-L amber	2	none	1,4-Dioxane	

Comments \_\_\_\_\_

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





RESOLUTION  
CONSULTANTS

Well ID: RE 121 D2

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 12/21/15 Time: Start 1115 am/pm  
 Project No: 60266526 Finish 1340 am/pm  
 Site Location: Veely @ Union  
 Weather Conds: Cloudy 47°F Collector(s): JC

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

a. Total Well Length 575 c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material  
4-inch PVC  
 b. Water Table Depth 34.84 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%
  - pH ± 0.1 unit
  - Sp. Cond. ± 3%
  - Turbidity ± 10%
  - ORP ± 10mV
  - Drawdown < 0.3'
  - D.O. ± 10% (values >0.5 mg/L)
- Remove a minimum 1 screen volume

#### c. Field Testing Equipment used:

Make	Model	Serial Number
<u>Hanna</u>	<u>HI98703</u>	<u>U80211X</u>
<u>YSI</u>	<u>556 mPS</u>	<u>54577</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>1120</u>									<u>34.39</u>	
<u>1150</u>		<u>15.03</u>	<u>5.29</u>	<u>0.049</u>	<u>7.85</u>	<u>6.7</u>	<u>369</u>	<u>425</u>	<u>34.39</u>	<u>Cloudy</u>
<u>1155</u>		<u>15.05</u>	<u>5.27</u>	<u>0.049</u>	<u>6.63</u>	<u>-1.8</u>	<u>357</u>	<u>425</u>	<u>34.39</u>	<u>"</u>
<u>1200</u>		<u>15.11</u>	<u>5.24</u>	<u>0.063</u>	<u>5.29</u>	<u>-11.0</u>	<u>366</u>	<u>425</u>	<u>34.39</u>	<u>"</u>
<u>1205</u>		<u>15.12</u>	<u>5.25</u>	<u>0.063</u>	<u>5.13</u>	<u>-13.2</u>	<u>362</u>	<u>450</u>	<u>34.40</u>	<u>"</u>
<u>1210</u>		<u>15.12</u>	<u>5.27</u>	<u>0.068</u>	<u>4.40</u>	<u>-22.4</u>	<u>106</u>	<u>450</u>	<u>34.40</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"/>

(continued on back)

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>RE121D2-GW-12212015</u>	<u>40-mL vials</u>	<u>3</u>	<u>HCl</u>	<u>VOCs</u>	<u>1331</u>
<u>RE121D2-GW-12212015</u>	<u>1-L amber</u>	<u>2</u>	<u>none</u>	<u>1,4-Dioxane</u>	<u>1331</u>

Comments \_\_\_\_\_

Signature \_\_\_\_\_

Date 12/21/2015







Well ID: RE12201

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 12/15/15 Time: Start 1720 am/pm  
 Project No: 60266526 Finish 1600 am/pm  
 Site Location: Hayden & Curtis  
 Weather Conds: Sunny 60° Collector(s): DL

### 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 42.79 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% - Turbidity ± 10% - D.O. ± 10% (values >0.5 mg/L)  
 - 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 \_\_\_\_\_

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										Pump on
1340		15.11	6.06	0.080	9.57	112.6	0.57	1.00	43.70	Clear
1345		15.20	5.63	0.077	4.91	103.1		600	43.71	clear.
1350		15.12	5.63	0.078	4.51	100.7				
1355		15.08	5.64	0.078	4.27	107.8				
1400		15.09	5.69	0.080	3.87	95.6			42.61	

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
<u>RE12201-GW-12/15/16</u>	40-mL vials	3	HCl	VOCs	1455
	1-L amber	2	none	1,4-Dioxane	1455

Comments \_\_\_\_\_

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





Well ID: RE12202

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 12/15/15 Time: Start 1220 am/pm  
 Project No: 60266526 Finish 1600 am/pm  
 Site Location: Haydon & Curtis  
 Weather Conds: sunny 60° 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 43.03 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% - Turbidity ± 10% - D.O. ± 10% (values >0.5 mg/L)  
 - 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>173042X</u>
<u>Hanna</u>	<u>#1 98703</u>	<u>1180211 X</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>1325</u>										<u>0N</u>
<u>1330</u>								<u>600</u>		
<u>1335</u>		<u>15.46</u>	<u>5.37</u>	<u>0.090</u>	<u>5.82</u>	<u>165.2</u>			<u>43.02</u>	
<u>1340</u>		<u>15.39</u>	<u>5.31</u>	<u>0.072</u>	<u>4.47</u>	<u>164.9</u>	<u>4.15</u>			
<u>1345</u>		<u>15.22</u>	<u>5.28</u>	<u>0.071</u>	<u>4.46</u>	<u>166.9</u>				
<u>1350</u>		<u>15.32</u>	<u>5.22</u>	<u>0.071</u>	<u>4.93</u>	<u>170.7</u>				

d. Acceptance criteria pass/fail

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

If no or N/A - Explain below.

### 3. SAMPLE COLLECTION:

Method: Geotech bladder pump with drop tube assembly

Sample ID	Container Type	No. of Containers	Preservation	Analysis Req.	Time
<u>RE12202-GW-121515</u>	<u>40-mL vials</u>	<u>3</u>	<u>HCl</u>	<u>VOCs</u>	<u>1510</u>
	<u>1-L amber</u>	<u>2</u>	<u>none</u>	<u>1,4-Dioxane</u>	

Comments \_\_\_\_\_

Signature Paul Kautz Date 12/15/15





Well ID: RE12203

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 12/15/15 Time: Start 1225 am/pm  
 Project No: 60266526 Finish 1600 am/pm  
 Site Location: Hayden & Curtis  
 Weather Conds: Sunny 60° Collector(s): \_\_\_\_\_

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

a. Total Well Length 240 c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material 4-inch PVC  
 b. Water Table Depth 43.36 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%
  - pH ± 0.1 unit
  - Sp. Cond. ± 3%
  - Turbidity ± 10%
  - ORP ± 10mV
  - Drawdown < 0.3'
  - D.O. ± 10% (values >0.5 mg/L)
- Remove a minimum 1 screen volume

c. Field Testing Equipment used:

Make	Model	Serial Number
<u>YSI</u>	<u>556 MPS</u>	<u>05H1965AR</u>
<u>Hanna</u>	<u>HI9870J</u>	<u>F0018653</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>13:40</u>	<u>0</u>	<u>15.15</u>	<u>4.95</u>	<u>0.024</u>	<u>3.74</u>	<u>174.2</u>	<u>86.1</u>	<u>500</u>	<u>42.90</u>	<u>Cloudy</u>
<u>13:45</u>	<u>-</u>	<u>14.90</u>	<u>4.87</u>	<u>0.023</u>	<u>3.60</u>	<u>178.7</u>	<u>-</u>	<u>-</u>	<u>-</u>	<u>Cloudy</u>
<u>13:50</u>		<u>14.91</u>	<u>4.68</u>	<u>0.021</u>	<u>3.59</u>	<u>176.5</u>				
<u>13:55</u>		<u>14.90</u>	<u>4.63</u>	<u>0.020</u>	<u>3.60</u>	<u>177.8</u>				
<u>14:00</u>	<u>5 gal</u>	<u>14.91</u>	<u>4.66</u>	<u>0.020</u>	<u>3.52</u>	<u>176.4</u>				
<u>14:05</u>		<u>14.96</u>	<u>4.80</u>	<u>0.020</u>	<u>3.32</u>	<u>171.7</u>			<u>42.85</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
<u>RE12203-GW-121515</u>	<u>40-mL vials</u>	<u>3</u>	<u>HCl</u>	<u>VOCs</u>	<u>1500</u>
<u>LE12203-GW-121515</u>	<u>1-L amber</u>	<u>2</u>	<u>none</u>	<u>1,4-Dioxane</u>	

Comments \_\_\_\_\_

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





RESOLUTION  
CONSULTANTS

Well ID: RE12301

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 12/2/15 Time: Start 7:30 am/pm  
 Project No: 60266526 Finish 10:30 am/pm  
 Site Location: LIRK Lot  
 Weather Conds: Sunny 40° Collector(s): \_\_\_\_\_

### 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 47.65 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%
  - pH ± 0.1 unit
  - Sp. Cond. ± 3%
  - Turbidity ± 10%
  - ORP ± 10mV
  - Drawdown < 0.3'
  - D.O. ± 10% (values >0.5 mg/L)
- Remove a minimum 1 screen volume

c. Field Testing Equipment used: Make \_\_\_\_\_ Model \_\_\_\_\_ Serial Number \_\_\_\_\_

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
8:30										ON
8:40		15.08	6.64	0.087	21.35	89.2		550	47.58	
8:45		15.20	5.91	0.087	11.32	102.1	15.4	550	47.54	clear
8:50		15.29	5.67	0.087	10.27	110.4				
9:00	5gal	15.38	5.58	0.087	10.24	113.3				
9:05		15.45	5.41	0.087	9.94	121.7		700		

d. Acceptance criteria pass/fail (continued on back)

	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>RE12301-GW-122115</u>	<u>40-mL vials</u>	<u>3</u>	<u>HCl</u>	<u>VOCs</u>	
	<u>1-L amber</u>	<u>2</u>	<u>none</u>	<u>1,4-Dioxane</u>	

Comments \_\_\_\_\_

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









Well ID: RL12302

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 12/21/15 Time: Start 730 am/pm  
 Project No: 60266526 Finish 1030 am/pm  
 Site Location: LIRR 600  
 Weather Conds: sunny 40 Collector(s): E. ACS

**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 48.90 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%
  - pH ± 0.1 unit
  - Sp. Cond. ± 3%
  - Turbidity ± 10%
  - ORP ± 10mV
  - Drawdown < 0.3'
  - D.O. ± 10% (values >0.5 mg/L)
- Remove a minimum 1 screen volume

c. Field Testing Equipment used:

Make	Model	Serial Number
<u>YSI</u>	<u>556</u>	<u>0561942</u>
<u>HANNA</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>810</u>								<u>550</u>		<u>OK</u>
<u>830</u>		<u>13.81</u>	<u>5.93</u>	<u>0.030</u>	<u>6.00</u>	<u>162.1</u>		<u>550</u>	<u>48.98</u>	<u>clear / no odor</u>
<u>835</u>		<u>13.87</u>	<u>5.77</u>	<u>0.025</u>	<u>6.52</u>	<u>167.8</u>				
<u>840</u>		<u>13.82</u>	<u>5.69</u>	<u>0.024</u>	<u>7.32</u>	<u>170.7</u>				<u>clear / no odor</u>
<u>845</u>		<u>13.97</u>	<u>5.54</u>	<u>0.022</u>	<u>8.26</u>	<u>179.2</u>	<u>4.22</u>	<u>550</u>	<u>48.98</u>	<u>" "</u>
<u>850</u>	<u>56L</u>	<u>13.82</u>	<u>5.54</u>	<u>0.022</u>	<u>8.31</u>	<u>183.8</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>RL123 D2-GW-122115</u>	<u>40-mL vials</u>	<u>3</u>	<u>HCl</u>	<u>VOCs</u>	<u>9:55</u>
	<u>1-L amber</u>	<u>2</u>	<u>none</u>	<u>1,4-Dioxane</u>	<u>9:55</u>

Comments \_\_\_\_\_

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





RESOLUTION  
CONSULTANTS

Well ID: RE12303

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 12/21/15 Time: Start 7:30 am/pm  
 Project No: 60266526 Finish \_\_\_\_\_ am/pm  
 Site Location: LIRR Lot  
 Weather Conds: Sunny 40° Collector(s): \_\_\_\_\_

### 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 48.74 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  $\pm 3\%$  - Turbidity  $\pm 10\%$  - D.O.  $\pm 10\%$  (values  $>0.5$  mg/L)  
 - pH  $\pm 0.1$  unit - ORP  $\pm 10$ mV  
 - Sp. Cond.  $\pm 3\%$  - Drawdown  $< 0.3'$  Remove a minimum 1 screen volume  
 c. Field Testing Equipment used:

Make	Model	Serial Number

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
8:20		10.59	7.33	0.067	11.47	97.1	<del>550</del>	550	48.74	Clear
8:25		14.30	5.82	0.029	2.64	39.1	<del>550</del>	550	48.95	mucky.
8:30		14.33	4.98	0.028	1.44	6.4	-	-		
8:35		14.37	4.90	0.029	1.23	-4.2	-	-		
8:40		14.38	5.01	0.035	1.26	-9.1	-	-	48.95	
8:45	5 gal	14.43	5.68	0.069	1.19	-24.3	-	600		

d. Acceptance criteria pass/fail

	Yes	No	N/A
Has required volume been removed	<input type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Has required turbidity been reached	<input type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Have parameters stabilized	<input 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>RE12303-6W-122115</u>	<u>40-mL vials</u>	<u>3</u>	<u>HCl</u>	<u>VOCs</u>	<u>9:45</u>
	<u>1-L amber</u>	<u>2</u>	<u>none</u>	<u>1,4-Dioxane</u>	

Comments: Sample time - 9:45

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





RESOLUTION CONSULTANTS

Well ID: TT1010

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 12/17/15 Time: Start 1045 am/pm  
 Project No: 60266526 Finish 1145 am/pm  
 Site Location: Emerson  
 Weather Conds: Overcast, drizzle 50° 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 4-inch PVC  
 b. Water Table Depth 33.85 d. Calculated System Volume (see back) 13.1

### 2. WELL PURGE DATA

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

#### b. Acceptance Criteria defined (see workplan)

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

#### c. Field Testing Equipment used:

Make	Model	Serial Number

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
10:50		15.39	4.50	0.081	0.73	266.4		850		
10:55		15.36	4.67	0.078	0.66	239.6		850	34.05	
11:00		15.33	4.66	0.078	0.61	236.9		850		
11:05		15.30	4.66	0.078	0.58	231.9		850	34.00	
11:10		15.30	4.66	0.078	0.57	231.5		850	33.98	
11:15		15.31	4.66	0.078	0.56	230.7		875	34.00	Started Reading

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>TT1010-GW-121715</u>	<u>40-mL vials</u>	<u>3</u>	<u>HCl</u>	<u>VOCs</u>	<u>1145</u>
<u>DUPLICATE 2 - GW-121715</u>	<u>1-L amber</u>	<u>2</u>	<u>none</u>	<u>1,4-Dioxane</u>	
		<u>5</u>		<u>VOCs, 1,4-Dioxane</u>	<u>1155</u>

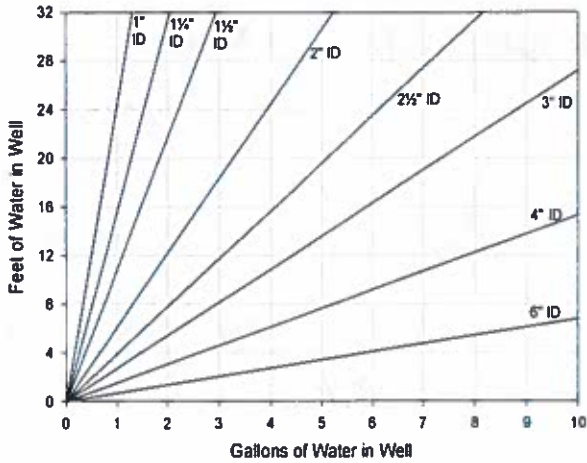
Comments: Duplicate

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

One screen volume  
(4-inch well)

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

Well ID:

(continued from front)

Time (24 hr)	Volume Removed (Liters)	Temp (°C)	pH	Specific Cond. (mS/cm)	DO (mg/L)	ORP (mV)	Turbidity (NTU)	Flow Rate (ml/min)	Depth to water (ft)	Color/Odor
11:20		15.28	4.65	0.078	0.51	225.4	0.28	850		Clear
11:25		15.28	4.65	0.078	0.50	225.2				
11:30	10 Gal	15.28	4.65	0.078	0.50	224.9		850	34.00	Clear
11:35		15.28	4.64	0.078	0.49	224.8	0.02			
11:40		15.27	4.63	0.078	0.47	224.2	0.08	850		
11:45	13.5 Gal	15.26	4.63	0.078	0.47	225.0	0.10	850	34.00	Clear
	Sample 11:45									
	Duplicate 11:55									
	Note: Grouting around casing subsided about 2.5 ft recommend more sand <del>1.0</del> 2.0 / sand									



Well ID: TT10101

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 12/17/15 Time: Start 1145 am/pm  
 Project No: 60266526 Finish 1330 am/pm  
 Site Location: Emerson  
 Weather Conds: Overcast drizzle, 50° Collector(s): \_\_\_\_\_

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

a. Total Well Length 595 c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material 4-inch PVC  
 b. Water Table Depth 35.00 d. Calculated System Volume (see back) 13.1

### 2. WELL PURGE DATA

a. Purge Method: Geotech bladder pump with drop tube assembly dedicated  
 b. Acceptance Criteria defined (see workplan)  
 - Temperature ± 3% - Turbidity ± 10% - D.O. ± 10% (values >0.5 mg/L)  
 - 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

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
12:00	-	15.37	5.10	0.081	0.60	239.4		850		Bleeding Clear
12:05		15.36	5.09	0.081	0.61	239.3		850		
12:10		15.36	5.08	0.081	0.59	239.3				
12:15		15.30	5.02	0.082	0.92	238.1		850		
12:20	5 Gal	15.29	5.02	0.082	0.93	238.0				
12:25		15.30	5.03	0.082	0.91	237.4	0.21	850	35.21	Clear

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
<u>TT 10101-GW-121715</u>	40-mL vials	3	HCl	VOCs	1300
	1-L amber	2	none	1,4-Dioxane	

Comments \_\_\_\_\_

Signature \_\_\_\_\_

Date \_\_\_\_\_







RESOLUTION  
CONSULTANTS

Well ID: TT 10102

# Low Flow Ground Water Sample Collection Record

Client: Navy NWIRP Bethpage Date: 12/21/15 Time: Start 1345 am/pm  
 Project No: 60266526 Finish 1630 am/pm  
 Site Location: Emerson  
 Weather Conds: \_\_\_\_\_ Collector(s): Paul Kurek

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

a. Total Well Length 765 c. Length of Water Column \_\_\_\_\_ (a-b) Casing Diameter/Material  
 4-inch PVC  
 b. Water Table Depth 3538 d. Calculated System Volume (see back) 13.1

### 2. WELL PURGE DATA

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

b. Acceptance Criteria defined (see workplan)

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

c. Field Testing Equipment used:

Make	Model	Serial Number
<u>YST</u>	<u>556</u>	<u>0541965 050231X</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>1355</u>										<u>0.4</u>
<u>1400</u>		<u>15.42</u>	<u>5.61</u>	<u>0.033</u>	<u>4.10</u>	<u>152.6</u>	<u>0.62</u>	<u>700</u>	<u>35.38</u>	
<u>1405</u>		<u>15.36</u>	<u>5.54</u>	<u>0.033</u>	<u>3.38</u>	<u>145.7</u>	<u>0.00</u>			
<u>1410</u>		<u>15.34</u>	<u>5.41</u>	<u>0.033</u>	<u>4.04</u>	<u>142.9</u>			<u>35.38</u>	
<u>1415</u>		<u>15.34</u>	<u>5.31</u>	<u>0.034</u>	<u>4.44</u>	<u>142.4</u>	<u>0.00</u>	<u>700</u>		
<u>1420</u>	<u>5.6AL</u>	<u>15.32</u>	<u>5.25</u>	<u>0.034</u>	<u>5.14</u>	<u>146.0</u>	<u>0.00</u>	<u>700</u>	<u>35.38</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>TT 10102-GW-12215</u>	<u>40-mL vials</u>	<u>3</u>	<u>HCl</u>	<u>VOCs</u>	<u>1510</u>
	<u>1-L amber</u>	<u>2</u>	<u>none</u>	<u>1,4-Dioxane</u>	

DUPLICATE-GW-12215 1600

Comments: Duplicate sample

Signature: Paul Kurek Date: 12/21/15



## Appendix B

### Analytical Data Validation – Resolution Consultants



DATA VALIDATION REPORT

Project:	Regional Groundwater Investigation — NWIRP Bethpage	
Laboratory:	Katahdin Analytical	
Sample Delivery Group:	BETHPAGE-3	
Analyses/Method:	Volatile Organic Compounds by U.S. EPA SW-846 Method 8260C 1,4-Dioxane by U.S. EPA SW-846 Method 8270D via Selective Ion Monitoring (SIM)	
Validation Level:	3	
Project Number:	0888812477.SA.DV	
Prepared by:	Dana Miller/Resolution Consultants	Completed on: 02/5/2016
Reviewed by:	Tina Clemmey/Resolution Consultants	File Name: BETHPAGE 3_8260C_8270D

SUMMARY

This report summarizes data review findings for samples listed below, collected by Resolution Consultants from the Regional Groundwater Investigation — NWIRP Bethpage Site on 14 to 29 December 2015 in accordance with the following Sampling and Analysis Plans:

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

Sample ID	Matrix/Sample Type	Analysis
FIELD1-FB-121615	Field Blank	8260C/8270D_SIM
RE103D1-GW-121415	Groundwater	8260C/8270D_SIM
RE103D2-GW-121415	Groundwater	8260C/8270D_SIM
RE103D3-GW-121415	Groundwater	8260C/8270D_SIM
RE104D1-GW-121515	Groundwater	8260C/8270D_SIM
RE104D2-GW-121515	Groundwater	8260C/8270D_SIM
DUPLICATE1-GW-121515	Field Duplicate of RE104D2-GW-121515	8260C/8270D_SIM
RE104D3-GW-121515	Groundwater	8260C/8270D_SIM
RE105D1-GW-121715	Groundwater	8260C/8270D_SIM
RE105D2-GW-121715	Groundwater	8260C/8270D_SIM
RE107D1-GW-121815	Groundwater	8260C/8270D_SIM

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

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

## REVIEW ELEMENTS

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

- ✓ Data completeness (chain-of-custody)/sample integrity
- ✓ Holding times and sample preservation
- ✓ Gas chromatography/Mass spectrometer performance checks
- ✗ Initial calibration verification (ICV)/continuing calibration verification (CCV)
- ✓ Laboratory blanks/trip blanks/field blanks
- ✗ Surrogate spike recoveries
- ✗ 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 (✗) 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 Tables A-1, A-2, and A-3.

Surrogate Spike Recoveries

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

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

Surrogate Recovery Non-conformance Chart:

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

Notes:

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

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

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

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

MS/MSD Non-conformances Chart:

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

Notes:

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

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



### Field Duplicate

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

### Field Duplicate Non-conformances Chart:

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

*Notes:*

LOQ = Limit of quantitation  
 J = Estimated  
 UJ = Undetected and estimated

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

### Qualifications Actions

The data were reviewed independently from the laboratory to assess data quality. All compounds detected at concentrations less than the limit of quantitation but greater than the method detection limit were qualified by the laboratory as estimated (J). This "J" qualifier was retained during data validation. Any sample that was analyzed at a dilution because of high concentrations of target or non-target analytes was checked to confirm that the results and/or sample-specific limit of quantitation and limit of detections were adjusted accordingly by the laboratory. Trichloroethene in sample RE108D2-GW-121215 result value reported above the calibration range and was qualified estimated "J" because the value was off-scale.

No results were rejected; therefore, analytical completeness was calculated to be 100 percent. Data not qualified during data review are considered usable by the project. The remaining results qualified as estimated may be high or low, but the data are usable for their intended purpose,

according to U.S. EPA and Department of Defense guidelines. Final data review qualifiers used to describe results and how they should be interpreted by the end data user are provided in Attachment B and Attachment C. Attachment D provides final results after data review.

#### ATTACHMENTS

Attachment A: Non-Conformance Summary Tables

Attachment B: Qualifier Codes and Explanations

Attachment C: Reason Codes and Explanations

Attachment D: Final Results after Data Review

Attachment A  
Non-Conformance Summary Table

Table A-1 Initial Calibration Non-Conformance					
Method	Analyte	%RSD	Limit	Associated Samples	Qualifier
8260C	CHLOROETHANE	17.80841	<15%	TI0330-1, -2, -5, -6, -11, -4, -7, -3DL, -10RA, -12, -13, -14, -17, -4DL, -6DL, and 7DL	Detects: J Non-detects: UJ
8260C	TETRACHLOROETHENE	15.13611	<15%	TI0428-4, -1, -3RA, -1DL, -2RA, and TI0330-17DL	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	TETRACHLOROETHENE	P3840.D	150.8	80-120	TI0116-1, -2, -3, -4, -5, -6, -7, -8, -9, -11, -10RA, -2DL, -4DL, -9DL, -3DL, -11DL, TI0214-1, -4, -5, -6, -7, -9, -8, -10, 12, -8DL, and -7DL	Detects: J Non-detects: UJ
8260C	TRANS-1,3-DICHLOROPROPENE	C6396A.D	123.47	80-120	TI0330-1, -2, -5, -6, -11, -4, -7, -3DL, -10RA, -12, -13, -14, -17, -4DL, -6DL, and 7DL	Detects: J Non-detects: UJ
8260C	ACETONE	P4114A.D	72.62	80-120	TI0428-3RA, -1DL, -2RA, -4, -1, and TI0330-17DL	Detects: J Non-detects: UJ

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

*Notes:*

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

Table A-4  
Surrogate Non-Conformance

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

*Notes:*

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

Table A-5  
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-122115	METHYLENE CHLORIDE	<2.5	50.0	53.4	60.2	55-140	UJ
TT101D2-GW-122115	CIS-1,2-DICHLOROETHENE	1.7	50.0	54.6	59.8	70-125	J
TT101D2-GW-122115	1,2,4-TRICHLOROBENZENE	<0.50	50.0	50.8	56.2	65-135	UJ
TT101D2-GW-122115	CHLOROBENZENE	<0.50	50.0	55.8	60.2	80-120	UJ
TT101D2-GW-122115	1,1-DICHLOROETHANE	<0.50	50.0	64.4	70.8	70-135	UJ
TT101D2-GW-122115	CIS-1,3-DICHLOROPROPENE	<0.50	50.0	60.8	64.4	70-130	UJ
TT101D2-GW-122115	1,2-DIBROMO-3-CHLOROPROPANE	<0.75	50.0	46.4	57	50-130	UJ
TT101D2-GW-122115	ISOPROPYL BENZENE	<0.50	50.0	56.2	61	75-125	UJ
TT101D2-GW-122115	TRANS-1,2-DICHLOROETHENE	<0.50	50.0	58.4	65.2	60-140	UJ
TT101D2-GW-122115	BENZENE	<0.50	50.0	62.4	67.6	80-120	UJ
TT101D2-GW-122115	1,2-DICHLOROPROPANE	<0.50	50.0	63.8	68.8	75-125	UJ
TT101D2-GW-122115	O-XYLENE	<0.50	50.0	58.8	62.2	80-120	UJ
TT101D2-GW-122115	1,3-DICHLOROBENZENE	<0.50	50.0	51.4	56.8	75-125	UJ
TT101D2-GW-122115	1,1-DICHLOROETHENE	3.6	50.0	56.8	63.4	70-130	J
TT101D2-GW-122115	1,1,2-TRICHLOROETHANE	0.50	50.0	63.4	65.4	75-125	J
TT101D2-GW-122115	CYCLOHEXANE	<0.50	50.0	63.8	69	71-133	UJ
TT101D2-GW-122115	TOLUENE	<0.50	50.0	65	69.6	75-120	UJ
TT101D2-GW-122115	CARBON TETRACHLORIDE	1.3	50.0	63.6	66.6	65-140	J
TT101D2-GW-122115	1,2-DICHLOROETHANE	<0.50	50.0	61.6	66	70-130	UJ
TT101D2-GW-122115	1,2-DICHLOROETHENE, TOTAL	1.7	100	56.5	62.5	84-121	J
TT101D2-GW-122115	XYLENES, TOTAL	<1.5	150	59.8	63.2	89-116	UJ
TT101D2-GW-122115	STYRENE	<0.50	50.0	60.4	63.4	65-135	UJ
TT101D2-GW-122115	DIBROMOCHLOROMETHANE	<0.50	50.0	59.4	63.6	60-135	UJ
TT101D2-GW-122115	1,4-DICHLOROBENZENE	<0.50	50.0	50.8	54.6	75-125	UJ
TT101D2-GW-122115	ETHYLBENZENE	<0.50	50.0	56.2	60.8	75-125	UJ
TT101D2-GW-122115	M- AND P-XYLENE	<1.0	100	60.2	63.8	75-130	UJ
TT101D2-GW-122115	BROMOFORM	<0.50	50.0	55.6	57	70-130	UJ
TT101D2-GW-122115	BROMODICHLOROMETHANE	<0.50	50.0	67.6	70	75-120	UJ
TT101D2-GW-122115	CHLOROFORM	0.90	50.0	57	62.4	65-135	J
TT101D2-GW-122115	1,1,1-TRICHLOROETHANE	0.34	50.0	58.9	64.3	65-130	J
TT101D2-GW-122115	1,2-DIBROMOETHANE	<0.50	50.0	61.2	66.6	80-120	UJ
TT101D2-GW-122115	1,1,2,2-TETRACHLOROETHANE	<0.50	50.0	51.6	59.2	65-130	UJ
TT101D2-GW-122115	1,2-DICHLOROBENZENE	<0.50	50.0	52	57.6	70-120	UJ

*Notes:*

- MS = Matrix spike
- MSD = Matrix spike duplicate
- %R = Percent recovery
- Bold** = Percent recovery not within control limit
- UJ = Nondetect analyte in associated sample qualified estimated "UJ" because the %R is lower than the control limit.
- J = Detected analyte in associated sample qualified estimated "J" because %R is lower than the control limit.



Table A-6  
Field Duplicate  
(Micrograms per liter)

Sample ID	Duplicate ID	Analyte	Sample Result	Duplicate Result	RPD	Qualifiers
TT101D2-GW-122115	DUPLICATE-GW-122115	1,1-DICHLOROETHENE	3.6	5	32.6	J - both results

*Notes:*

RPD = Relative percent difference  
 J = Estimated value

Attachment B  
Qualifier Codes and Explanations

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

Attachment C  
Reason Codes and Explanations

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

Attachment D  
Final Results after Data Review

Sample Delivery Group				BETHPAGE-3		
Lab ID				TI0116-1		
Sample ID				TRIP BLANK 121415		
Sample Date				12/14/2015		
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	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	UJ	c
8260C	CARBON DISULFIDE	75-15-0	UG_L	0.5	U	
8260C	CARBON TETRACHLORIDE	56-23-5	UG_L	0.5	U	
8260C	CHLOROETHANE	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	U	
8260C	ETHYLBENZENE	100-41-4	UG_L	0.5	U	
8260C	ISOPROPYLBENZENE	98-82-8	UG_L	0.5	U	
8260C	M- AND P-XYLENE	108-38-3/106-42	UG_L	1	U	
8260C	METHYL ACETATE	79-20-9	UG_L	0.75	U	
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	U	
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	U	
8260C	METHYLENE CHLORIDE	75-09-2	UG_L	2.5	U	
8260C	O-XYLENE	95-47-6	UG_L	0.5	U	
8260C	STYRENE	100-42-5	UG_L	0.5	U	
8260C	TETRACHLOROETHENE	127-18-4	UG_L	0.5	UJ	c
8260C	TOLUENE	108-88-3	UG_L	0.33	J	
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	U	
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	U	
8260C	TRICHLOROETHENE	79-01-6	UG_L	0.5	U	
8260C	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-3		
				Lab ID	TI0116-10RA		
				Sample ID	RE122D3-GW-121515		
				Sample Date	12/15/2015		
				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	U		
8260C	2-HEXANONE	591-78-6	UG L	2.5	U		
8260C	4-METHYL-2-PENTANONE	108-10-1	UG L	2.5	UJ		c
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	UJ		c
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	U		
8260C	ETHYLBENZENE	100-41-4	UG L	0.5	U		
8260C	ISOPROPYLBENZENE	98-82-8	UG L	0.5	U		
8260C	M- AND P-XYLENE	108-38-3/106-42	UG L	1	U		
8260C	METHYL ACETATE	79-20-9	UG L	0.75	U		
8260C	METHYL CYCLOHEXANE	108-87-2	UG L	0.5	U		
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG L	0.5	U		
8260C	METHYLENE CHLORIDE	75-09-2	UG L	2.5	U		
8260C	O-XYLENE	95-47-6	UG L	0.5	U		
8260C	STYRENE	100-42-5	UG L	0.5	U		
8260C	TETRACHLOROETHENE	127-18-4	UG L	0.5	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	2.5			
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-3		
				Lab ID	TI0116-11		
				Sample ID	RE122D2-GW-121515		
				Sample Date	12/15/2015		
				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	3.1			
8260C	1,1-DICHLOROETHANE	75-34-3	UG L	1.5			
8260C	1,1-DICHLOROETHENE	75-35-4	UG L	8.9			
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG L	0.5	U		
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG L	0.75	U		
8260C	1,2-DIBROMOETHANE	106-93-4	UG L	0.5	U		
8260C	1,2-DICHLOROBENZENE	95-50-1	UG L	0.5	U		
8260C	1,2-DICHLOROETHANE	107-06-2	UG L	0.5	U		
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG L	5.7			
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	UJ		c
8260C	CARBON DISULFIDE	75-15-0	UG L	0.5	U		
8260C	CARBON TETRACHLORIDE	56-23-5	UG L	1.9			
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	2.6			
8260C	CHLOROMETHANE	74-87-3	UG L	1	U		
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG L	5.7			
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	U		
8260C	ETHYLBENZENE	100-41-4	UG L	0.5	U		
8260C	ISOPROPYLBENZENE	98-82-8	UG L	0.5	U		
8260C	M- AND P-XYLENE	108-38-3/106-42	UG L	1	U		
8260C	METHYL ACETATE	79-20-9	UG L	0.75	U		
8260C	METHYL CYCLOHEXANE	108-87-2	UG L	0.5	U		
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG L	0.5	U		
8260C	METHYLENE CHLORIDE	75-09-2	UG L	2.5	U		
8260C	O-XYLENE	95-47-6	UG L	0.5	U		
8260C	STYRENE	100-42-5	UG L	0.5	U		
8260C	TETRACHLOROETHENE	127-18-4	UG L	2.3	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	4700			
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	11			

				Sample Delivery Group		BETHPAGE-3	
				Lab ID	TI0116-2		
				Sample ID	RE103D3-GW-121415		
				Sample Date	12/14/2015		
				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	2.5			
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG L	0.5	U		
8260C	1,1-DICHLOROETHANE	75-34-3	UG L	0.5	U		
8260C	1,1-DICHLOROETHENE	75-35-4	UG L	0.62	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	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	UJ	c	
8260C	CARBON DISULFIDE	75-15-0	UG L	0.5	U		
8260C	CARBON TETRACHLORIDE	56-23-5	UG L	0.24	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	0.79	J		
8260C	CHLOROMETHANE	74-87-3	UG L	1	U		
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG L	1			
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG L	0.5	U		
8260C	CYCLOHEXANE	110-82-7	UG L	0.5	U		
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG L	0.5	U		
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG L	1	U		
8260C	ETHYLBENZENE	100-41-4	UG L	0.5	U		
8260C	ISOPROPYLBENZENE	98-82-8	UG L	0.5	U		
8260C	M- AND P-XYLENE	108-38-3/106-42	UG L	1	U		
8260C	METHYL ACETATE	79-20-9	UG L	0.75	U		
8260C	METHYL CYCLOHEXANE	108-87-2	UG L	0.5	U		
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG L	0.5	U		
8260C	METHYLENE CHLORIDE	75-09-2	UG L	2.5	U		
8260C	O-XYLENE	95-47-6	UG L	0.5	U		
8260C	STYRENE	100-42-5	UG L	0.5	U		
8260C	TETRACHLOROETHENE	127-18-4	UG L	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	510			
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	0.81			

				Sample Delivery Group		BETHPAGE-3	
				Lab ID	TI0116-3		
				Sample ID	RE103D1-GW-121415		
				Sample Date	12/14/2015		
				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	0.62	J		
8260C	1,1-DICHLOROETHANE	75-34-3	UG L	1.1			
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	3.2			
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	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	UJ	c	
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	3.2			
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	U		
8260C	ETHYLBENZENE	100-41-4	UG L	0.5	U		
8260C	ISOPROPYLBENZENE	98-82-8	UG L	0.5	U		
8260C	M- AND P-XYLENE	108-38-3/106-42	UG L	1	U		
8260C	METHYL ACETATE	79-20-9	UG L	0.75	U		
8260C	METHYL CYCLOHEXANE	108-87-2	UG L	0.5	U		
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG L	0.5	U		
8260C	METHYLENE CHLORIDE	75-09-2	UG L	2.5	U		
8260C	O-XYLENE	95-47-6	UG L	0.5	U		
8260C	STYRENE	100-42-5	UG L	0.5	U		
8260C	TETRACHLOROETHENE	127-18-4	UG L	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	930			
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	12			

				Sample Delivery Group		BETHPAGE-3	
				Lab ID	TI0116-4		
				Sample ID	RE103D2-GW-121415		
				Sample Date	12/14/2015		
				Sample Type	Groundwater		
Method	Analyte	CAS No	Units	Result	Qual	RC	
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG L	0.5	U		
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG L	0.5	U		
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG L	3.2			
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	0.77	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	1.1	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	UJ	c	
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	1.1			
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG L	0.5	U		
8260C	CYCLOHEXANE	110-82-7	UG L	0.5	U		
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG L	0.5	U		
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG L	1	U		
8260C	ETHYLBENZENE	100-41-4	UG L	0.5	U		
8260C	ISOPROPYLBENZENE	98-82-8	UG L	0.5	U		
8260C	M- AND P-XYLENE	108-38-3/106-42	UG L	1	U		
8260C	METHYL ACETATE	79-20-9	UG L	0.75	U		
8260C	METHYL CYCLOHEXANE	108-87-2	UG L	0.5	U		
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG L	0.5	U		
8260C	METHYLENE CHLORIDE	75-09-2	UG L	2.5	U		
8260C	O-XYLENE	95-47-6	UG L	0.5	U		
8260C	STYRENE	100-42-5	UG L	0.5	U		
8260C	TETRACHLOROETHENE	127-18-4	UG L	0.72	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	620			
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-3	
				Lab ID	TI0116-5		
				Sample ID	RE104D1-GW-121515		
				Sample Date	12/15/2015		
				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	0.8	J		
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG L	0.5	U		
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG L	0.75	U		
8260C	1,2-DIBROMOETHANE	106-93-4	UG L	0.5	U		
8260C	1,2-DICHLOROBENZENE	95-50-1	UG L	0.5	U		
8260C	1,2-DICHLOROETHANE	107-06-2	UG L	0.5	U		
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG L	1.1	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	UJ	c	
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	1.1			
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG L	0.5	U		
8260C	CYCLOHEXANE	110-82-7	UG L	0.5	U		
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG L	0.5	U		
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG L	0.68	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	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	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	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	6.9			

				Sample Delivery Group		BETHPAGE-3	
				Lab ID	TI0116-6		
				Sample ID	RE104D2-GW-121515		
				Sample Date	12/15/2015		
				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	2.7			
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	UJ	c	
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	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	U		
8260C	ETHYLBENZENE	100-41-4	UG L	0.5	U		
8260C	ISOPROPYLBENZENE	98-82-8	UG L	0.5	U		
8260C	M- AND P-XYLENE	108-38-3/106-42	UG L	1	U		
8260C	METHYL ACETATE	79-20-9	UG L	0.75	U		
8260C	METHYL CYCLOHEXANE	108-87-2	UG L	0.5	U		
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG L	0.5	U		
8260C	METHYLENE CHLORIDE	75-09-2	UG L	2.5	U		
8260C	O-XYLENE	95-47-6	UG L	0.5	U		
8260C	STYRENE	100-42-5	UG L	0.5	U		
8260C	TETRACHLOROETHENE	127-18-4	UG L	0.5	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	6.8			
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	0.22	J		

				Sample Delivery Group		BETHPAGE-3	
				Lab ID	TI0116-7		
				Sample ID	RE104D3-GW-121515		
				Sample Date	12/15/2015		
				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	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	UJ	c	
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	U		
8260C	ETHYLBENZENE	100-41-4	UG L	0.5	U		
8260C	ISOPROPYLBENZENE	98-82-8	UG L	0.5	U		
8260C	M- AND P-XYLENE	108-38-3/106-42	UG L	1	U		
8260C	METHYL ACETATE	79-20-9	UG L	0.75	U		
8260C	METHYL CYCLOHEXANE	108-87-2	UG L	0.5	U		
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG L	0.5	U		
8260C	METHYLENE CHLORIDE	75-09-2	UG L	2.5	U		
8260C	O-XYLENE	95-47-6	UG L	0.5	U		
8260C	STYRENE	100-42-5	UG L	0.5	U		
8260C	TETRACHLOROETHENE	127-18-4	UG L	0.5	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-3	
				Lab ID	TI0116-8		
				Sample ID	DUPLICATE1-GW-121515		
				Sample Date	12/15/2015		
				Sample Type	Field Duplicate		
Method	Analyte	CAS No	Units	Result	Qual	RC	
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG L	0.5	U		
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG L	0.5	U		
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG L	0.5	U		
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG L	0.5	U		
8260C	1,1-DICHLOROETHANE	75-34-3	UG L	0.5	U		
8260C	1,1-DICHLOROETHENE	75-35-4	UG L	0.5	U		
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG L	0.5	U		
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG L	0.75	U		
8260C	1,2-DIBROMOETHANE	106-93-4	UG L	0.5	U		
8260C	1,2-DICHLOROBENZENE	95-50-1	UG L	0.5	U		
8260C	1,2-DICHLOROETHANE	107-06-2	UG L	0.5	U		
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG L	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	UJ		c
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	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	U		
8260C	ETHYLBENZENE	100-41-4	UG L	0.5	U		
8260C	ISOPROPYLBENZENE	98-82-8	UG L	0.5	U		
8260C	M- AND P-XYLENE	108-38-3/106-42	UG L	1	U		
8260C	METHYL ACETATE	79-20-9	UG L	0.75	U		
8260C	METHYL CYCLOHEXANE	108-87-2	UG L	0.5	U		
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG L	0.5	U		
8260C	METHYLENE CHLORIDE	75-09-2	UG L	2.5	U		
8260C	O-XYLENE	95-47-6	UG L	0.5	U		
8260C	STYRENE	100-42-5	UG L	0.5	U		
8260C	TETRACHLOROETHENE	127-18-4	UG L	0.5	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	6.8			
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	0.28			



				Sample Delivery Group	BETHPAGE-3		
				Lab ID	TI0116-9		
				Sample ID	RE122D1-GW-121515		
				Sample Date	12/15/2015		
				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	4.7			
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	0.63	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	1.9	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	UJ		c
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	1	U		
8260C	ETHYLBENZENE	100-41-4	UG L	0.5	U		
8260C	ISOPROPYLBENZENE	98-82-8	UG L	0.5	U		
8260C	M- AND P-XYLENE	108-38-3/106-42	UG L	1	U		
8260C	METHYL ACETATE	79-20-9	UG L	0.75	U		
8260C	METHYL CYCLOHEXANE	108-87-2	UG L	0.5	U		
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG L	0.5	U		
8260C	METHYLENE CHLORIDE	75-09-2	UG L	2.5	U		
8260C	O-XYLENE	95-47-6	UG L	0.5	U		
8260C	STYRENE	100-42-5	UG L	0.5	U		
8260C	TETRACHLOROETHENE	127-18-4	UG L	1.5	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	600			
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	8.7			

				Sample Delivery Group	BETHPAGE-3		
				Lab ID	TI0214-1		
				Sample ID	TRIP BLANK 121615		
				Sample Date	12/16/2015		
				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	U		
8260C	2-HEXANONE	591-78-6	UG L	2.5	U		
8260C	4-METHYL-2-PENTANONE	108-10-1	UG L	2.5	UJ		c
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	UJ		c
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	U		
8260C	ETHYLBENZENE	100-41-4	UG L	0.5	U		
8260C	ISOPROPYLBENZENE	98-82-8	UG L	0.5	U		
8260C	M- AND P-XYLENE	108-38-3/106-42	UG L	1	U		
8260C	METHYL ACETATE	79-20-9	UG L	0.75	U		
8260C	METHYL CYCLOHEXANE	108-87-2	UG L	0.5	U		
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG L	0.5	U		
8260C	METHYLENE CHLORIDE	75-09-2	UG L	2.5	U		
8260C	O-XYLENE	95-47-6	UG L	0.5	U		
8260C	STYRENE	100-42-5	UG L	0.5	U		
8260C	TETRACHLOROETHENE	127-18-4	UG L	0.5	UJ		c
8260C	TOLUENE	108-88-3	UG L	0.37	J		
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG L	0.5	U		
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG L	0.5	U		
8260C	TRICHLOROETHENE	79-01-6	UG L	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-3	
				Lab ID		TI0214-10	
				Sample ID		TT101D1-GW-121715	
				Sample Date		12/17/2015	
				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	16			
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG L	0.48	J		
8260C	1,1-DICHLOROETHANE	75-34-3	UG L	0.5	U		
8260C	1,1-DICHLOROETHENE	75-35-4	UG L	4.6			
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG L	0.5	U		
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG L	0.75	U		
8260C	1,2-DIBROMOETHANE	106-93-4	UG L	0.5	U		
8260C	1,2-DICHLOROBENZENE	95-50-1	UG L	0.5	U		
8260C	1,2-DICHLOROETHANE	107-06-2	UG L	0.5	U		
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG L	1.9	J		
8260C	1,2-DICHLOROPROPANE	78-87-5	UG L	0.5	U		
8260C	1,3-DICHLOROBENZENE	541-73-1	UG L	0.5	U		
8260C	1,4-DICHLOROBENZENE	106-46-7	UG L	0.5	U		
8260C	2-BUTANONE	78-93-3	UG L	2.5	U		
8260C	2-HEXANONE	591-78-6	UG L	2.5	U		
8260C	4-METHYL-2-PENTANONE	108-10-1	UG L	2.5	UJ	c	
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	UJ	c	
8260C	CARBON DISULFIDE	75-15-0	UG L	0.5	U		
8260C	CARBON TETRACHLORIDE	56-23-5	UG L	1.9			
8260C	CHLOROBENZENE	108-90-7	UG L	0.5	U		
8260C	CHLOROETHANE	75-00-3	UG L	1	UJ	c	
8260C	CHLOROFORM	67-66-3	UG L	1			
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	1.8	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	U		
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG L	0.5	U		
8260C	METHYLENE CHLORIDE	75-09-2	UG L	2.5	U		
8260C	O-XYLENE	95-47-6	UG L	0.5	U		
8260C	STYRENE	100-42-5	UG L	0.5	U		
8260C	TETRACHLOROETHENE	127-18-4	UG L	0.5	UJ	c	
8260C	TOLUENE	108-88-3	UG L	0.5	U		
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG L	0.5	U		
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG L	0.5	U		
8260C	TRICHLOROETHENE	79-01-6	UG L	200			
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	11			

				Sample Delivery Group	BETHPAGE-3		
				Lab ID	TI0214-12		
				Sample ID	FIELD1-FB-121615		
				Sample Date	12/16/2015		
				Sample Type	Field 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	U		
8260C	2-HEXANONE	591-78-6	UG L	2.5	U		
8260C	4-METHYL-2-PENTANONE	108-10-1	UG L	2.5	UJ		c
8260C	ACETONE	67-64-1	UG L	2.2	J		
8260C	BENZENE	71-43-2	UG L	0.5	U		
8260C	BROMODICHLOROMETHANE	75-27-4	UG L	0.5	U		
8260C	BROMOFORM	75-25-2	UG L	0.5	U		
8260C	BROMOMETHANE	74-83-9	UG L	1	UJ		c
8260C	CARBON DISULFIDE	75-15-0	UG L	0.5	U		
8260C	CARBON TETRACHLORIDE	56-23-5	UG L	0.5	U		
8260C	CHLOROBENZENE	108-90-7	UG L	0.5	U		
8260C	CHLOROETHANE	75-00-3	UG L	1	UJ		c
8260C	CHLOROFORM	67-66-3	UG L	0.5	U		
8260C	CHLOROMETHANE	74-87-3	UG L	1	U		
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG L	0.5	U		
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG L	0.5	U		
8260C	CYCLOHEXANE	110-82-7	UG L	0.5	U		
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG L	0.5	U		
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG L	1	U		
8260C	ETHYLBENZENE	100-41-4	UG L	0.5	U		
8260C	ISOPROPYLBENZENE	98-82-8	UG L	0.5	U		
8260C	M- AND P-XYLENE	108-38-3/106-42	UG L	1	U		
8260C	METHYL ACETATE	79-20-9	UG L	0.75	U		
8260C	METHYL CYCLOHEXANE	108-87-2	UG L	0.5	U		
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG L	0.5	U		
8260C	METHYLENE CHLORIDE	75-09-2	UG L	2.5	U		
8260C	O-XYLENE	95-47-6	UG L	0.5	U		
8260C	STYRENE	100-42-5	UG L	0.5	U		
8260C	TETRACHLOROETHENE	127-18-4	UG L	0.5	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	0.36	J		
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			

				Sample Delivery Group		BETHPAGE-3	
				Lab ID	TI0214-4		
				Sample ID	RE114D3-GW-121615		
				Sample Date	12/16/2015		
				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	13			
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	1.1			
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG L	0.5	U		
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG L	0.75	U		
8260C	1,2-DIBROMOETHANE	106-93-4	UG L	0.5	U		
8260C	1,2-DICHLOROBENZENE	95-50-1	UG L	0.5	U		
8260C	1,2-DICHLOROETHANE	107-06-2	UG L	0.5	U		
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG L	0.67	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	UJ	c	
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	UJ	c	
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	0.67	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	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	43			
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-3	
				Lab ID	TI0214-5		
				Sample ID	RE114D2-GW-121615		
				Sample Date	12/16/2015		
				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	0.5	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	0.82	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	UJ	c	
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	UJ	c	
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	0.4	J		
8260C	CHLOROMETHANE	74-87-3	UG L	1	U		
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG L	0.82	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	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	70			
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	2.5			

				Sample Delivery Group		BETHPAGE-3	
				Lab ID	TI0214-6		
				Sample ID	RE105D1-GW-121715		
				Sample Date	12/17/2015		
				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	8.7			
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	1.3			
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	1.7	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	UJ	c	
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	UJ	c	
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	0.38	J		
8260C	CHLOROMETHANE	74-87-3	UG L	1	U		
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG L	1.7			
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	0.58	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	U		
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG L	0.5	U		
8260C	METHYLENE CHLORIDE	75-09-2	UG L	2.5	U		
8260C	O-XYLENE	95-47-6	UG L	0.5	U		
8260C	STYRENE	100-42-5	UG L	0.5	U		
8260C	TETRACHLOROETHENE	127-18-4	UG L	0.5	UJ	c	
8260C	TOLUENE	108-88-3	UG L	0.5	U		
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG L	0.5	U		
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG L	0.5	U		
8260C	TRICHLOROETHENE	79-01-6	UG L	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	10			

				Sample Delivery Group		BETHPAGE-3	
				Lab ID	TI0214-7		
				Sample ID	RE105D2-GW-121715		
				Sample Date	12/17/2015		
				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	26			
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG L	1.3			
8260C	1,1-DICHLOROETHANE	75-34-3	UG L	1.9			
8260C	1,1-DICHLOROETHENE	75-35-4	UG L	7			
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG L	0.5	U		
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG L	0.75	U		
8260C	1,2-DIBROMOETHANE	106-93-4	UG L	0.5	U		
8260C	1,2-DICHLOROBENZENE	95-50-1	UG L	0.5	U		
8260C	1,2-DICHLOROETHANE	107-06-2	UG L	0.5	U		
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG L	4			
8260C	1,2-DICHLOROPROPANE	78-87-5	UG L	0.5	U		
8260C	1,3-DICHLOROBENZENE	541-73-1	UG L	0.5	U		
8260C	1,4-DICHLOROBENZENE	106-46-7	UG L	0.5	U		
8260C	2-BUTANONE	78-93-3	UG L	2.5	U		
8260C	2-HEXANONE	591-78-6	UG L	2.5	U		
8260C	4-METHYL-2-PENTANONE	108-10-1	UG L	2.5	UJ	c	
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	UJ	c	
8260C	CARBON DISULFIDE	75-15-0	UG L	0.5	U		
8260C	CARBON TETRACHLORIDE	56-23-5	UG L	3			
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	2			
8260C	CHLOROMETHANE	74-87-3	UG L	1	U		
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG L	4			
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	0.45	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	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	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	1800			
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	5.8			



				Sample Delivery Group		BETHPAGE-3	
				Lab ID	TI0214-8		
				Sample ID	TT101D2-GW-122115		
				Sample Date	12/17/2015		
				Sample Type	Groundwater		
Method	Analyte	CAS No	Units	Result	Qual	RC	
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG L	0.34	J	m	
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG L	0.5	UJ	m	
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG L	19			
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG L	0.5	J	m	
8260C	1,1-DICHLOROETHANE	75-34-3	UG L	0.5	UJ	m	
8260C	1,1-DICHLOROETHENE	75-35-4	UG L	3.6	J	m,fd	
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG L	0.5	UJ	m	
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG L	0.75	UJ	m	
8260C	1,2-DIBROMOETHANE	106-93-4	UG L	0.5	UJ	m	
8260C	1,2-DICHLOROBENZENE	95-50-1	UG L	0.5	UJ	m	
8260C	1,2-DICHLOROETHANE	107-06-2	UG L	0.5	UJ	m	
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG L	1.7	J	m	
8260C	1,2-DICHLOROPROPANE	78-87-5	UG L	0.5	UJ	m	
8260C	1,3-DICHLOROBENZENE	541-73-1	UG L	0.5	UJ	m	
8260C	1,4-DICHLOROBENZENE	106-46-7	UG L	0.5	UJ	m	
8260C	2-BUTANONE	78-93-3	UG L	2.5	U		
8260C	2-HEXANONE	591-78-6	UG L	2.5	U		
8260C	4-METHYL-2-PENTANONE	108-10-1	UG L	2.5	UJ	c	
8260C	ACETONE	67-64-1	UG L	2.5	U		
8260C	BENZENE	71-43-2	UG L	0.5	UJ	m	
8260C	BROMODICHLOROMETHANE	75-27-4	UG L	0.5	UJ	m	
8260C	BROMOFORM	75-25-2	UG L	0.5	UJ	m	
8260C	BROMOMETHANE	74-83-9	UG L	1	UJ	c	
8260C	CARBON DISULFIDE	75-15-0	UG L	0.5	U		
8260C	CARBON TETRACHLORIDE	56-23-5	UG L	1.3	J	m	
8260C	CHLOROBENZENE	108-90-7	UG L	0.5	UJ	m	
8260C	CHLOROETHANE	75-00-3	UG L	1	UJ	c	
8260C	CHLOROFORM	67-66-3	UG L	0.9	J	m	
8260C	CHLOROMETHANE	74-87-3	UG L	1	U		
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG L	1.7	J	m	
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG L	0.5	UJ	m	
8260C	CYCLOHEXANE	110-82-7	UG L	0.5	UJ	m	
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG L	0.5	UJ	m	
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG L	1	U		
8260C	ETHYLBENZENE	100-41-4	UG L	0.5	UJ	m	
8260C	ISOPROPYLBENZENE	98-82-8	UG L	0.5	UJ	m	
8260C	M- AND P-XYLENE	108-38-3/106-42	UG L	1	UJ	m	
8260C	METHYL ACETATE	79-20-9	UG L	0.75	U		
8260C	METHYL CYCLOHEXANE	108-87-2	UG L	0.5	U		
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG L	0.5	U		
8260C	METHYLENE CHLORIDE	75-09-2	UG L	2.5	UJ	m	
8260C	O-XYLENE	95-47-6	UG L	0.5	UJ	m	
8260C	STYRENE	100-42-5	UG L	0.5	UJ	m	
8260C	TETRACHLOROETHENE	127-18-4	UG L	0.5	UJ	c	
8260C	TOLUENE	108-88-3	UG L	0.5	UJ	m	
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG L	0.5	UJ	m	
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG L	0.5	U		
8260C	TRICHLOROETHENE	79-01-6	UG L	510			
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	UJ	m	
8270D SIM	1,4-DIOXANE	123-91-1	UG L	1.7			

				Sample Delivery Group		BETHPAGE-3	
				Lab ID	TI0214-9		
				Sample ID	TT101D-GW-121715		
				Sample Date	12/17/2015		
				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	16			
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG L	0.5	U		
8260C	1,1-DICHLOROETHANE	75-34-3	UG L	0.84	J		
8260C	1,1-DICHLOROETHENE	75-35-4	UG L	3.4			
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	3.1			
8260C	1,2-DICHLOROPROPANE	78-87-5	UG L	0.5	U		
8260C	1,3-DICHLOROBENZENE	541-73-1	UG L	0.5	U		
8260C	1,4-DICHLOROBENZENE	106-46-7	UG L	0.5	U		
8260C	2-BUTANONE	78-93-3	UG L	2.5	U		
8260C	2-HEXANONE	591-78-6	UG L	2.5	U		
8260C	4-METHYL-2-PENTANONE	108-10-1	UG L	2.5	UJ	c	
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	UJ	c	
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	0.55	J		
8260C	CHLOROMETHANE	74-87-3	UG L	1	U		
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG L	3.1			
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG L	0.5	U		
8260C	CYCLOHEXANE	110-82-7	UG L	0.5	U		
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG L	0.5	U		
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG L	2.2			
8260C	ETHYLBENZENE	100-41-4	UG L	0.5	U		
8260C	ISOPROPYLBENZENE	98-82-8	UG L	0.5	U		
8260C	M- AND P-XYLENE	108-38-3/106-42	UG L	1	U		
8260C	METHYL ACETATE	79-20-9	UG L	0.75	U		
8260C	METHYL CYCLOHEXANE	108-87-2	UG L	0.5	U		
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG L	0.5	U		
8260C	METHYLENE CHLORIDE	75-09-2	UG L	2.5	U		
8260C	O-XYLENE	95-47-6	UG L	0.5	U		
8260C	STYRENE	100-42-5	UG L	0.5	U		
8260C	TETRACHLOROETHENE	127-18-4	UG L	0.5	UJ	c	
8260C	TOLUENE	108-88-3	UG L	0.5	U		
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG L	0.5	U		
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG L	0.5	U		
8260C	TRICHLOROETHENE	79-01-6	UG L	74			
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	8.4			

				Sample Delivery Group		BETHPAGE-3	
				Lab ID	TI0330-1		
				Sample ID	TRIP BLANK-121815		
				Sample Date	12/18/2015		
				Sample Type	Trip Blank		
Method	Analyte	CAS No	Units	Result	Qual	RC	
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG L	0.5	U		
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG L	0.5	U		
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG L	0.5	U		
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG L	0.5	U		
8260C	1,1-DICHLOROETHANE	75-34-3	UG L	0.5	U		
8260C	1,1-DICHLOROETHENE	75-35-4	UG L	0.5	U		
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG L	0.5	U		
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG L	0.75	U		
8260C	1,2-DIBROMOETHANE	106-93-4	UG L	0.5	U		
8260C	1,2-DICHLOROBENZENE	95-50-1	UG L	0.5	U		
8260C	1,2-DICHLOROETHANE	107-06-2	UG L	0.5	U		
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG L	1	U		
8260C	1,2-DICHLOROPROPANE	78-87-5	UG L	0.5	U		
8260C	1,3-DICHLOROBENZENE	541-73-1	UG L	0.5	U		
8260C	1,4-DICHLOROBENZENE	106-46-7	UG L	0.5	U		
8260C	2-BUTANONE	78-93-3	UG L	2.5	UJ	c	
8260C	2-HEXANONE	591-78-6	UG L	2.5	UJ	c	
8260C	4-METHYL-2-PENTANONE	108-10-1	UG L	2.5	U		
8260C	ACETONE	67-64-1	UG L	2.5	UJ	c	
8260C	BENZENE	71-43-2	UG L	0.5	U		
8260C	BROMODICHLOROMETHANE	75-27-4	UG L	0.5	U		
8260C	BROMOFORM	75-25-2	UG L	0.5	U		
8260C	BROMOMETHANE	74-83-9	UG L	1	U		
8260C	CARBON DISULFIDE	75-15-0	UG L	0.5	U		
8260C	CARBON TETRACHLORIDE	56-23-5	UG L	0.5	U		
8260C	CHLOROBENZENE	108-90-7	UG L	0.5	U		
8260C	CHLOROETHANE	75-00-3	UG L	1	UJ	c	
8260C	CHLOROFORM	67-66-3	UG L	0.5	U		
8260C	CHLOROMETHANE	74-87-3	UG L	0.45	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	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	UJ	c	
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-3	
				Lab ID		TI0330-10RA	
				Sample ID		RE107D1-GW-121815	
				Sample Date		12/18/2015	
				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	0.95	J		
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG L	0.5	U		
8260C	1,1-DICHLOROETHANE	75-34-3	UG L	0.5	U		
8260C	1,1-DICHLOROETHENE	75-35-4	UG L	0.5	U		
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG L	0.5	U		
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG L	0.75	U		
8260C	1,2-DIBROMOETHANE	106-93-4	UG L	0.5	U		
8260C	1,2-DICHLOROBENZENE	95-50-1	UG L	0.5	U		
8260C	1,2-DICHLOROETHANE	107-06-2	UG L	0.5	U		
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG L	0.21	J		
8260C	1,2-DICHLOROPROPANE	78-87-5	UG L	0.5	U		
8260C	1,3-DICHLOROBENZENE	541-73-1	UG L	0.5	U		
8260C	1,4-DICHLOROBENZENE	106-46-7	UG L	0.5	U		
8260C	2-BUTANONE	78-93-3	UG L	2.5	U		
8260C	2-HEXANONE	591-78-6	UG L	2.5	U		
8260C	4-METHYL-2-PENTANONE	108-10-1	UG L	2.5	U		
8260C	ACETONE	67-64-1	UG L	2.5	UJ	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	U		
8260C	CARBON TETRACHLORIDE	56-23-5	UG L	0.5	U		
8260C	CHLOROBENZENE	108-90-7	UG L	0.5	U		
8260C	CHLOROETHANE	75-00-3	UG L	1	UJ	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	0.21	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	UJ	c	
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	1.6	J	c	
8260C	TOLUENE	108-88-3	UG L	0.5	U		
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG L	0.5	U		
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG L	0.5	UJ	c	
8260C	TRICHLOROETHENE	79-01-6	UG L	17			
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	6.9			

				Sample Delivery Group		BETHPAGE-3	
				Lab ID	T10330-11		
				Sample ID	RE107D2-GW-121815		
				Sample Date	12/18/2015		
				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	15			
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	2.7			
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	UJ		c
8260C	2-HEXANONE	591-78-6	UG L	2.5	UJ		c
8260C	4-METHYL-2-PENTANONE	108-10-1	UG L	2.5	U		
8260C	ACETONE	67-64-1	UG L	2.5	UJ		c
8260C	BENZENE	71-43-2	UG L	0.5	U		
8260C	BROMODICHLOROMETHANE	75-27-4	UG L	0.5	U		
8260C	BROMOFORM	75-25-2	UG L	0.5	U		
8260C	BROMOMETHANE	74-83-9	UG L	1	U		
8260C	CARBON DISULFIDE	75-15-0	UG L	0.5	U		
8260C	CARBON TETRACHLORIDE	56-23-5	UG L	0.5	U		
8260C	CHLOROBENZENE	108-90-7	UG L	0.5	U		
8260C	CHLOROETHANE	75-00-3	UG L	1	UJ		c
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	U		
8260C	ETHYLBENZENE	100-41-4	UG L	0.5	U		
8260C	ISOPROPYLBENZENE	98-82-8	UG L	0.5	U		
8260C	M- AND P-XYLENE	108-38-3/106-42	UG L	1	U		
8260C	METHYL ACETATE	79-20-9	UG L	0.75	U		
8260C	METHYL CYCLOHEXANE	108-87-2	UG L	0.5	U		
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG L	0.5	U		
8260C	METHYLENE CHLORIDE	75-09-2	UG L	2.5	U		
8260C	O-XYLENE	95-47-6	UG L	0.5	U		
8260C	STYRENE	100-42-5	UG L	0.5	U		
8260C	TETRACHLOROETHENE	127-18-4	UG L	6.4			
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	UJ		c
8260C	TRICHLOROETHENE	79-01-6	UG L	140			
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	9.3			

				Sample Delivery Group	BETHPAGE-3		
				Lab ID	TI0330-12		
				Sample ID	RE123D3-GW-122115		
				Sample Date	12/21/2015		
				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	U		
8260C	2-HEXANONE	591-78-6	UG L	2.5	U		
8260C	4-METHYL-2-PENTANONE	108-10-1	UG L	2.5	U		
8260C	ACETONE	67-64-1	UG L	2.5	UJ		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	0.56	J		
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	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	UJ		c
8260C	METHYL CYCLOHEXANE	108-87-2	UG L	0.5	U		
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG L	0.5	U		
8260C	METHYLENE CHLORIDE	75-09-2	UG L	2.5	U		
8260C	O-XYLENE	95-47-6	UG L	0.5	U		
8260C	STYRENE	100-42-5	UG L	0.5	U		
8260C	TETRACHLOROETHENE	127-18-4	UG L	0.5	UJ		c
8260C	TOLUENE	108-88-3	UG L	0.5	U		
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG L	0.5	U		
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG L	0.5	UJ		c
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-3		
				Lab ID	TI0330-13		
				Sample ID	RE123D1-GW-122115		
				Sample Date	12/21/2015		
				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	U		
8260C	2-HEXANONE	591-78-6	UG L	2.5	U		
8260C	4-METHYL-2-PENTANONE	108-10-1	UG L	2.5	U		
8260C	ACETONE	67-64-1	UG L	2.5	UJ		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	U		
8260C	CARBON TETRACHLORIDE	56-23-5	UG L	0.5	U		
8260C	CHLOROBENZENE	108-90-7	UG L	0.5	U		
8260C	CHLOROETHANE	75-00-3	UG L	1	UJ		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	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	UJ		c
8260C	METHYL CYCLOHEXANE	108-87-2	UG L	0.5	U		
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG L	0.5	U		
8260C	METHYLENE CHLORIDE	75-09-2	UG L	2.5	U		
8260C	O-XYLENE	95-47-6	UG L	0.5	U		
8260C	STYRENE	100-42-5	UG L	0.5	U		
8260C	TETRACHLOROETHENE	127-18-4	UG L	0.5	UJ		c
8260C	TOLUENE	108-88-3	UG L	0.5	U		
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG L	0.5	U		
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG L	0.5	UJ		c
8260C	TRICHLOROETHENE	79-01-6	UG L	6.1			
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	5			

				Sample Delivery Group	BETHPAGE-3		
				Lab ID	TI0330-14		
				Sample ID	RE123D2-GW-122115		
				Sample Date	12/21/2015		
				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	U		
8260C	2-HEXANONE	591-78-6	UG L	2.5	U		
8260C	4-METHYL-2-PENTANONE	108-10-1	UG L	2.5	U		
8260C	ACETONE	67-64-1	UG L	2.5	UJ		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	U		
8260C	CARBON TETRACHLORIDE	56-23-5	UG L	0.5	U		
8260C	CHLOROBENZENE	108-90-7	UG L	0.5	U		
8260C	CHLOROETHANE	75-00-3	UG L	1	UJ		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	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	UJ		c
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	0.59	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	UJ		c
8260C	TRICHLOROETHENE	79-01-6	UG L	1.5			
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	0.7			



				Sample Delivery Group		BETHPAGE-3	
				Lab ID		TI0330-17	
				Sample ID		RE120D1-GW-121815	
				Sample Date		12/18/2015	
				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	42			
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG L	1.4			
8260C	1,1-DICHLOROETHANE	75-34-3	UG L	3.2			
8260C	1,1-DICHLOROETHENE	75-35-4	UG L	23			
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	4			
8260C	1,2-DICHLOROPROPANE	78-87-5	UG L	0.5	U		
8260C	1,3-DICHLOROBENZENE	541-73-1	UG L	0.5	U		
8260C	1,4-DICHLOROBENZENE	106-46-7	UG L	0.5	U		
8260C	2-BUTANONE	78-93-3	UG L	2.5	U		
8260C	2-HEXANONE	591-78-6	UG L	2.5	U		
8260C	4-METHYL-2-PENTANONE	108-10-1	UG L	2.5	U		
8260C	ACETONE	67-64-1	UG L	2.5	UJ	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	U		
8260C	CARBON TETRACHLORIDE	56-23-5	UG L	0.79	J		
8260C	CHLOROBENZENE	108-90-7	UG L	0.5	U		
8260C	CHLOROETHANE	75-00-3	UG L	1	UJ	c	
8260C	CHLOROFORM	67-66-3	UG L	0.99	J		
8260C	CHLOROMETHANE	74-87-3	UG L	1	UJ	c	
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG L	4			
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	0.38	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	UJ	c	
8260C	METHYL CYCLOHEXANE	108-87-2	UG L	0.5	U		
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG L	0.5	U		
8260C	METHYLENE CHLORIDE	75-09-2	UG L	2.5	U		
8260C	O-XYLENE	95-47-6	UG L	0.5	U		
8260C	STYRENE	100-42-5	UG L	0.5	U		
8260C	TETRACHLOROETHENE	127-18-4	UG L	2.1	J	c	
8260C	TOLUENE	108-88-3	UG L	0.5	U		
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG L	0.5	U		
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG L	0.5	UJ	c	
8260C	TRICHLOROETHENE	79-01-6	UG L	1300			
8260C	TRICHLOROFUOROMETHANE	75-69-4	UG L	0.39	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	12			

				Sample Delivery Group		BETHPAGE-3	
				Lab ID	TI0330-2		
				Sample ID	RE108D1-GW-122215		
				Sample Date	12/22/2015		
				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	1.4			
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	0.44	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	0.61	J		
8260C	1,2-DICHLOROPROPANE	78-87-5	UG L	0.5	U		
8260C	1,3-DICHLOROBENZENE	541-73-1	UG L	0.5	U		
8260C	1,4-DICHLOROBENZENE	106-46-7	UG L	0.5	U		
8260C	2-BUTANONE	78-93-3	UG L	2.5	UJ		c
8260C	2-HEXANONE	591-78-6	UG L	2.5	UJ		c
8260C	4-METHYL-2-PENTANONE	108-10-1	UG L	2.5	U		
8260C	ACETONE	67-64-1	UG L	2.5	UJ		c
8260C	BENZENE	71-43-2	UG L	0.5	U		
8260C	BROMODICHLOROMETHANE	75-27-4	UG L	0.5	U		
8260C	BROMOFORM	75-25-2	UG L	0.5	U		
8260C	BROMOMETHANE	74-83-9	UG L	1	U		
8260C	CARBON DISULFIDE	75-15-0	UG L	0.5	U		
8260C	CARBON TETRACHLORIDE	56-23-5	UG L	0.5	U		
8260C	CHLOROBENZENE	108-90-7	UG L	0.5	U		
8260C	CHLOROETHANE	75-00-3	UG L	1	UJ		c
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	0.61	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	U		
8260C	STYRENE	100-42-5	UG L	0.5	U		
8260C	TETRACHLOROETHENE	127-18-4	UG L	1.2			
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	UJ		c
8260C	TRICHLOROETHENE	79-01-6	UG L	110			
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	6.7			

				Sample Delivery Group		BETHPAGE-3	
				Lab ID		TI0330-3DL	
				Sample ID		RE108D2-GW-122215	
				Sample Date		12/22/2015	
				Sample Type		Groundwater	
Method	Analyte	CAS No	Units	Result	Qual	RC	
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG L	1.4	J		
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG L	2.5	U		
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG L	6.2			
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG L	2.5	U		
8260C	1,1-DICHLOROETHANE	75-34-3	UG L	5.1			
8260C	1,1-DICHLOROETHENE	75-35-4	UG L	9			
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG L	2.5	U		
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG L	3.8	U		
8260C	1,2-DIBROMOETHANE	106-93-4	UG L	2.5	U		
8260C	1,2-DICHLOROBENZENE	95-50-1	UG L	2.5	U		
8260C	1,2-DICHLOROETHANE	107-06-2	UG L	2.5	U		
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG L	9	J		
8260C	1,2-DICHLOROPROPANE	78-87-5	UG L	2.5	U		
8260C	1,3-DICHLOROBENZENE	541-73-1	UG L	2.5	U		
8260C	1,4-DICHLOROBENZENE	106-46-7	UG L	2.5	U		
8260C	2-BUTANONE	78-93-3	UG L	12	UJ	c	
8260C	2-HEXANONE	591-78-6	UG L	12	UJ	c	
8260C	4-METHYL-2-PENTANONE	108-10-1	UG L	12	U		
8260C	ACETONE	67-64-1	UG L	12	UJ	c	
8260C	BENZENE	71-43-2	UG L	2.5	U		
8260C	BROMODICHLOROMETHANE	75-27-4	UG L	2.5	U		
8260C	BROMOFORM	75-25-2	UG L	2.5	U		
8260C	BROMOMETHANE	74-83-9	UG L	5	U		
8260C	CARBON DISULFIDE	75-15-0	UG L	2.5	U		
8260C	CARBON TETRACHLORIDE	56-23-5	UG L	1.8	J		
8260C	CHLOROBENZENE	108-90-7	UG L	2.5	U		
8260C	CHLOROETHANE	75-00-3	UG L	5	UJ	c	
8260C	CHLOROFORM	67-66-3	UG L	4.4	J		
8260C	CHLOROMETHANE	74-87-3	UG L	5	U		
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG L	9			
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG L	2.5	U		
8260C	CYCLOHEXANE	110-82-7	UG L	2.5	U		
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG L	2.5	U		
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG L	5	U		
8260C	ETHYLBENZENE	100-41-4	UG L	2.5	U		
8260C	ISOPROPYLBENZENE	98-82-8	UG L	2.5	U		
8260C	M- AND P-XYLENE	108-38-3/106-42	UG L	5	U		
8260C	METHYL ACETATE	79-20-9	UG L	3.8	U		
8260C	METHYL CYCLOHEXANE	108-87-2	UG L	2.5	U		
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG L	2.5	U		
8260C	METHYLENE CHLORIDE	75-09-2	UG L	12	U		
8260C	O-XYLENE	95-47-6	UG L	2.5	U		
8260C	STYRENE	100-42-5	UG L	2.5	U		
8260C	TETRACHLOROETHENE	127-18-4	UG L	2.5	U		
8260C	TOLUENE	108-88-3	UG L	2.5	U		
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG L	2.5	U		
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG L	2.5	UJ	c	
8260C	TRICHLOROETHENE	79-01-6	UG L	2900	J	ej	
8260C	TRICHLOROFUOROMETHANE	75-69-4	UG L	5	U		
8260C	VINYL CHLORIDE	75-01-4	UG L	5	U		
8260C	XYLENES, TOTAL	1330-20-7	UG L	7.5	U		
8270D SIM	1,4-DIOXANE	123-91-1	UG L	8.8			

				Sample Delivery Group		BETHPAGE-3	
				Lab ID	TI0330-4		
				Sample ID	RE114D1-GW-122115		
				Sample Date	12/21/2015		
				Sample Type	Groundwater		
Method	Analyte	CAS No	Units	Result	Qual	RC	
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG L	0.64	J	s	
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	20	J	s	
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG L	1.6	J	s	
8260C	1,1-DICHLOROETHANE	75-34-3	UG L	1.5	J	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	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	5.1	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	U		
8260C	ACETONE	67-64-1	UG L	2.5	UJ	c	
8260C	BENZENE	71-43-2	UG L	0.5	U		
8260C	BROMODICHLOROMETHANE	75-27-4	UG L	0.5	U		
8260C	BROMOFORM	75-25-2	UG L	0.5	U		
8260C	BROMOMETHANE	74-83-9	UG L	1	U		
8260C	CARBON DISULFIDE	75-15-0	UG L	0.5	U		
8260C	CARBON TETRACHLORIDE	56-23-5	UG L	2.5	J	s	
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	2.9	J	s	
8260C	CHLOROMETHANE	74-87-3	UG L	1	U		
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG L	5.1	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	J	s	
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	0.3	J	s	
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	UJ	c	
8260C	TRICHLOROETHENE	79-01-6	UG L	370			
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	5.5			

				Sample Delivery Group	BETHPAGE-3		
				Lab ID	TI0330-5		
				Sample ID	RE121D1-GW-122115		
				Sample Date	12/21/2015		
				Sample Type	Groundwater		
Method	Analyte	CAS No	Units	Result	Qual	RC	
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG L	0.38	J		
8260C	1,1,2,2-TETRACHLOROETHANE	79-34-5	UG L	0.5	U		
8260C	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	76-13-1	UG L	8.3			
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	2.1			
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG L	0.5	U		
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG L	0.75	U		
8260C	1,2-DIBROMOETHANE	106-93-4	UG L	0.5	U		
8260C	1,2-DICHLOROBENZENE	95-50-1	UG L	0.5	U		
8260C	1,2-DICHLOROETHANE	107-06-2	UG L	0.38	J		
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG L	0.96	J		
8260C	1,2-DICHLOROPROPANE	78-87-5	UG L	0.5	U		
8260C	1,3-DICHLOROBENZENE	541-73-1	UG L	0.5	U		
8260C	1,4-DICHLOROBENZENE	106-46-7	UG L	0.5	U		
8260C	2-BUTANONE	78-93-3	UG L	2.5	UJ		c
8260C	2-HEXANONE	591-78-6	UG L	2.5	UJ		c
8260C	4-METHYL-2-PENTANONE	108-10-1	UG L	2.5	U		
8260C	ACETONE	67-64-1	UG L	2.5	UJ		c
8260C	BENZENE	71-43-2	UG L	0.5	U		
8260C	BROMODICHLOROMETHANE	75-27-4	UG L	0.5	U		
8260C	BROMOFORM	75-25-2	UG L	0.5	U		
8260C	BROMOMETHANE	74-83-9	UG L	1	U		
8260C	CARBON DISULFIDE	75-15-0	UG L	0.5	U		
8260C	CARBON TETRACHLORIDE	56-23-5	UG L	0.34	J		
8260C	CHLOROBENZENE	108-90-7	UG L	0.5	U		
8260C	CHLOROETHANE	75-00-3	UG L	1	UJ		c
8260C	CHLOROFORM	67-66-3	UG L	0.47	J		
8260C	CHLOROMETHANE	74-87-3	UG L	1	U		
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG L	0.96	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	2.2			
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	UJ		c
8260C	TRICHLOROETHENE	79-01-6	UG L	29			
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	6.8			

				Sample Delivery Group	BETHPAGE-3		
				Lab ID	TI0330-6		
				Sample ID	RE121D2-GW-122115		
				Sample Date	12/21/2015		
				Sample Type	Groundwater		
Method	Analyte	CAS No	Units	Result	Qual	RC	
8260C	1,1,1-TRICHLOROETHANE	71-55-6	UG L	0.48	J	s	
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	17	J	s	
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG L	0.64	J	s	
8260C	1,1-DICHLOROETHANE	75-34-3	UG L	0.51	J	s	
8260C	1,1-DICHLOROETHENE	75-35-4	UG L	3.1	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	2.1	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	U		
8260C	ACETONE	67-64-1	UG L	2.5	UJ	c	
8260C	BENZENE	71-43-2	UG L	0.5	U		
8260C	BROMODICHLOROMETHANE	75-27-4	UG L	0.5	U		
8260C	BROMOFORM	75-25-2	UG L	0.5	U		
8260C	BROMOMETHANE	74-83-9	UG L	1	U		
8260C	CARBON DISULFIDE	75-15-0	UG L	0.5	U		
8260C	CARBON TETRACHLORIDE	56-23-5	UG L	3.1	J	s	
8260C	CHLOROBENZENE	108-90-7	UG L	0.5	U		
8260C	CHLOROETHANE	75-00-3	UG L	1	UJ	c	
8260C	CHLOROFORM	67-66-3	UG L	1.7	J	s	
8260C	CHLOROMETHANE	74-87-3	UG L	1	U		
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG L	2.1	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	0.85	J	s	
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	UJ	c	
8260C	TRICHLOROETHENE	79-01-6	UG L	480			
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	4.9			

				Sample Delivery Group	BETHPAGE-3		
				Lab ID	TI0330-7		
				Sample ID	DUPLICATE-GW-122115		
				Sample Date	12/21/2015		
				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	24			
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG L	0.65	J		
8260C	1,1-DICHLOROETHANE	75-34-3	UG L	0.81	J		
8260C	1,1-DICHLOROETHENE	75-35-4	UG L	5	J		fd
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG L	0.5	U		
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG L	0.75	U		
8260C	1,2-DIBROMOETHANE	106-93-4	UG L	0.5	U		
8260C	1,2-DICHLOROBENZENE	95-50-1	UG L	0.5	U		
8260C	1,2-DICHLOROETHANE	107-06-2	UG L	0.5	U		
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG L	2			
8260C	1,2-DICHLOROPROPANE	78-87-5	UG L	0.5	U		
8260C	1,3-DICHLOROBENZENE	541-73-1	UG L	0.5	U		
8260C	1,4-DICHLOROBENZENE	106-46-7	UG L	0.5	U		
8260C	2-BUTANONE	78-93-3	UG L	2.5	UJ		c
8260C	2-HEXANONE	591-78-6	UG L	2.5	UJ		c
8260C	4-METHYL-2-PENTANONE	108-10-1	UG L	2.5	U		
8260C	ACETONE	67-64-1	UG L	2.5	UJ		c
8260C	BENZENE	71-43-2	UG L	0.5	U		
8260C	BROMODICHLOROMETHANE	75-27-4	UG L	0.5	U		
8260C	BROMOFORM	75-25-2	UG L	0.5	U		
8260C	BROMOMETHANE	74-83-9	UG L	1	U		
8260C	CARBON DISULFIDE	75-15-0	UG L	0.5	U		
8260C	CARBON TETRACHLORIDE	56-23-5	UG L	1.4			
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	0.92	J		
8260C	CHLOROMETHANE	74-87-3	UG L	1	U		
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	U		
8260C	ETHYLBENZENE	100-41-4	UG L	0.5	U		
8260C	ISOPROPYLBENZENE	98-82-8	UG L	0.5	U		
8260C	M- AND P-XYLENE	108-38-3/106-42	UG L	1	U		
8260C	METHYL ACETATE	79-20-9	UG L	0.75	U		
8260C	METHYL CYCLOHEXANE	108-87-2	UG L	0.5	U		
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG L	0.5	U		
8260C	METHYLENE CHLORIDE	75-09-2	UG L	2.5	U		
8260C	O-XYLENE	95-47-6	UG L	0.5	U		
8260C	STYRENE	100-42-5	UG L	0.5	U		
8260C	TETRACHLOROETHENE	127-18-4	UG L	0.94	J		
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	UJ		c
8260C	TRICHLOROETHENE	79-01-6	UG L	590			
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	2.2			

				Sample Delivery Group		BETHPAGE-3	
				Lab ID	TI0428-1		
				Sample ID	RE120D2-GW-122915		
				Sample Date	12/29/2015		
				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	25			
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG L	0.64	J		
8260C	1,1-DICHLOROETHANE	75-34-3	UG L	1.1			
8260C	1,1-DICHLOROETHENE	75-35-4	UG L	5.7			
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG L	0.5	U		
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG L	0.75	U		
8260C	1,2-DIBROMOETHANE	106-93-4	UG L	0.5	U		
8260C	1,2-DICHLOROBENZENE	95-50-1	UG L	0.5	U		
8260C	1,2-DICHLOROETHANE	107-06-2	UG L	0.5	U		
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG L	3.4			
8260C	1,2-DICHLOROPROPANE	78-87-5	UG L	0.5	U		
8260C	1,3-DICHLOROBENZENE	541-73-1	UG L	0.5	U		
8260C	1,4-DICHLOROBENZENE	106-46-7	UG L	0.5	U		
8260C	2-BUTANONE	78-93-3	UG L	2.5	U		
8260C	2-HEXANONE	591-78-6	UG L	2.5	U		
8260C	4-METHYL-2-PENTANONE	108-10-1	UG L	2.5	U		
8260C	ACETONE	67-64-1	UG L	2.5	UJ	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	U		
8260C	CARBON TETRACHLORIDE	56-23-5	UG L	0.69	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	0.77	J		
8260C	CHLOROMETHANE	74-87-3	UG L	1	U		
8260C	CIS-1,2-DICHLOROETHENE	156-59-2	UG L	3.4			
8260C	CIS-1,3-DICHLOROPROPENE	10061-01-5	UG L	0.5	U		
8260C	CYCLOHEXANE	110-82-7	UG L	0.5	U		
8260C	DIBROMOCHLOROMETHANE	124-48-1	UG L	0.5	U		
8260C	DICHLORODIFLUOROMETHANE	75-71-8	UG L	1	U		
8260C	ETHYLBENZENE	100-41-4	UG L	0.5	U		
8260C	ISOPROPYLBENZENE	98-82-8	UG L	0.5	U		
8260C	M- AND P-XYLENE	108-38-3/106-42	UG L	1	U		
8260C	METHYL ACETATE	79-20-9	UG L	0.75	U		
8260C	METHYL CYCLOHEXANE	108-87-2	UG L	0.5	U		
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG L	0.5	U		
8260C	METHYLENE CHLORIDE	75-09-2	UG L	2.5	U		
8260C	O-XYLENE	95-47-6	UG L	0.5	U		
8260C	STYRENE	100-42-5	UG L	0.5	U		
8260C	TETRACHLOROETHENE	127-18-4	UG L	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	680			
8260C	TRICHLOROFUOROMETHANE	75-69-4	UG L	0.26	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	8.8			



				Sample Delivery Group		BETHPAGE-3	
				Lab ID		TI0428-2RA	
				Sample ID		RE120D3-GW-122915	
				Sample Date		12/29/2015	
				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	3.1			
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG L	0.5	U		
8260C	1,1-DICHLOROETHANE	75-34-3	UG L	0.5	U		
8260C	1,1-DICHLOROETHENE	75-35-4	UG L	0.5	U		
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG L	0.5	U		
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG L	0.75	U		
8260C	1,2-DIBROMOETHANE	106-93-4	UG L	0.5	U		
8260C	1,2-DICHLOROBENZENE	95-50-1	UG L	0.5	U		
8260C	1,2-DICHLOROETHANE	107-06-2	UG L	0.5	U		
8260C	1,2-DICHLOROETHENE, TOTAL	540-59-0	UG L	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	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	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	U		
8260C	ETHYLBENZENE	100-41-4	UG L	0.5	U		
8260C	ISOPROPYLBENZENE	98-82-8	UG L	0.5	U		
8260C	M- AND P-XYLENE	108-38-3/106-42	UG L	1	U		
8260C	METHYL ACETATE	79-20-9	UG L	0.75	U		
8260C	METHYL CYCLOHEXANE	108-87-2	UG L	0.5	U		
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG L	0.5	U		
8260C	METHYLENE CHLORIDE	75-09-2	UG L	2.5	U		
8260C	O-XYLENE	95-47-6	UG L	0.5	U		
8260C	STYRENE	100-42-5	UG L	0.5	U		
8260C	TETRACHLOROETHENE	127-18-4	UG L	0.5	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	29			
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	0.28			

				Sample Delivery Group		BETHPAGE-3	
				Lab ID		TI0428-3RA	
				Sample ID		RE107D3-GW-122915	
				Sample Date		12/29/2015	
				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	4.9			
8260C	1,1,2-TRICHLOROETHANE	79-00-5	UG L	0.5	U		
8260C	1,1-DICHLOROETHANE	75-34-3	UG L	0.5	U		
8260C	1,1-DICHLOROETHENE	75-35-4	UG L	0.5	U		
8260C	1,2,4-TRICHLOROBENZENE	120-82-1	UG L	0.5	U		
8260C	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	UG L	0.75	U		
8260C	1,2-DIBROMOETHANE	106-93-4	UG L	0.5	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	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	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	U		
8260C	ETHYLBENZENE	100-41-4	UG L	0.5	U		
8260C	ISOPROPYLBENZENE	98-82-8	UG L	0.5	U		
8260C	M- AND P-XYLENE	108-38-3/106-42	UG L	1	U		
8260C	METHYL ACETATE	79-20-9	UG L	0.75	U		
8260C	METHYL CYCLOHEXANE	108-87-2	UG L	0.5	U		
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG L	0.5	U		
8260C	METHYLENE CHLORIDE	75-09-2	UG L	2.5	U		
8260C	O-XYLENE	95-47-6	UG L	0.5	U		
8260C	STYRENE	100-42-5	UG L	0.5	U		
8260C	TETRACHLOROETHENE	127-18-4	UG L	0.5	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	0.36	J		
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-3		
				Lab ID	T10428-4		
				Sample ID	TRIP BLANK-122915		
				Sample Date	12/29/2015		
				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	U		
8260C	2-HEXANONE	591-78-6	UG_L	2.5	U		
8260C	4-METHYL-2-PENTANONE	108-10-1	UG_L	2.5	U		
8260C	ACETONE	67-64-1	UG_L	2.5	UJ	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	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	U		
8260C	ETHYLBENZENE	100-41-4	UG_L	0.5	U		
8260C	ISOPROPYLBENZENE	98-82-8	UG_L	0.5	U		
8260C	M- AND P-XYLENE	108-38-3/106-42	UG_L	1	U		
8260C	METHYL ACETATE	79-20-9	UG_L	0.75	U		
8260C	METHYL CYCLOHEXANE	108-87-2	UG_L	0.5	U		
8260C	METHYL TERT-BUTYL ETHER	1634-04-4	UG_L	0.5	U		
8260C	METHYLENE CHLORIDE	75-09-2	UG_L	2.5	U		
8260C	O-XYLENE	95-47-6	UG_L	0.5	U		
8260C	STYRENE	100-42-5	UG_L	0.5	U		
8260C	TETRACHLOROETHENE	127-18-4	UG_L	0.5	UJ	c	
8260C	TOLUENE	108-88-3	UG_L	0.5	U		
8260C	TRANS-1,2-DICHLOROETHENE	156-60-5	UG_L	0.5	U		
8260C	TRANS-1,3-DICHLOROPROPENE	10061-02-6	UG_L	0.5	U		
8260C	TRICHLOROETHENE	79-01-6	UG_L	0.5	U		
8260C	TRICHLOROFLUOROMETHANE	75-69-4	UG_L	1	U		
8260C	VINYL CHLORIDE	75-01-4	UG_L	1	U		
8260C	XYLENES, TOTAL	1330-20-7	UG_L	1.5	U		
8270D_SIM	1,4-DIOXANE	123-91-1	UG_L	NA			

**Notes:**

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

Appendix C  
Analytical Data Validation – ARCADIS

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

### **Data Review**

BETHPAGE, NEW YORK

Volatile and Semivolatile Analyses

SDGs #JC9792 and JC9923

Analyses Performed By:  
Accutest Laboratories  
Dayton, New Jersey

Report #24833R  
December 28, 2015  
Review Level: Tier II  
Project #NY001496.1514.NAVI4

## SUMMARY

This data quality assessment summarizes the review of Sample Delivery Groups (SDGs) # JC9792 and JC9923 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
JC9792	TB120215PP1	JC9792-1	Water	12/02/2015		X				
	FB120215PP1	JC9792-2	Water	12/02/2015		X	X			
	BPOW 6-5	JC9792-3	Water	12/02/2015		X	X			
	BPOW 6-6	JC9792-4	Water	12/02/2015		X	X			
JC9923	BPOW5-3	JC9923-1	Water	12/03/2015		X	X			
	FB120315PP1	JC9923-2	Water	12/03/2015		X	X			
	TB120315PP1	JC9923-3	Water	12/03/2015		X				

Note:

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 in SDGs JC9792 or JC9923.

### 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 SDGs JC9792 and JC9923.

## **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 SDGs JC9792 or JC9923.

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

Tentatively identified compounds (TICs) were identified in SDG JC9792 in sample location FB120215PP1. The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification. TICs are qualified as estimated (JN).

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

## DATA VALIDATION CHECKLIST FOR VOCs

VOCs: 524.2	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 SDGs JC9792 or JC9923.

## **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: December 28, 2015

PEER REVIEW BY: Todd Church

DATE: December 29, 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

FED-EX Tracking # <b>#5</b>		Bottle Order Control #	
Account Order #		Account Job # <b>509792</b>	
<b>Client / Reporting Information</b>		<b>Project Information</b>	
Company Name <b>Arcadis</b>		Project Name: <b>AGMNYM62235 // OU2 Monitoring Wells Northrop Grumman OU2 Hydro</b>	
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>		Street Address <b>630 Plaza Drive, Suite 600</b>	
Phone # Fax # <b>631-249-7600 631-249-7610</b>		City State Zip <b>Highlands Ranch, CO 80129</b>	
Sample(s) Name(s) <b>Pat Proroki 516-297-0247</b>		Client Purchase Order # <b>NY01496.1514.NAVI3</b>	
Project Manager <b>Carlo San Giovanni</b>		Attention: <b>Soma Das</b>	
<b>Turnaround Time (Business days)</b>		<b>Data Deliverable Information</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		<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 LUMML+	
Approved By (Accutest PM) / Date:		Commercial "A" = Results Only Commercial "B" = Results + QC Summary NJ Reduced = Results + QC Summary + Partial Raw data	
RL reporting for metals <b>VOCS 524.2 Full list</b> <b>√ 524 SL4 + 40, VMS + F113</b>		Comments / Special Instructions	
Sample Custody must be documented below each time samples change possession, including courier delivery.			
Relinquished by Sampler: <b>1 Pat Proroki</b>	Date Time: <b>12-12-15 1945</b>	Received By: <b>3 [Signature]</b>	Date Time: <b>12/15/15 11:30</b>
Relinquished by Sampler: <b>3</b>	Date Time:	Received By: <b>4</b>	Date Time:
Relinquished by:	Date Time:	Received By:	Date Time:
Custody Seal #		Preserved where applicable	
<input type="checkbox"/> Intact <input type="checkbox"/> Not Intact		On Ice <input checked="" type="checkbox"/> Cooler Temp: <b>1.90</b>	

5.1  
5



## Report of Analysis

Client Sample ID:	TB120215PP1	Date Sampled:	12/02/15
Lab Sample ID:	JC9792-1	Date Received:	12/03/15
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	1B100904.D	1	12/04/15	MD	n/a	n/a	V1B4777
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	1.5	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> TB120215PP1		<b>Date Sampled:</b> 12/02/15
<b>Lab Sample ID:</b> JC9792-1		<b>Date Received:</b> 12/03/15
<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	102%		78-114%
460-00-4	4-Bromofluorobenzene	103%		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

4.1  
4

Report of Analysis

Client Sample ID: FB120215PP1	Date Sampled: 12/02/15
Lab Sample ID: JC9792-2	Date Received: 12/03/15
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	1B100905.D	1	12/04/15	MD	n/a	n/a	V1B4777
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	2.1	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

4.2  
**4**

## Report of Analysis

Client Sample ID: FB120215PP1		Date Sampled: 12/02/15
Lab Sample ID: JC9792-2		Date Received: 12/03/15
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	101%		78-114%
460-00-4	4-Bromofluorobenzene	102%		77-115%

CAS No.	Tentatively Identified Compounds	R.T.	Est. Conc.	Units	Q
	alkene	8.24	.81	ug/l	J N
	Total TIC, Volatile		.81	ug/l	J N

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

Client Sample ID:	BPOW 6-5	Date Sampled:	12/02/15
Lab Sample ID:	JC9792-3	Date Received:	12/03/15
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	1B100906.D	1	12/04/15	MD	n/a	n/a	V1B4777
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	0.89	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: BPOW 6-5		Date Sampled: 12/02/15
Lab Sample ID: JC9792-3		Date Received: 12/03/15
Matrix: AQ - Ground 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	101%		78-114%
460-00-4	4-Bromofluorobenzene	102%		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

Client Sample ID:	BPOW 6-6	Date Sampled:	12/02/15
Lab Sample ID:	JC9792-4	Date Received:	12/03/15
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	1B100907.D	1	12/04/15	MD	n/a	n/a	V1B4777
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	0.40	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

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 6-6		<b>Date Sampled:</b> 12/02/15
<b>Lab Sample ID:</b> JC9792-4		<b>Date Received:</b> 12/03/15
<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	103%		78-114%
460-00-4	4-Bromofluorobenzene	104%		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

4.4  
4

## Report of Analysis

<b>Client Sample ID:</b> FB120215PP1	<b>Date Sampled:</b> 12/02/15
<b>Lab Sample ID:</b> JC9792-2	<b>Date Received:</b> 12/03/15
<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>	I99455.D	1	12/16/15	AMA	12/05/15	M:OP45645	M:MSI3717
Run #2							

Run #	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	66%		26-121%		
321-60-8	2-Fluorobiphenyl	60%		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

4.2  
4

## Report of Analysis

<b>Client Sample ID:</b> BPOW 6-5		<b>Date Sampled:</b> 12/02/15
<b>Lab Sample ID:</b> JC9792-3		<b>Date Received:</b> 12/03/15
<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>	I99456.D	1	12/16/15	AMA	12/05/15	M:OP45645	M:MSI3717
Run #2							

Run #	Initial Volume	Final Volume
Run #1	940 ml	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
123-91-1	1,4-Dioxane	ND	0.21	0.081	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	72%		26-121%		
321-60-8	2-Fluorobiphenyl	65%		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> BPOW 6-6		<b>Date Sampled:</b> 12/02/15
<b>Lab Sample ID:</b> JC9792-4		<b>Date Received:</b> 12/03/15
<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>	I99457.D	1	12/16/15	AMA	12/05/15	M:OP45645	M:MSI3717
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	ND	0.22	0.083	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

4.4  
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

Impulse

<b>FED-EX Tracking #</b> #5 <b>Accutest Quote #</b>		<b>Bottle Order Control #</b> Jc9923	
<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>631-249-7600</b> Fax #: <b>631-249-7610</b> Sample Name(s): <b>PA Perovki 516 297-627</b>		<b>Project Information</b> Project Name: <b>AGMNYM62235 // OU2 Monitoring 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> Zip: <b>80129</b> Client Purchase Order #: <b>NY001496.154, 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> B8270SIM14DIOX VOCs 524,2 Full list		<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		<b>Collection</b> MECH/DI/Vol # Date Time Sampled by Matrix # of bottles Number of preserved bottles HDI HNO3 HNO2 H2SO4 H2SO4 NONE ED Water MESH ENCORE	
Field ID / Point of Collection 1 BPOW 5-3 2 FB120315PP2 3 TB120315PP2		LAB USE ONLY 1 V1047	
<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"/> 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 CUMMIL+ Commercial "A" = Results Only Commercial "B" = Results + QC Summary NJ Reduced = Results + QC Summary + Partial Raw data	
<b>Comments / Special Instructions</b> INITIAL ASSESSMENT: <b>3808</b> LABEL VERIFICATION: <b>JA</b> RL reporting for metals VOCs 524,2 Full list 524,2 + 40, VMS + F113			
<b>Sample Custody must be documented below each time samples change possession, including courier delivery.</b>			
Relinquished by Sampler: 1 <b>Carlo San Giovanni</b> Relinquished by Sampler: 3 Relinquished by: 5	Date Time: 12/3/15 1945 Date Time: Date Time: Date Time:	Received By: 1 <b>James Kwag</b> Received By: 3 Received By: 5	Date Time: 12/4/15 1946 Date Time: Date Time: Date Time:
Relinquished by: 4 <b>Ken Arcadis</b>		Relinquished By: 4 Date Time: Date Time: Date Time:	
Custody Seal # Intact <input checked="" type="checkbox"/> Not Intact <input type="checkbox"/>		Preserved where applicable <input type="checkbox"/> On Ice <input checked="" type="checkbox"/> Cooler Temp. <b>2.5°C</b>	

5.1  
5







## Report of Analysis

<b>Client Sample ID:</b> BPOW5-3		<b>Date Sampled:</b> 12/03/15
<b>Lab Sample ID:</b> JC9923-1		<b>Date Received:</b> 12/04/15
<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	98%		78-114%
460-00-4	4-Bromofluorobenzene	97%		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> FB120315PP1		<b>Date Sampled:</b> 12/03/15
<b>Lab Sample ID:</b> JC9923-2		<b>Date Received:</b> 12/04/15
<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	98%		78-114%
460-00-4	4-Bromofluorobenzene	98%		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> TB120315PP1		<b>Date Sampled:</b> 12/03/15
<b>Lab Sample ID:</b> JC9923-3		<b>Date Received:</b> 12/04/15
<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	1B100961.D	1	12/08/15	MD	n/a	n/a	V1B4779
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	1.4	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> TB120315PP1		<b>Date Sampled:</b> 12/03/15
<b>Lab Sample ID:</b> JC9923-3		<b>Date Received:</b> 12/04/15
<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	100%		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> BPOW5-3		<b>Date Sampled:</b> 12/03/15
<b>Lab Sample ID:</b> JC9923-1		<b>Date Received:</b> 12/04/15
<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>	I99391.D	1	12/14/15	AMA	12/07/15	M:OP45658	M:MSI3715
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	0.39	0.21	0.080	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	77%		26-121%		
321-60-8	2-Fluorobiphenyl	68%		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> FB120315PP1		<b>Date Sampled:</b> 12/03/15
<b>Lab Sample ID:</b> JC9923-2		<b>Date Received:</b> 12/04/15
<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>	I99392.D	1	12/14/15	AMA	12/07/15	M:OP45658	M:MSI3715
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	71%		26-121%		
321-60-8	2-Fluorobiphenyl	64%		28-107%		
1718-51-0	Terphenyl-d14	83%		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

4.2  
4

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

### **Data Review**

BETHPAGE, NEW YORK

Volatile and Semivolatile Analyses

SDGs #JC8577, JC8603 and JC8685

Analyses Performed By:  
Accutest Laboratories  
Dayton, New Jersey

Report #24830R  
December 22, 2015  
Review Level: Tier II  
Project #NY001496.1514.NAVI4



## SUMMARY

This data quality assessment summarizes the review of Sample Delivery Groups (SDGs) # JC8577, JC8603, and JC8685 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
JC8577	BPOW5-6	JC8577-1	Water	11/13/2015		X	X			
	BPOW5-5	JC8577-2	Water	11/13/2015		X	X			
	FB111315PP1	JC8577-3	Water	11/13/2015		X	X			
	TB111315PP1	JC8577-4	Water	11/13/2015		X				
JC8603	BPOW5-2	JC8603-1	Water	11/12/2015		X	X			
	BPOW5-1	JC8603-2	Water	11/12/2015		X	X			
	FB111215PP1	JC8603-3	Water	11/12/2015		X	X			
	TB111215PP1	JC8603-4	Water	11/12/2015		X				
JC8685	BPOW 5-4	JC8685-1	Water	11/16/2015		X	X			
	FB111615PP1	JC8685-2	Water	11/16/2015		X	X			
	TB111615PP1	JC8685-3	Water	11/16/2015		X				

**Notes:**

1. Semivolatile (1,4-Dioxane) analysis was performed at Accutest Laboratories, Marlborough, MA (Accutest-NE) facility.
2. SDG JC8577: Matrix spike analysis was performed on sample location BPOW5-6 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.

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
<u>SDG JC8577:</u>			
BPOW5-5	Acetone	Detected sample results <RL and <BAL	"UB" at the RL
<u>SDG JC8603:</u>			
BPOW5-1	Acetone	Detected sample results <RL and <BAL	"UB" at the RL
<u>SDG JC8685:</u>			
BPOW 5-4	Acetone Methylene Chloride	Detected sample results <RL and <BAL	"UB" at the RL

RL Reporting limit

### 3. Surrogates/System Monitoring Compounds

All samples to be analyzed for organic compounds are spiked with surrogate compounds prior to sample preparation to evaluate overall laboratory performance and efficiency of the analytical technique. VOC

analysis requires that all surrogates associated with the analysis exhibit recoveries within the laboratory-established acceptance limits.

All surrogate recoveries were within control limits.

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

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

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

The MS exhibited acceptable recoveries in SDG JC8577.

A MS/MSD analysis was not performed on a sample location associated with SDGs JC8603 and JC8685.

#### 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 in SDG JC8685.

Sample locations associated with LCS analysis exhibiting recoveries outside of the control limits presented in the following table.

Sample Locations	Compound	LCS Recovery
<u>SDG JC8577:</u>		
BPOW5-6 BPOW5-5 FB111315PP1 TB111315PP1	2-Butanone	>UL
<u>SDG JC8603:</u>		
BPOW5-2 BPOW5-1 TB111215PP1	2-Butanone	>UL

The criteria used to evaluate the LCS recoveries are presented in the following table. In the case of an LCS 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

Control Limit	Sample Result	Qualification
< the lower control limit (LL) but > 10%	Non-detect	UJ
	Detect	J
< 10%	Non-detect	R
	Detect	J

## 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 laboratory duplicate sample results exhibited RPD within the control limit in SDG JC8577.

A laboratory duplicate sample was not performed on a sample location associated with SDGs JC8603 and JC8685.

## 8. System Performance and Overall Assessment

Tentatively identified compounds (TICs) were identified in SDG JC8577 in sample locations: BPOW5-5, FB111315PP1 and TB111315PP1; SDG JC8603 in sample locations: BPOW5-1, FB111215PP1 and TB111215PP1; and, SDG JC8685 sample locations: BPOW 5-4 and FB111615PP1. The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification. TICs are qualified as estimated (JN).

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

## DATA VALIDATION CHECKLIST FOR VOCs

VOCs: 524.2	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 SDGs JC8577, JC8603 or JC8685.



## **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 for SDGs JC8577, JC8603 and JC8685.

## **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: December 22, 2015

PEER REVIEW BY: Todd Church

DATE: December 29, 2015

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

GW  
FB  
WB

<b>FED-EX Tracking #</b> #4 <b>Accutest Quote #</b>		<b>Bottle Order Control #</b> <b>Accutest Job #</b> JC8577	
<b>Client / Reporting Information</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>631-249-7600</b> Fax #: <b>631-249-7610</b> Sample(s) Name(s): <b>Pat Proroki</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: <b>Highlands Ranch, CO</b> State: <b>CO</b> Zip: <b>80129</b> Project #: <b>1614</b> Client Purchase Order #: <b>NY001496.3144.NAVI3</b> Work Authorization #: <b>NY001496_2015</b> Attention: <b>* Soma Das</b> Project Manager: <b>Carlo San Giovanni</b>	
<b>Requested Analysis (see TEST CODE sheet)</b> VOCs 524 Full List B827051M 14 D10X		<b>Matrix Codes</b> 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	
<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>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 Format <input type="checkbox"/> NJ Reduced <input type="checkbox"/> EDD Format <input type="checkbox"/> Commercial "C" <input checked="" type="checkbox"/> Other CUMMUL+ Commercial "A" = Results Only Commercial "B" = Results + QC Summary NJ Reduced = Results + QC Summary + Partial Raw data	
<b>Comments / Special Instructions</b> VOCs 524 Full List V 524SL4 + 40, VMS + F113		<b>INITIAL ASSESSMENT</b> JA Don <b>LABEL VERIFICATION</b>	
<b>Sample Custody must be documented below each time samples change possession, including courier delivery.</b>			
Relinquished by: <b>CA Caporal</b> Date Time: <b>11/13/15 15:03</b>	Received By: <b>Robert Chambers</b> Date Time: <b>11-12-15 15:32</b>	Relinquished By: <b>Robert Chambers</b> Date Time: <b>11-13-15</b>	Received By: <b>[Signature]</b> Date Time: <b>11-13-15</b>
Relinquished by: <b>[Signature]</b> Date Time:	Received By: <b>[Signature]</b> Date Time:	Relinquished By: <b>[Signature]</b> Date Time:	Received By: <b>[Signature]</b> Date Time:
Relinquished by: <b>[Signature]</b> Date Time:	Received By: <b>[Signature]</b> Date Time:	Relinquished By: <b>[Signature]</b> Date Time:	Received By: <b>[Signature]</b> Date Time:
Custody Seal # <b>853</b> <input type="checkbox"/> Intact <input type="checkbox"/> Not Intact Preserved where applicable <input type="checkbox"/>		On Ice <input checked="" type="checkbox"/> Cooler Temp: <b>4.7°C</b>	

5.1  
5





## Report of Analysis

<b>Client Sample ID:</b> BPOW5-6		<b>Date Sampled:</b> 11/13/15
<b>Lab Sample ID:</b> JC8577-1		<b>Date Received:</b> 11/13/15
<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	97%		78-114%
460-00-4	4-Bromofluorobenzene	99%		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> BPOW5-5		<b>Date Sampled:</b> 11/13/15
<b>Lab Sample ID:</b> JC8577-2		<b>Date Received:</b> 11/13/15
<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	1B100653.D	1	11/17/15	MD	n/a	n/a	V1B4766
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 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

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> BPOW5-5		<b>Date Sampled:</b> 11/13/15
<b>Lab Sample ID:</b> JC8577-2		<b>Date Received:</b> 11/13/15
<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	101%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
78-84-2	Propanal, 2-methyl-	8.86	.54	ug/l	JN
	Total TIC, Volatile		.54	ug/l	J N

ND = Not detected      MDL = Method Detection Limit  
 RL = Reporting Limit

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> FB111315PP1		<b>Date Sampled:</b> 11/13/15
<b>Lab Sample ID:</b> JC8577-3		<b>Date Received:</b> 11/13/15
<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	101%		77-115%		
CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q	
287-92-3	Cyclopentane	8.24	24	ug/l	JN	
	Total TIC, Volatile		24	ug/l	J N	

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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> TB111315PP1		<b>Date Sampled:</b> 11/13/15
<b>Lab Sample ID:</b> JC8577-4		<b>Date Received:</b> 11/13/15
<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	1B100658.D	1	11/17/15	MD	n/a	n/a	V1B4766
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	1.7	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> TB111315PP1		<b>Date Sampled:</b> 11/13/15
<b>Lab Sample ID:</b> JC8577-4		<b>Date Received:</b> 11/13/15
<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	101%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	unknown	7.61	1.1	ug/l	J N
	Total TIC, Volatile		1.1	ug/l	J N

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

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

4.4  
4

## Report of Analysis

<b>Client Sample ID:</b> BPOW5-6		<b>Date Sampled:</b> 11/13/15
<b>Lab Sample ID:</b> JC8577-1		<b>Date Received:</b> 11/13/15
<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>	I99234.D	1	11/28/15	AMA	11/16/15	M:OP45433	M:MSI3705
Run #2							

Run #	Initial Volume	Final Volume
Run #1	970 ml	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
123-91-1	1,4-Dioxane	ND	0.21	0.078	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	81%		26-121%		
321-60-8	2-Fluorobiphenyl	71%		28-107%		
1718-51-0	Terphenyl-d14	87%		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> BPOW5-5	<b>Date Sampled:</b> 11/13/15
<b>Lab Sample ID:</b> JC8577-2	<b>Date Received:</b> 11/13/15
<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>	I99235.D	1	11/28/15	AMA	11/16/15	M:OP45433	M:MSI3705
Run #2							

Run #	Initial Volume	Final Volume
Run #1	970 ml	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
123-91-1	1,4-Dioxane	0.42	0.21	0.078	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	84%		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

4.2  
4



## Report of Analysis

<b>Client Sample ID:</b> FB111315PP1	<b>Date Sampled:</b> 11/13/15
<b>Lab Sample ID:</b> JC8577-3	<b>Date Received:</b> 11/13/15
<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>	I99236.D	1	11/28/15	AMA	11/16/15	M:OP45433	M:MSI3705
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	72%		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  
4



FED-EX Tracking # 774      Bottle Order Control #  
Accutest Order #      Accutest Job # ARC JC8603

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> 631-249-7600 <b>Fax #</b> 631-249-7610 <b>Sample(s) Name(s)</b> Pat Przekacki      516      287-6247      Phone #		<b>Project Name:</b> AGMNYM62235 // OU2 Outpost Wells Northrop Grumman OU2 Hydro <b>Street</b> <b>City State</b> Bathpage NY <b>Project #</b> 1614 NY001496-STAT-NAV13 <b>Client Purchase Order #</b> Work Authorization #: 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 State Zip</b> Highlands Ranch, CO 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 1 AIR - Air SOL - Other Solid WP - Wipe FB - Field Blank ED - Equipment Blank RB - Ribase Blank TB - Trip Blank					
<b>Accutest Sample #</b> -1 -2	<b>Field ID / Point of Collection</b> BPOW 5-2 BPOW 5-1	<b>MECH/DIVIAL #</b>	<b>Collection</b> <b>Date</b> <b>Time</b> <b>Sampled by</b> 11/12/15      1425      PP 11/12/15      1510      PP			<b>Mark</b> GW GW	<b># of bottles</b> 2 2	<b>Number of preserved bottles</b> VOCs 524.2 Full List B827051M 14 DIOX										<b>LAB USE ONLY</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>Approved By (Accutest P#):</b>     	<input type="checkbox"/> Commercial "A" (Level 1) <input type="checkbox"/> Commercial "B" (Level 2) <input type="checkbox"/> FULLT1 (Level 3+1) <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 CUMMUL+	<b>Comments / Special Instructions</b> VOCs 524.2 Full List      110C										
<b>Sample Custody must be documented below each time samples change possession, including courier delivery.</b>																			
<b>Relinquished by Sampler:</b> 1 <i>[Signature]</i>	<b>Date/Time:</b> 11/12/15 1845	<b>Received By:</b> 1 <i>[Signature]</i>	<b>Date/Time:</b> 11/13/15 9:25	<b>Relinquished by:</b> 2 <i>[Signature]</i>	<b>Date/Time:</b> 11/13/15 2	<b>Received By:</b> 2 <i>[Signature]</i>	<b>Relinquished by:</b> 3 <i>[Signature]</i>										<b>Date/Time:</b> 3	<b>Received By:</b> 3	
<b>Relinquished by:</b> 4	<b>Date/Time:</b> 4	<b>Received By:</b> 4	<b>Date/Time:</b> 4	<b>Relinquished by:</b> 5	<b>Date/Time:</b> 5	<b>Received By:</b> 5	<b>Date/Time:</b> 5	<b>Relinquished by:</b> 5										<b>Date/Time:</b> 5	<b>Received By:</b> 5
<b>Emergency &amp; Rush TIA date available VIA Lablink</b>																			
<b>Sample Custody must be documented below each time samples change possession, including courier delivery.</b>																			
<b>Relinquished by:</b> <input type="checkbox"/> Intact <input type="checkbox"/> Not Intact <b>Preserved where applicable:</b> <input type="checkbox"/> On Ice <input type="checkbox"/> Cooler Temp.																			

5.1  
5



## Report of Analysis

<b>Client Sample ID:</b> BPOW5-2		<b>Date Sampled:</b> 11/12/15
<b>Lab Sample ID:</b> JC8603-1		<b>Date Received:</b> 11/14/15
<b>Matrix:</b> AQ - 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	101%		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> BPOW5-1		<b>Date Sampled:</b> 11/12/15
<b>Lab Sample ID:</b> JC8603-2		<b>Date Received:</b> 11/14/15
<b>Matrix:</b> AQ - 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	102%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
78-84-2	Propanal, 2-methyl-	8.85	.56	ug/l	JN
	Total TIC, Volatile		.56	ug/l	JN

ND = Not detected      MDL = Method Detection Limit  
 RL = Reporting Limit

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> FB111215PP1		<b>Date Sampled:</b> 11/12/15
<b>Lab Sample ID:</b> JC8603-3		<b>Date Received:</b> 11/14/15
<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	100%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	alkene	8.26	24	ug/l	J N
	Total TIC, Volatile		24	ug/l	J N

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

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

4.3  
4



## Report of Analysis

<b>Client Sample ID:</b> TB111215PP1		<b>Date Sampled:</b> 11/12/15
<b>Lab Sample ID:</b> JC8603-4		<b>Date Received:</b> 11/14/15
<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	100%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
67-63-0	Isopropyl Alcohol	7.61	1.2	ug/l	JN
	Total TIC, Volatile		1.2	ug/l	JN

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  
4

## Report of Analysis

<b>Client Sample ID:</b> BPOW5-2	<b>Date Sampled:</b> 11/12/15
<b>Lab Sample ID:</b> JC8603-1	<b>Date Received:</b> 11/14/15
<b>Matrix:</b> AQ - 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>	I99223.D	1	11/27/15	AMA	11/17/15	M:OP45450	M:MSI3705
Run #2							

Run #	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	74%		26-121%		
321-60-8	2-Fluorobiphenyl	65%		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

4.1  
4

## Report of Analysis

<b>Client Sample ID:</b> BPOW5-1		<b>Date Sampled:</b> 11/12/15
<b>Lab Sample ID:</b> JC8603-2		<b>Date Received:</b> 11/14/15
<b>Matrix:</b> AQ - 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>	I99224.D	1	11/27/15	AMA	11/17/15	M:OP45450	M:MSI3705
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	ND	0.22	0.083	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	68%		26-121%		
321-60-8	2-Fluorobiphenyl	66%		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  
4

## Report of Analysis

<b>Client Sample ID:</b> FB111215PP1	<b>Date Sampled:</b> 11/12/15
<b>Lab Sample ID:</b> JC8603-3	<b>Date Received:</b> 11/14/15
<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>	R46141.D	1	11/19/15	AMA	11/17/15	M:OP45457	M:MSR1697
Run #2							

Run #	Initial Volume	Final Volume
Run #1	970 ml	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
123-91-1	1,4-Dioxane	ND	0.21	0.078	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	91%		26-121%		
321-60-8	2-Fluorobiphenyl	71%		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

4.3  
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

<b>Client / Reporting Information</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>631-249-7600</b> Fax #: <b>631-249-7610</b> Sample(s) Name(s): <b>Pat Proroki 516 287-0247</b>		<b>Project Information</b> Project Name: <b>AGMNYM62235 // OU2 Outpost Wells Northrop Grumman OU2 Hydro</b> Street: _____ City: _____ State: _____ Zip: _____ 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 #: <b>NY001496.5741.NAV13</b> Work Authorization #: <b>NY001496_2015</b> Project Manager: <b>Carlo San Giovanni</b>		FED-EX Tracking # <b>444</b> Bottle Order Control # <b>JL8685</b> Accutest Quote # _____ Accutest Job # _____ <b>Requested Analysis (see TEST CODE sheet)</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 WIP - Wipe FB - Field Blank EB - Equipment Blank RB - Rinse Blank TB - Trip Blank LAB USE ONLY <b>V855</b>	
<b>Collection</b> MECH/UDI Vial # _____ Date _____ Time _____ Sampled by: _____ Matrix: _____ # of bottles: _____ Number of preserved bottles: HCl _____ HNO3 _____ H2SO4 _____ NONE _____ DI Water _____ MICH _____ ENDORNE _____		<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 CUMMCL+ Commercial "A" = Results Only Commercial "B" = Results + QC Summary NJ Reduced = Results + QC Summary + Partial Raw data		Comments / Special Instructions: <b>505 524.2 full list</b> <b>V 524SL4+V0, VM5+FI13</b>	
<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"/> 9 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>Sample Custody must be documented below each time samples change possession, including courier delivery.</b>	
Relinquished by Sampler: <b>Pat Proroki</b> Date Time: <b>11/16/15 1900</b>		Received By: <b>Robert deubers</b> Date Time: <b>11-17-15 10:30</b>		Relinquished By: _____ Date Time: _____	
Relinquished by: _____ Date Time: _____		Received By: _____ Date Time: _____		Relinquished By: _____ Date Time: _____	
Relinquished by: _____ Date Time: _____		Received By: _____ Date Time: _____		Relinquished By: _____ Date Time: _____	
Custody Seal # <b>352</b> <input type="checkbox"/> Intact <input type="checkbox"/> Not Intact		Preserved where applicable <input type="checkbox"/>		On Ice <input checked="" type="checkbox"/> Cooler Temp <b>7.0°C</b>	

VOCs 524 Full List  
 B827051M 14 P10X

INITIAL ASSESSMENT **ZVS**  
 LABEL VERIFICATION **09**

5.1  
 5





# Report of Analysis

<b>Client Sample ID:</b> BPOW 5-4		<b>Date Sampled:</b> 11/16/15
<b>Lab Sample ID:</b> JC8685-1		<b>Date Received:</b> 11/17/15
<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	1B100684.D	1	11/19/15	MD	n/a	n/a	V1B4767
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	<del>1.7</del> 5.0	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	<del>0.096</del> 0.50	0.50	0.047	ug/l	J UB
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> BPOW 5-4		<b>Date Sampled:</b> 11/16/15
<b>Lab Sample ID:</b> JC8685-1		<b>Date Received:</b> 11/17/15
<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	99%		78-114%
460-00-4	4-Bromofluorobenzene	101%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
67-63-0	Isopropyl Alcohol	7.65	1.6	ug/l	JN
	Total TIC, Volatile		1.6	ug/l	J N

ND = Not detected      MDL = Method Detection Limit  
 RL = Reporting Limit

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

4.1  
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## Report of Analysis

<b>Client Sample ID:</b> FB111615PP1		
<b>Lab Sample ID:</b> JC8685-2		<b>Date Sampled:</b> 11/16/15
<b>Matrix:</b> AQ - Field Blank Water		<b>Date Received:</b> 11/17/15
<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	1B100685.D	1	11/19/15	MD	n/a	n/a	V1B4767
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	4.0	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	0.15	0.50	0.047	ug/l	J
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> FB111615PP1		<b>Date Sampled:</b> 11/16/15
<b>Lab Sample ID:</b> JC8685-2		<b>Date Received:</b> 11/17/15
<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	99%		78-114%
460-00-4	4-Bromofluorobenzene	101%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	alkene	8.28	24	ug/l	J N
78-84-2	Propanal, 2-methyl-	8.89	.71	ug/l	JN
	Total TIC, Volatile		24.71	ug/l	J N

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

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

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## Report of Analysis

<b>Client Sample ID:</b> TB111615PP1	<b>Date Sampled:</b> 11/16/15
<b>Lab Sample ID:</b> JC8685-3	<b>Date Received:</b> 11/17/15
<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 #1	1B100686.D	1	11/19/15	MD	n/a	n/a	V1B4767
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	1.7	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> TB111615PP1		<b>Date Sampled:</b> 11/16/15
<b>Lab Sample ID:</b> JC8685-3		<b>Date Received:</b> 11/17/15
<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	103%		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 5-4		<b>Date Sampled:</b> 11/16/15
<b>Lab Sample ID:</b> JC8685-1		<b>Date Received:</b> 11/17/15
<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		

Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 <sup>a</sup>	I99179.D	1	11/25/15	AMA	11/18/15	M:OP45466	M:MSI3703
Run #2							

Run #	Initial Volume	Final Volume
Run #1	1000 ml	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
123-91-1	1,4-Dioxane	0.28	0.20	0.076	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	87%		26-121%		
321-60-8	2-Fluorobiphenyl	83%		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

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## Report of Analysis

<b>Client Sample ID:</b> FB111615PP1	<b>Date Sampled:</b> 11/16/15
<b>Lab Sample ID:</b> JC8685-2	<b>Date Received:</b> 11/17/15
<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	

Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 <sup>a</sup>	I99180.D	1	11/25/15	AMA	11/18/15	M:OP45466	M:MSI3703
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	77%		26-121%		
321-60-8	2-Fluorobiphenyl	75%		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

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

### **Data Review**

BETHPAGE, NEW YORK

Volatile and Semivolatile Analyses

SDGs #JC8939, JC9090 and JC9091

Analyses Performed By:  
Accutest Laboratories  
Dayton, New Jersey

Report #24831R  
December 22, 2015  
Review Level: Tier II  
Project #NY001496.1514.NAVI4

## SUMMARY

This data quality assessment summarizes the review of Sample Delivery Groups (SDGs) # JC8939, JC9090 and JC9091 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
JC8939	RE 117D1	JC8939-1	Water	11/18/2015		X	X			
	RE 117D2	JC8939-2	Water	11/18/2015		X	X			
	FB111815PP1	JC8939-3	Water	11/18/2015		X	X			
	TB111815PP1	JC8939-4	Water	11/18/2015		X				
JC9090	BPOW5-7	JC9090-1	Water	11/20/2015		X	X			
	FB112015PP1	JC9090-2	Water	11/20/2015		X	X			
	TB112015PP2	JC9090-3	Water	11/20/2015		X				
JC9091	RE119D1	JC9091-1	Water	11/20/2015		X	X			
	TB112015PP1	JC9091-2	Water	11/20/2015		X				

Notes:

1. Semivolatile (1,4-Dioxane) analysis was performed at Accutest Laboratories, Marlborough, MA (Accutest-NE) facility.
2. SDG JC8939: Matrix spike/matrix spike duplicate analysis was performed on sample location RE 117D2 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 8260C, 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..
SW-846 8260C			

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 not detected above the MDL in the associated blanks; therefore detected sample results were not associated with blank contamination in SDGs JC8939 and JC9091.

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
SDG JC9090:			
BPOW5-7	Acetone	Detected sample results >RL and <BAL	"UB" at the RL
	TIC: Unknown (RT7.60)	Detected sample results less than 5 times blank result	R

RL Reporting limit

## 3. Surrogates/System Monitoring Compounds

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

All surrogate recoveries were within control limits.

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

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

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

A MS/MSD analysis was not performed on a sample location associated with SDGs JC9090 or JC9091.

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

Sample Locations	Compound
<u>SDG JC8939:</u>	
RE 117D2	All compounds , except Freon 113

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 in SDGs JC8939, JC9090 and JC9091.

#### 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 SDGs JC8939, JC9090 or JC9091.

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

Tentatively identified compounds (TICs) were identified in SDG JC9090 in sample locations: BPOW5-7 and FB112015PP1. The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification. TICs are qualified as estimated (JN).

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

## DATA VALIDATION CHECKLIST FOR VOCs

VOCs: 524.2 and 8260C	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 in SDG JC8939.

A MS/MSD analysis was not performed on a sample location associated with SDGs JC9090 and JC9091.

## **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 for SDGs JC8939, JC9090 and JC9091.

## **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: December 22, 2015

PEER REVIEW BY: Todd Church

DATE: December 29, 2015

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

GW  
WFB  
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>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>631-249-7600</b> Fax: <b>631-249-7610</b> Sampler(s) Name(s): <b>Pat Prerato 516-642-5166</b>		<b>Project Information</b> Project Name: <b>AGMNYM62235 // OU2 Monitoring Wells</b> Street: <b>Northrop Grumman OU2 Hydro</b> Billing Information (if different from Report to): Company Name: <b>Bethpage NY</b> Street Address: <b>Arcadis, U.S., Inc. Attn: Accts Payable</b> Client Purchase Order #: <b>NY001496_2015</b> Work Authorization #: <b>NY001496_2015</b> Project Manager: <b>Carlo San Giovanni</b>		<b>Requested Analysis (see TEST CODE sheet)</b> FED-EX Tracking #: <b>#5</b> Accutest Quote #: <b>VC82602NG36GW-40</b> Bottle Order Control #: <b>B8270S1M14D10X</b> Accutest Job #: <b>JC8939</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 - Wipes FB - Field Blank EB - Equipment Blank RB - Rinse Blank TB - Trip Blank
<b>Field ID / Point of Collection</b> MEQHDV1 Vial # Date Time Sampled by Matrix # of bottles HCl NICK MNO3 HRSO4 NH4 DI Water MICH ENCORE		<b>Number of preserved bottles</b> VC82602NG36GW-40 B8270S1M14D10X		<b>LAB USE ONLY</b> V900		
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		<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 CUMMUL+		<b>Comments / Special Instructions</b> RL reporting for metals Please use RE 11702 as a QA/QC MS/MSO sample		
<b>Sample Custody must be documented below each time samples change possession, including courier delivery.</b>						
Requisitioned by: <b>Pat Prerato</b> Date Time: <b>11/18/15 1900</b>	Received By: <b>[Signature]</b> Date Time: <b>11/18/15 17:50</b>	Requisitioned By: <b>[Signature]</b> Date Time: <b>11/18/15 16:11</b>	Received By: <b>[Signature]</b> Date Time: <b>11/18/15 16:11</b>	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>3.300</b>		

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FED-EX Tracking # <b>#5</b>		Bottle Order Control #	
Accutest Quote #		Accutest Job # <b>JC8939</b>	
<b>Client / Reporting Information</b>		<b>Project Information</b>	
Company Name <b>Arcadis</b>		Project Name: <b>AGMNYM62235 // OU2 Monitoring Wells Northrop Grumman OU2 Hydro</b>	
Street Address <b>2 Huntington Quad, Suite 1S10</b>		Billing Information (if different from Report to) Company Name <b>Arcadis, U.S., Inc. Attn: Accts Payable</b>	
City State Zip <b>Melville NY 11747</b>		Street Address <b>630 Plaza Drive, Suite 600</b>	
Project Contact <b>Soma Das, soma.das@arcadis-us.com</b>		City State Zip <b>Highlands Ranch, CO 80129</b>	
Phone # <b>631-249-7600</b>		Work Authorization #: NY001496_2015	
Sample(s) Name(s) <b>RE 117D1, RE 117D2, FB 111815PP2</b>		Project Manager <b>Carlo San Giovanni</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 T/A date 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" <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 LUN/MLL+ Commercial "A" = Results Only Commercial "B" = Results + QC Summary NJ Reduced = Results + QC Summary + Partial Raw data	
Approved By (Accutest PMP) / Date:		RL-reporting for metals <b>Please use RE 117D2 as a QA/QC MS/MSD sample</b>	
Sample Custody must be documented below each time samples change possession, including courier delivery.			
Relinquished by: <b>[Signature]</b>	Date/Time: <b>11/18/15 1830</b>	Received By: <b>FDEX</b>	Date/Time: <b>09:30</b>
Relinquished by: <b>[Signature]</b>	Date/Time:	Received By:	Date/Time: <b>11/19/15</b>
Relinquished by:	Date/Time:	Received By:	Date/Time:
Relinquished by:	Date/Time:	Received By:	Date/Time:
Custody Seal #		On Ice / Cooler Temp. <b>15, 20, 25</b>	

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**JC8939: Chain of Custody**

**Page 3 of 4**

## Report of Analysis

<b>Client Sample ID:</b> RE 117D1		<b>Date Sampled:</b> 11/18/15
<b>Lab Sample ID:</b> JC8939-1		<b>Date Received:</b> 11/19/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8260C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	2D150433.D	1	11/24/15	BK	n/a	n/a	V2D6321
Run #2							

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

## VOA OU2 GW List

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	10	3.3	ug/l	
71-43-2	Benzene	ND	0.50	0.24	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.23	ug/l	
75-25-2	Bromoform	ND	1.0	0.23	ug/l	
74-83-9	Bromomethane	ND	2.0	0.42	ug/l	
78-93-3	2-Butanone (MEK)	ND	10	5.6	ug/l	
75-15-0	Carbon disulfide	ND	2.0	0.25	ug/l	
56-23-5	Carbon tetrachloride	ND	1.0	0.22	ug/l	
108-90-7	Chlorobenzene	ND	1.0	0.19	ug/l	
75-00-3	Chloroethane	ND	1.0	0.34	ug/l	
67-66-3	Chloroform	ND	1.0	0.19	ug/l	
74-87-3	Chloromethane	ND	1.0	0.41	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.15	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.17	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.18	ug/l	
75-35-4	1,1-Dichloroethene	ND	1.0	0.51	ug/l	
156-59-2	cis-1,2-Dichloroethene	ND	1.0	0.27	ug/l	
156-60-5	trans-1,2-Dichloroethene	ND	1.0	0.65	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.39	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.21	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.19	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.27	ug/l	
76-13-1	Freon 113	ND	5.0	0.52	ug/l	
591-78-6	2-Hexanone	ND	5.0	1.7	ug/l	
108-10-1	4-Methyl-2-pentanone(MIBK)	ND	5.0	1.0	ug/l	
75-09-2	Methylene chloride	ND	2.0	0.73	ug/l	
100-42-5	Styrene	ND	1.0	0.27	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	1.0	0.21	ug/l	
127-18-4	Tetrachloroethene	ND	1.0	0.40	ug/l	
108-88-3	Toluene	0.75	1.0	0.16	ug/l	J
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.25	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.21	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> RE 117D1		<b>Date Sampled:</b> 11/18/15
<b>Lab Sample ID:</b> JC8939-1		<b>Date Received:</b> 11/19/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8260C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**VOA OU2 GW List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethene	9.4	1.0	0.22	ug/l	
75-01-4	Vinyl chloride	ND	1.0	0.15	ug/l	
	m,p-Xylene	ND	1.0	0.38	ug/l	
95-47-6	o-Xylene	ND	1.0	0.17	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	102%		76-120%
17060-07-0	1,2-Dichloroethane-D4	100%		73-122%
2037-26-5	Toluene-D8	100%		84-119%
460-00-4	4-Bromofluorobenzene	102%		78-117%

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> RE 117D2		<b>Date Sampled:</b> 11/18/15
<b>Lab Sample ID:</b> JC8939-2		<b>Date Received:</b> 11/19/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8260C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**VOA OU2 GW List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethene	ND	1.0	0.22	ug/l	J
75-01-4	Vinyl chloride	ND	1.0	0.15	ug/l	J
	m,p-Xylene	ND	1.0	0.38	ug/l	J
95-47-6	o-Xylene	ND	1.0	0.17	ug/l	J

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	100%		76-120%
17060-07-0	1,2-Dichloroethane-D4	98%		73-122%
2037-26-5	Toluene-D8	100%		84-119%
460-00-4	4-Bromofluorobenzene	100%		78-117%

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> FB111815PP1		
<b>Lab Sample ID:</b> JC8939-3		<b>Date Sampled:</b> 11/18/15
<b>Matrix:</b> AQ - Field Blank Water		<b>Date Received:</b> 11/19/15
<b>Method:</b> SW846 8260C		<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	2D150331.D	1	11/21/15	BK	n/a	n/a	V2D6317
Run #2							

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

## VOA OU2 GW List

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	10	3.3	ug/l	
71-43-2	Benzene	ND	0.50	0.24	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.23	ug/l	
75-25-2	Bromoform	ND	1.0	0.23	ug/l	
74-83-9	Bromomethane	ND	2.0	0.42	ug/l	
78-93-3	2-Butanone (MEK)	ND	10	5.6	ug/l	
75-15-0	Carbon disulfide	ND	2.0	0.25	ug/l	
56-23-5	Carbon tetrachloride	ND	1.0	0.22	ug/l	
108-90-7	Chlorobenzene	ND	1.0	0.19	ug/l	
75-00-3	Chloroethane	ND	1.0	0.34	ug/l	
67-66-3	Chloroform	ND	1.0	0.19	ug/l	
74-87-3	Chloromethane	ND	1.0	0.41	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.15	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.17	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.18	ug/l	
75-35-4	1,1-Dichloroethene	ND	1.0	0.51	ug/l	
156-59-2	cis-1,2-Dichloroethene	ND	1.0	0.27	ug/l	
156-60-5	trans-1,2-Dichloroethene	ND	1.0	0.65	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.39	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.21	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.19	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.27	ug/l	
76-13-1	Freon 113	ND	5.0	0.52	ug/l	
591-78-6	2-Hexanone	ND	5.0	1.7	ug/l	
108-10-1	4-Methyl-2-pentanone(MIBK)	ND	5.0	1.0	ug/l	
75-09-2	Methylene chloride	ND	2.0	0.73	ug/l	
100-42-5	Styrene	ND	1.0	0.27	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	1.0	0.21	ug/l	
127-18-4	Tetrachloroethene	ND	1.0	0.40	ug/l	
108-88-3	Toluene	ND	1.0	0.16	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.25	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.21	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> FB111815PP1		<b>Date Sampled:</b> 11/18/15
<b>Lab Sample ID:</b> JC8939-3		<b>Date Received:</b> 11/19/15
<b>Matrix:</b> AQ - Field Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8260C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**VOA OU2 GW List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethene	ND	1.0	0.22	ug/l	
75-01-4	Vinyl chloride	ND	1.0	0.15	ug/l	
	m,p-Xylene	ND	1.0	0.38	ug/l	
95-47-6	o-Xylene	ND	1.0	0.17	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	99%		76-120%
17060-07-0	1,2-Dichloroethane-D4	97%		73-122%
2037-26-5	Toluene-D8	99%		84-119%
460-00-4	4-Bromofluorobenzene	101%		78-117%

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> TB111815PP1		<b>Date Sampled:</b> 11/18/15
<b>Lab Sample ID:</b> JC8939-4		<b>Date Received:</b> 11/19/15
<b>Matrix:</b> AQ - Trip Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8260C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	2D150332.D	1	11/21/15	BK	n/a	n/a	V2D6317
Run #2							

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

## VOA OU2 GW List

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	10	3.3	ug/l	
71-43-2	Benzene	ND	0.50	0.24	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.23	ug/l	
75-25-2	Bromoform	ND	1.0	0.23	ug/l	
74-83-9	Bromomethane	ND	2.0	0.42	ug/l	
78-93-3	2-Butanone (MEK)	ND	10	5.6	ug/l	
75-15-0	Carbon disulfide	ND	2.0	0.25	ug/l	
56-23-5	Carbon tetrachloride	ND	1.0	0.22	ug/l	
108-90-7	Chlorobenzene	ND	1.0	0.19	ug/l	
75-00-3	Chloroethane	ND	1.0	0.34	ug/l	
67-66-3	Chloroform	ND	1.0	0.19	ug/l	
74-87-3	Chloromethane	ND	1.0	0.41	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.15	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.17	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.18	ug/l	
75-35-4	1,1-Dichloroethene	ND	1.0	0.51	ug/l	
156-59-2	cis-1,2-Dichloroethene	ND	1.0	0.27	ug/l	
156-60-5	trans-1,2-Dichloroethene	ND	1.0	0.65	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.39	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.21	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.19	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.27	ug/l	
76-13-1	Freon 113	ND	5.0	0.52	ug/l	
591-78-6	2-Hexanone	ND	5.0	1.7	ug/l	
108-10-1	4-Methyl-2-pentanone(MIBK)	ND	5.0	1.0	ug/l	
75-09-2	Methylene chloride	ND	2.0	0.73	ug/l	
100-42-5	Styrene	ND	1.0	0.27	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	1.0	0.21	ug/l	
127-18-4	Tetrachloroethene	ND	1.0	0.40	ug/l	
108-88-3	Toluene	ND	1.0	0.16	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.25	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.21	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> TB111815PP1		<b>Date Sampled:</b> 11/18/15
<b>Lab Sample ID:</b> JC8939-4		<b>Date Received:</b> 11/19/15
<b>Matrix:</b> AQ - Trip Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8260C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**VOA OU2 GW List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethene	ND	1.0	0.22	ug/l	
75-01-4	Vinyl chloride	ND	1.0	0.15	ug/l	
	m,p-Xylene	ND	1.0	0.38	ug/l	
95-47-6	o-Xylene	ND	1.0	0.17	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	100%		76-120%
17060-07-0	1,2-Dichloroethane-D4	97%		73-122%
2037-26-5	Toluene-D8	99%		84-119%
460-00-4	4-Bromofluorobenzene	101%		78-117%

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  
4

## Report of Analysis

<b>Client Sample ID:</b> RE 117D1		<b>Date Sampled:</b> 11/18/15
<b>Lab Sample ID:</b> JC8939-1		<b>Date Received:</b> 11/19/15
<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>	I99203.D	1	11/27/15	AMA	11/22/15	M:OP45518	M:MSI3704
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	77%		26-121%		
321-60-8	2-Fluorobiphenyl	64%		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

4.1  
4



## Report of Analysis

<b>Client Sample ID:</b> RE 117D2		<b>Date Sampled:</b> 11/18/15
<b>Lab Sample ID:</b> JC8939-2		<b>Date Received:</b> 11/19/15
<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>	I99186.D	1	11/26/15	AMA	11/22/15	M:OP45518	M:MSI3703
Run #2							

Run #	Initial Volume	Final Volume
Run #1	1000 ml	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
123-91-1	1,4-Dioxane	ND	0.20	0.076	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	75%		26-121%		
321-60-8	2-Fluorobiphenyl	66%		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> FB111815PP1		<b>Date Sampled:</b> 11/18/15
<b>Lab Sample ID:</b> JC8939-3		<b>Date Received:</b> 11/19/15
<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>	I99204.D	1	11/27/15	AMA	11/22/15	M:OP45518	M:MSI3704
Run #2							

	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	ND	0.21	0.079	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	79%		26-121%		
321-60-8	2-Fluorobiphenyl	68%		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



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>Satellite (if Name(s))</b> V. Prasad 516.877.6477		<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> Arcadis, U.S., Inc. Attn: Accts Payable <b>Project #</b> NY001496.1614.NAVI3 <b>Street Address</b> 630 Plaza Drive, Suite 600 <b>City State Zip</b> Highlands Ranch, CO 80129 <b>Client Purchase Order #</b> <b>Work Authorization #:</b> NY001496_2015 <b>Project Manager</b> Carlo San Giovanni <b>Attention:</b> Soma Das		<b>Accutest Tracking #</b> #4 <b>Accutest Quote #</b> <b>Accutest Job #</b> JC9090		<b>Matrix Codes</b> DW - Drinking Water GW - Ground Water WW - Water SW - Surface Water SO - Soil SL - Sludge SEP - 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 (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+		<b>Comments / Special Instructions</b> VMS 524.2 Full list V524SL4 + 40 VMS + F113	
Sample Custody must be documented below each time sample change possession, including courier delivery.							
<b>Relinquished by Sampler:</b> 1 [Signature] <b>Relinquished by Sampler:</b> 3 <b>Relinquished by:</b> 5	<b>Date Time:</b> 11/20/15 1730 <b>Date Time:</b> <b>Date Time:</b>	<b>Received By:</b> 1 [Signature] <b>Received By:</b> 3 <b>Received By:</b> 5	<b>Date Time:</b> <b>Date Time:</b> <b>Date Time:</b>	<b>Relinquished By:</b> 2 [Signature] <b>Relinquished By:</b> 4 <b>Custody Seal #</b>	<b>Date Time:</b> 11/20/15 <b>Date Time:</b> <b>Date Time:</b>	<b>Received By:</b> 2 [Signature] <b>Received By:</b> 4 <b>Received By:</b>	<b>Preserved where applicable:</b> <input type="checkbox"/> Exact <input type="checkbox"/> Not Exact Preserved where applicable: 2x 600ml On Ice: <input type="checkbox"/> 2.0°C

5.1  
5

17B

# Report of Analysis

<b>Client Sample ID:</b> BPOW5-7		<b>Date Sampled:</b> 11/20/15
<b>Lab Sample ID:</b> JC9090-1		<b>Date Received:</b> 11/20/15
<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	1B100758.D	1	11/24/15	MD	n/a	n/a	V1B4770
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	5.0	1.1	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	0.49	0.50	0.044	ug/l	J

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> BPOW5-7		<b>Date Sampled:</b> 11/20/15
<b>Lab Sample ID:</b> JC9090-1		<b>Date Received:</b> 11/20/15
<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	102%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	unknown	7.60	<del>.65</del>	ug/l	<del>J</del> R
	Total TIC, Volatile		<del>.65</del>	ug/l	<del>J</del>

ND = Not detected      MDL = Method Detection Limit  
 RL = Reporting Limit

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> FB112015PP1		
<b>Lab Sample ID:</b> JC9090-2		<b>Date Sampled:</b> 11/20/15
<b>Matrix:</b> AQ - Field Blank Water		<b>Date Received:</b> 11/20/15
<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	1B100759.D	1	11/24/15	MD	n/a	n/a	V1B4770
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	0.20	0.50	0.047	ug/l	J
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> FB112015PP1	<b>Date Sampled:</b> 11/20/15
<b>Lab Sample ID:</b> JC9090-2	<b>Date Received:</b> 11/20/15
<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	102%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	unknown	7.60	.59	ug/l	J N
	Total TIC, Volatile		.59	ug/l	J N

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

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





## Report of Analysis

<b>Client Sample ID:</b> TB112015PP2		<b>Date Sampled:</b> 11/20/15
<b>Lab Sample ID:</b> JC9090-3		<b>Date Received:</b> 11/20/15
<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	103%		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> BPOW5-7		<b>Date Sampled:</b> 11/20/15
<b>Lab Sample ID:</b> JC9090-1		<b>Date Received:</b> 11/20/15
<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>	I99205.D	1	11/27/15	AMA	11/22/15	M:OP45518	M:MSI3704
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	69%		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> FB112015PP1	<b>Date Sampled:</b> 11/20/15
<b>Lab Sample ID:</b> JC9090-2	<b>Date Received:</b> 11/20/15
<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>	I99206.D	1	11/27/15	AMA	11/22/15	M:OP45518	M:MSI3704
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	ND	0.22	0.083	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	82%		26-121%		
321-60-8	2-Fluorobiphenyl	74%		28-107%		
1718-51-0	Terphenyl-d14	87%		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> RE119D1		<b>Date Sampled:</b> 11/20/15
<b>Lab Sample ID:</b> JC9091-1		<b>Date Received:</b> 11/20/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8260C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	2B137063.D	1	12/01/15	EH	n/a	n/a	V2B6106
Run #2							

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

## VOA OU2 GW List

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	10	3.3	ug/l	
71-43-2	Benzene	ND	0.50	0.24	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.23	ug/l	
75-25-2	Bromoform	ND	1.0	0.23	ug/l	
74-83-9	Bromomethane	ND	2.0	0.42	ug/l	
78-93-3	2-Butanone (MEK)	ND	10	5.6	ug/l	
75-15-0	Carbon disulfide	ND	2.0	0.25	ug/l	
56-23-5	Carbon tetrachloride	ND	1.0	0.22	ug/l	
108-90-7	Chlorobenzene	ND	1.0	0.19	ug/l	
75-00-3	Chloroethane	ND	1.0	0.34	ug/l	
67-66-3	Chloroform	ND	1.0	0.19	ug/l	
74-87-3	Chloromethane	ND	1.0	0.41	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.15	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.17	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.18	ug/l	
75-35-4	1,1-Dichloroethene	ND	1.0	0.51	ug/l	
156-59-2	cis-1,2-Dichloroethene	ND	1.0	0.27	ug/l	
156-60-5	trans-1,2-Dichloroethene	ND	1.0	0.65	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.39	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.21	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.19	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.27	ug/l	
76-13-1	Freon 113	ND	5.0	0.52	ug/l	
591-78-6	2-Hexanone	ND	5.0	1.7	ug/l	
108-10-1	4-Methyl-2-pentanone(MIBK)	ND	5.0	1.0	ug/l	
75-09-2	Methylene chloride	ND	2.0	0.73	ug/l	
100-42-5	Styrene	ND	1.0	0.27	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	1.0	0.21	ug/l	
127-18-4	Tetrachloroethene	ND	1.0	0.40	ug/l	
108-88-3	Toluene	0.72	1.0	0.16	ug/l	J
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.25	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.21	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> RE119D1		<b>Date Sampled:</b> 11/20/15
<b>Lab Sample ID:</b> JC9091-1		<b>Date Received:</b> 11/20/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8260C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**VOA OU2 GW List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethene	ND	1.0	0.22	ug/l	
75-01-4	Vinyl chloride	ND	1.0	0.15	ug/l	
	m,p-Xylene	ND	1.0	0.38	ug/l	
95-47-6	o-Xylene	ND	1.0	0.17	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	105%		76-120%
17060-07-0	1,2-Dichloroethane-D4	115%		73-122%
2037-26-5	Toluene-D8	100%		84-119%
460-00-4	4-Bromofluorobenzene	100%		78-117%

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> TB112015PP1		<b>Date Sampled:</b> 11/20/15
<b>Lab Sample ID:</b> JC9091-2		<b>Date Received:</b> 11/20/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8260C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**VOA OU2 GW List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethene	ND	1.0	0.22	ug/l	
75-01-4	Vinyl chloride	ND	1.0	0.15	ug/l	
	m,p-Xylene	ND	1.0	0.38	ug/l	
95-47-6	o-Xylene	ND	1.0	0.17	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	102%		76-120%
17060-07-0	1,2-Dichloroethane-D4	103%		73-122%
2037-26-5	Toluene-D8	100%		84-119%
460-00-4	4-Bromofluorobenzene	102%		78-117%

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> RE119D1	<b>Date Sampled:</b> 11/20/15
<b>Lab Sample ID:</b> JC9091-1	<b>Date Received:</b> 11/20/15
<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>	I99207.D	1	11/27/15	AMA	11/22/15	M:OP45518	M:MSI3704
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	74%		26-121%		
321-60-8	2-Fluorobiphenyl	67%		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

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

### **Data Review**

BETHPAGE, NEW YORK

Volatile and Semivolatile Analyses

SDGs #JC9220, JC9566 and JC9689

Analyses Performed By:  
Accutest Laboratories  
Dayton, New Jersey

Report #24832R  
December 28, 2015  
Review Level: Tier II  
Project #NY001496.1514.NAVI4

## SUMMARY

This data quality assessment summarizes the review of Sample Delivery Groups (SDGs) # JC9220, JC9566 and JC9689 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
JC9220	FB112315PP1	JC9220-1	Water	11/23/2015		X	X			
	TB112315PP1	JC9220-2	Water	11/23/2015		X				
	RE118D1	JC9220-3	Water	11/23/2015		X	X			
JC9566	FB113015PP1	JC9566-1	Water	11/30/2015		X	X			
	BPOW 6-1	JC9566-2	Water	11/30/2015		X	X			
	BPOW 6-2	JC9566-3	Water	11/30/2015		X	X			
	TB113015PP1	JC9566-4	Water	11/30/2015		X				
JC9689	FB120115PP1	JC9689-1	Water	12/01/2015		X	X			
	BPOW 6-3	JC9689-2	Water	12/01/2015		X	X			
	BPOW 6-4	JC9689-3	Water	12/01/2015		X	X			
	REP120115PP1	JC9689-4	Water	12/01/2015	BPOW 6-4	X	X			
	TB120115PP1	JC9689-5	Water	12/01/2015		X				

Note:

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 8260C, 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..
SW-846 8260C			

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 not detected above the MDL in the associated blanks; therefore detected sample results were not associated with blank contamination in SDG JC9220.

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 in SDG JC9566.

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
SDG JC9689:			
BPOW 6-3	Acetone	Detected sample results >RL and <BAL	"UB" at the RL

RL Reporting limit

## 3. Surrogates/System Monitoring Compounds

All samples to be analyzed for organic compounds are spiked with surrogate compounds prior to sample preparation to evaluate overall laboratory performance and efficiency of the analytical technique. VOC



analysis requires that all surrogates associated with the analysis exhibit recoveries within the laboratory-established acceptance limits.

All surrogate recoveries were within control limits.

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

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

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

A MS/MSD analysis was not performed on a sample location associated with SDGs JC9220, JC9566 or JC9689.

#### 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 SDGs JC9220 or JC9566.

Results for duplicate samples are summarized in the following table.

Sample ID/Duplicate ID	Compound	Sample Result	Duplicate Result	RPD
SDG JC9689:				
BPOW 6-4/ REP120115PP1	All compounds	U	U	AC

AC Acceptable

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

#### 7. Laboratory Duplicate Analysis

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

concentrations are less than or equal to 5 times the RL, a control limit of two times the RL is applied for water matrices.

A laboratory duplicate sample was not performed on a sample location associated with SDGs JC9220, JC9566 or JC9689.

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

Tentatively identified compounds (TICs) were not identified in SDG JC9220. TICs were identified in SDG JC9566 in sample locations FB113015PP1; and, in SDG JC9689 in sample location FB12015PP1. The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification. TICs are qualified as estimated (JN).

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

## DATA VALIDATION CHECKLIST FOR VOCs

VOCs: 524.2 and 8260C	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		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 SDGs JC9220, JC9566 or JC9689.

## 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 SDG JC9220 and JC9566

Results for duplicate samples are summarized in the following table.

Sample ID/Duplicate ID	Compound	Sample Result	Duplicate Result	RPD
SDG JC9689:				
BPOW 6-4/ REP120115PP1	1,4-Dioxane	0.22 U	0.21 U	AC

AC Acceptable

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

## 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		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: December 28, 2015

PEER REVIEW BY: Todd Church

DATE: December 29, 2015

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



GW  
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

FED-EX Tracking # **#5**      Bottle Order Control #  
Accutest Quote # **JC9220**

Client / Reporting Information		Project Information				Requested Analysis ( see TEST CODE sheet)												Matrix Codes			
Company Name <b>Arcadis</b>		Project Name: <b>AGMNYM62235 // OU2 Monitoring Wells Northrop Grumman OU2 Hydro</b>				<div style="display: flex; justify-content: space-between;"> <div style="writing-mode: vertical-rl; transform: rotate(180deg);">VC82602NG36GW-40</div> <div style="writing-mode: vertical-rl; transform: rotate(180deg);">B8270SIM14DIOX</div> </div>												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		Billing Information ( if different from Report to)																	
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>																	
Project Contact <b>Soma Das, soma.das@arcadis-us.com</b>		Project # <b>NY001496.2015</b>		Street Address <b>630 Plaza Drive, Suite 600</b>																	
Phone # Fax # <b>631-249-7600 631-249-7810</b>		Client Purchase Order # <b>NY001496_2015</b>		City State Zip <b>Highlands Ranch, CO 80129</b>																	
Sampler(s) Name(s) <b>Pat Pruski 567-6477</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 <small>Emergency &amp; Rush T/A data available VIA Lablink</small>		Approved By (Accutest PM) / Date: _____ _____				<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 U/MML+				RL reporting for metals _____ _____											
Requisitioned by Sampler: <b>Carlo San Giovanni</b>		Date Time: <b>11/23/15 1830</b>		Received By: <b>[Signature]</b>				Requisitioned By: <b>[Signature]</b>				Date Time: <b>11/24/15 40</b>				Received By: <b>[Signature]</b>					
Requisitioned by: <b>3</b>		Date Time: <b>3</b>		Received By: <b>3</b>				Requisitioned By: <b>4</b>				Date Time: <b>4</b>				Received By: <b>4</b>					
Requisitioned by: <b>5</b>		Date Time: <b>5</b>		Received By: <b>5</b>				Requisitioned By: <b>5</b>				Date Time: <b>5</b>				Received By: <b>5</b>					
				Custody Seal # _____				<input type="checkbox"/> Intact      Preserved where applicable <input type="checkbox"/> Not Intact				On Ice <input type="checkbox"/> Cooler Temp. <b>0.9 C</b>									

5.1  
5





## Report of Analysis

<b>Client Sample ID:</b> FB112315PP1		<b>Date Sampled:</b> 11/23/15
<b>Lab Sample ID:</b> JC9220-1		<b>Date Received:</b> 11/24/15
<b>Matrix:</b> AQ - Field Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8260C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**VOA OU2 GW List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethene	ND	1.0	0.22	ug/l	
75-01-4	Vinyl chloride	ND	1.0	0.15	ug/l	
	m,p-Xylene	ND	1.0	0.38	ug/l	
95-47-6	o-Xylene	ND	1.0	0.17	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	100%		76-120%
17060-07-0	1,2-Dichloroethane-D4	97%		73-122%
2037-26-5	Toluene-D8	99%		84-119%
460-00-4	4-Bromofluorobenzene	100%		78-117%

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> TB112315PP1		<b>Date Sampled:</b> 11/23/15
<b>Lab Sample ID:</b> JC9220-2		<b>Date Received:</b> 11/24/15
<b>Matrix:</b> AQ - Trip Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8260C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1A156465.D	1	11/30/15	VC	n/a	n/a	V1A6717
Run #2							

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

## VOA OU2 GW List

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	10	3.3	ug/l	
71-43-2	Benzene	ND	0.50	0.24	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.23	ug/l	
75-25-2	Bromoform	ND	1.0	0.23	ug/l	
74-83-9	Bromomethane	ND	2.0	0.42	ug/l	
78-93-3	2-Butanone (MEK)	ND	10	5.6	ug/l	
75-15-0	Carbon disulfide	ND	2.0	0.25	ug/l	
56-23-5	Carbon tetrachloride	ND	1.0	0.22	ug/l	
108-90-7	Chlorobenzene	ND	1.0	0.19	ug/l	
75-00-3	Chloroethane	ND	1.0	0.34	ug/l	
67-66-3	Chloroform	ND	1.0	0.19	ug/l	
74-87-3	Chloromethane	ND	1.0	0.41	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.15	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.17	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.18	ug/l	
75-35-4	1,1-Dichloroethene	ND	1.0	0.51	ug/l	
156-59-2	cis-1,2-Dichloroethene	ND	1.0	0.27	ug/l	
156-60-5	trans-1,2-Dichloroethene	ND	1.0	0.65	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.39	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.21	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.19	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.27	ug/l	
76-13-1	Freon 113	ND	5.0	0.52	ug/l	
591-78-6	2-Hexanone	ND	5.0	1.7	ug/l	
108-10-1	4-Methyl-2-pentanone(MIBK)	ND	5.0	1.0	ug/l	
75-09-2	Methylene chloride	ND	2.0	0.73	ug/l	
100-42-5	Styrene	ND	1.0	0.27	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	1.0	0.21	ug/l	
127-18-4	Tetrachloroethene	ND	1.0	0.40	ug/l	
108-88-3	Toluene	ND	1.0	0.16	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.25	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.21	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> TB112315PP1		<b>Date Sampled:</b> 11/23/15
<b>Lab Sample ID:</b> JC9220-2		<b>Date Received:</b> 11/24/15
<b>Matrix:</b> AQ - Trip Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8260C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**VOA OU2 GW List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethene	ND	1.0	0.22	ug/l	
75-01-4	Vinyl chloride	ND	1.0	0.15	ug/l	
	m,p-Xylene	ND	1.0	0.38	ug/l	
95-47-6	o-Xylene	ND	1.0	0.17	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	100%		76-120%
17060-07-0	1,2-Dichloroethane-D4	97%		73-122%
2037-26-5	Toluene-D8	100%		84-119%
460-00-4	4-Bromofluorobenzene	99%		78-117%

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> RE 118D1		<b>Date Sampled:</b> 11/23/15
<b>Lab Sample ID:</b> JC9220-3		<b>Date Received:</b> 11/24/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8260C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1A156488.D	1	12/01/15	VC	n/a	n/a	V1A6718
Run #2							

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

### VOA OU2 GW List

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	10	3.3	ug/l	
71-43-2	Benzene	ND	0.50	0.24	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.23	ug/l	
75-25-2	Bromoform	ND	1.0	0.23	ug/l	
74-83-9	Bromomethane	ND	2.0	0.42	ug/l	
78-93-3	2-Butanone (MEK)	ND	10	5.6	ug/l	
75-15-0	Carbon disulfide	ND	2.0	0.25	ug/l	
56-23-5	Carbon tetrachloride	ND	1.0	0.22	ug/l	
108-90-7	Chlorobenzene	ND	1.0	0.19	ug/l	
75-00-3	Chloroethane	ND	1.0	0.34	ug/l	
67-66-3	Chloroform	ND	1.0	0.19	ug/l	
74-87-3	Chloromethane	ND	1.0	0.41	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.15	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.17	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.18	ug/l	
75-35-4	1,1-Dichloroethene	ND	1.0	0.51	ug/l	
156-59-2	cis-1,2-Dichloroethene	ND	1.0	0.27	ug/l	
156-60-5	trans-1,2-Dichloroethene	ND	1.0	0.65	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.39	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.21	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.19	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.27	ug/l	
76-13-1	Freon 113	ND	5.0	0.52	ug/l	
591-78-6	2-Hexanone	ND	5.0	1.7	ug/l	
108-10-1	4-Methyl-2-pentanone(MIBK)	ND	5.0	1.0	ug/l	
75-09-2	Methylene chloride	ND	2.0	0.73	ug/l	
100-42-5	Styrene	ND	1.0	0.27	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	1.0	0.21	ug/l	
127-18-4	Tetrachloroethene	ND	1.0	0.40	ug/l	
108-88-3	Toluene	0.57	1.0	0.16	ug/l	J
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.25	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.21	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> RE 118D1		<b>Date Sampled:</b> 11/23/15
<b>Lab Sample ID:</b> JC9220-3		<b>Date Received:</b> 11/24/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8260C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**VOA OU2 GW List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethene	ND	1.0	0.22	ug/l	
75-01-4	Vinyl chloride	ND	1.0	0.15	ug/l	
	m,p-Xylene	ND	1.0	0.38	ug/l	
95-47-6	o-Xylene	ND	1.0	0.17	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	98%		76-120%
17060-07-0	1,2-Dichloroethane-D4	97%		73-122%
2037-26-5	Toluene-D8	100%		84-119%
460-00-4	4-Bromofluorobenzene	101%		78-117%

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> FB112315PP1		<b>Date Sampled:</b> 11/23/15
<b>Lab Sample ID:</b> JC9220-1		<b>Date Received:</b> 11/24/15
<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>	I99247.D	1	12/01/15	AMA	11/24/15	M:OP45543	M:MSI3707
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	ND	0.21	0.079	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	88%		26-121%		
321-60-8	2-Fluorobiphenyl	75%		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

4.1  
4

## Report of Analysis

<b>Client Sample ID:</b> RE 118D1		<b>Date Sampled:</b> 11/23/15
<b>Lab Sample ID:</b> JC9220-3		<b>Date Received:</b> 11/24/15
<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>	I99248.D	1	12/01/15	AMA	11/24/15	M:OP45543	M:MSI3707
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	ND	0.22	0.083	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	67%		26-121%		
321-60-8	2-Fluorobiphenyl	60%		28-107%		
1718-51-0	Terphenyl-d14	87%		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

FED-EX Tracking # **#5** Bottle Order Control # **5C9566**  
Accutest Quote # \_\_\_\_\_

Client / Reporting Information		Project Information		Requested Analysis (see TEST CODE sheet)										Matrix Codes		
Company Name <b>Arcadis</b>		Project Name <b>AGMNYM62235 // OU2 Monitoring Wells</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>Northrop Grumman OU2 Hydro</b>														
City State Zip <b>Melville NY 11747</b>		City State <b>Bethpage NY</b>		Billing Information (If different from Report to)												
Project Contact <b>Soma Das, soma.das@arcadis-us.com</b>		Project # <b>NY001946.1514.NAVIS</b>		Company Name <b>Arcadis, U.S., Inc. Attn: Accts Payable</b>												
Phone # <b>631-249-7600</b>		Client Purchase Order #		Street Address <b>630 Plaza Drive, Suite 600</b>										LAB USE ONLY  <b>V988</b>		
Samples (Name(s)) <b>V1020701</b>		Work Authorization #: NY001496_2015		City State Zip <b>Highlands Ranch, CO 80129</b>												
Phone # <b>516 237-6247</b>		Project Manager <b>Carlo San Giovanni</b>		Attention: <b>Soma Das</b>												
Account Sample #	Field ID / Point of Collection	MECHDI Val #	Collection		Number of preserved bottles											
			Date	Time	Sampled by	Matrix	# of bottles	HC	HNOC	HRBC	HNOC	HRBC	NONE	ED Water	MESH	ENDORE
1	FB113015RPA		11/30/15	1310	AP	FB	3	3								
2	BOWL G-1		11/30/15	1635	AP	GH	3	3								
3	BOWL G-2		11/30/15	1635	AP	GH	3	3								
4	TB113015RPA		11/30/15	1300	-	TB	2	2								

VC62602NG6666WY740  
 B82705IM14DIOX  
 VDCS 524.2 Full list

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  
 Emergency & Rush T/A data available VIA Lablink

Approved By (Accutest PM): / Date: \_\_\_\_\_  
**INITIAL ASSESSMENT - NL3A**  
**LABEL VERIFICATION - COB**

Data Deliverable Information  
 Commercial "A" (Level 1)  
 Commercial "B" (Level 2)  
 FULLT (Level 3+4)  
 NJ Reduced  
 Commercial "C"  
 NYASP Category A  
 NYASP Category B  
 State Forms  
 EDD Format  
 Other LUMMLC+

Commercial "A" = Results Only  
 Commercial "B" = Results + QC Summary  
 NJ Reduced = Results + QC Summary + Partial Raw data

RL reporting for metals  
**VDCS 524.2 Full list**  
**V5245L4+40 VMS + F113**

Sample Custody must be documented below each time samples change possession, including courier delivery.

Relinquished by: <b>[Signature]</b>	Date Time: <b>11/30/15 2:30</b>	Received By: <b>Robert Chambers</b>	Date Time: <b>12-1-15 10:15</b>	Relinquished By: <b>[Signature]</b>	Date Time: <b>12-1-15</b>	Received By: <b>[Signature]</b>
Relinquished by: <b>[Signature]</b>	Date Time: <b>3</b>	Received By: <b>3</b>	Date Time: <b>4</b>	Relinquished By: <b>4</b>	Date Time: <b>4</b>	Received By: <b>4</b>

Custody Seal # **364**  
 Intact  
 Not Intact  
 Preserved where applicable  
 On Ice  Cooler Temp: **2.0°C**

5.1 5



## Report of Analysis

<b>Client Sample ID:</b> FB113015PP1		
<b>Lab Sample ID:</b> JC9566-1		<b>Date Sampled:</b> 11/30/15
<b>Matrix:</b> AQ - Field Blank Water		<b>Date Received:</b> 12/01/15
<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	1B100892.D	1	12/03/15	MD	n/a	n/a	V1B4776
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	2.2	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> FB113015PP1		<b>Date Sampled:</b> 11/30/15
<b>Lab Sample ID:</b> JC9566-1		<b>Date Received:</b> 12/01/15
<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	104%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	unknown	7.61	.65	ug/l	J N
	Total TIC, Volatile		.65	ug/l	J N

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

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

4.1  
4

## Report of Analysis

<b>Client Sample ID:</b>	BPOW 6-1	<b>Date Sampled:</b>	11/30/15
<b>Lab Sample ID:</b>	JC9566-2	<b>Date Received:</b>	12/01/15
<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		

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B100893.D	1	12/03/15	MD	n/a	n/a	V1B4776
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 6-1		<b>Date Sampled:</b> 11/30/15
<b>Lab Sample ID:</b> JC9566-2		<b>Date Received:</b> 12/01/15
<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	101%		78-114%
460-00-4	4-Bromofluorobenzene	102%		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> 11/30/15
<b>Lab Sample ID:</b> JC9566-3		<b>Date Received:</b> 12/01/15
<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	103%		78-114%
460-00-4	4-Bromofluorobenzene	104%		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> TB113015PP1		<b>Date Sampled:</b> 11/30/15
<b>Lab Sample ID:</b> JC9566-4		<b>Date Received:</b> 12/01/15
<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		

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B100895.D	1	12/03/15	MD	n/a	n/a	V1B4776
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> TB113015PP1		<b>Date Sampled:</b> 11/30/15
<b>Lab Sample ID:</b> JC9566-4		<b>Date Received:</b> 12/01/15
<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	102%		78-114%
460-00-4	4-Bromofluorobenzene	104%		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  
4

## Report of Analysis

<b>Client Sample ID:</b> FB113015PP1		<b>Date Sampled:</b> 11/30/15
<b>Lab Sample ID:</b> JC9566-1		<b>Date Received:</b> 12/01/15
<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>	I99338.D	1	12/10/15	AMA	12/02/15	M:OP45603	M:MSI3713
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	67%		26-121%		
321-60-8	2-Fluorobiphenyl	61%		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

4.1  
4

## Report of Analysis

<b>Client Sample ID:</b> BPOW 6-1	
<b>Lab Sample ID:</b> JC9566-2	<b>Date Sampled:</b> 11/30/15
<b>Matrix:</b> AQ - Ground Water	<b>Date Received:</b> 12/01/15
<b>Method:</b> SW846 8270D BY SIM SW846 3510C	<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 <sup>a</sup>	I99339.D	1	12/10/15	AMA	12/02/15	M:OP45603	M:MSI3713
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	72%		26-121%		
321-60-8	2-Fluorobiphenyl	64%		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

4.2  
4

## Report of Analysis

<b>Client Sample ID:</b> BPOW 6-2	<b>Date Sampled:</b> 11/30/15
<b>Lab Sample ID:</b> JC9566-3	<b>Date Received:</b> 12/01/15
<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>	I99340.D	1	12/10/15	AMA	12/02/15	M:OP45603	M:MSI3713
Run #2							

Run #	Initial Volume	Final Volume
Run #1	890 ml	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
123-91-1	1,4-Dioxane	ND	0.22	0.085	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	80%		26-121%		
321-60-8	2-Fluorobiphenyl	71%		28-107%		
1718-51-0	Terphenyl-d14	76%		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







**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 <b>Arcadis</b> 2 Huntington Quad, Suite 1S10 Melville NY 11747 Soma Das, soma.das@arcadis-us.com 631-249-7600 / 631-249-7610		Project Information <b>AGMNYM62235 // OU2 Monitoring Wells Northrop Grumman OU2 Hydro</b> Bethpage NY Arcadis, U.S., Inc. Attn: Accts Payable 630 Plaza Drive, Suite 600 Highlands Ranch, CO 80129 Carlo San Giovanni		Requested Analysis (see TEST CODE sheet) B8270 SIM14DIOX		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 WIP - Wipe FB-Field Blank EB-Equipment Blank RB-Rinsate 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"/> 5 Day RUSH <input type="checkbox"/> 3 Day EMERGENCY <input type="checkbox"/> 2 Day EMERGENCY <input type="checkbox"/> 1 Day EMERGENCY		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 LUM/MUC+		Comments / Special Instructions RL reporting for metals			
Sample Custody must be documented below each time samples change possession, including courier delivery.							
Relinquished By: [Signature] Date Time: 12/1/15 1810		Received By: [Signature] Date Time: 12-2-15 1045		Relinquished By: [Signature] Date Time: [Blank]		Received By: [Signature] Date Time: [Blank]	
Relinquished by: [Signature] Date Time: [Blank]		Received By: [Signature] Date Time: [Blank]		Relinquished By: [Signature] Date Time: [Blank]		Received By: [Signature] Date Time: [Blank]	

5.1  
5





## Report of Analysis

<b>Client Sample ID:</b> FB120115PP1		<b>Date Sampled:</b> 12/01/15
<b>Lab Sample ID:</b> JC9689-1		<b>Date Received:</b> 12/02/15
<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	104%		78-114%
460-00-4	4-Bromofluorobenzene	105%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	alkene	8.25	2	ug/l	J N
	Total TIC, Volatile		2	ug/l	J N

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

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

4.1  
4



## Report of Analysis

<b>Client Sample ID:</b> BPOW 6-3		<b>Date Sampled:</b> 12/01/15
<b>Lab Sample ID:</b> JC9689-2		<b>Date Received:</b> 12/02/15
<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	102%		78-114%
460-00-4	4-Bromofluorobenzene	103%		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 6-4		<b>Date Sampled:</b> 12/01/15
<b>Lab Sample ID:</b> JC9689-3		<b>Date Received:</b> 12/02/15
<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 #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B100887.D	1	12/03/15	MD	n/a	n/a	V1B4776
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

4.3  
 4

## Report of Analysis

<b>Client Sample ID:</b> BPOW 6-4		<b>Date Sampled:</b> 12/01/15
<b>Lab Sample ID:</b> JC9689-3		<b>Date Received:</b> 12/02/15
<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	102%		78-114%
460-00-4	4-Bromofluorobenzene	103%		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> REP120115PP1		<b>Date Sampled:</b> 12/01/15
<b>Lab Sample ID:</b> JC9689-4		<b>Date Received:</b> 12/02/15
<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	1B100885.D	1	12/03/15	MD	n/a	n/a	V1B4776
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

4.4  
4



## Report of Analysis

<b>Client Sample ID:</b> REP120115PP1		<b>Date Sampled:</b> 12/01/15
<b>Lab Sample ID:</b> JC9689-4		<b>Date Received:</b> 12/02/15
<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	103%		78-114%
460-00-4	4-Bromofluorobenzene	103%		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  
4



## Report of Analysis

<b>Client Sample ID:</b> TB120115PP1		<b>Date Sampled:</b> 12/01/15
<b>Lab Sample ID:</b> JC9689-5		<b>Date Received:</b> 12/02/15
<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	102%		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

4.5  
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## Report of Analysis

<b>Client Sample ID:</b> FB120115PP1	<b>Date Sampled:</b> 12/01/15
<b>Lab Sample ID:</b> JC9689-1	<b>Date Received:</b> 12/02/15
<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>	I99290.D	1	12/04/15	AMA	12/03/15	M:OP45613	M:MSI3710
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	74%		26-121%		
321-60-8	2-Fluorobiphenyl	74%		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

4.1  
4

## Report of Analysis

<b>Client Sample ID:</b> BPOW 6-3	<b>Date Sampled:</b> 12/01/15
<b>Lab Sample ID:</b> JC9689-2	<b>Date Received:</b> 12/02/15
<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>	I99291.D	1	12/04/15	AMA	12/03/15	M:OP45613	M:MSI3710
Run #2							

Run #	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	83%		26-121%		
321-60-8	2-Fluorobiphenyl	78%		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  
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## Report of Analysis

<b>Client Sample ID:</b> BPOW 6-4	<b>Date Sampled:</b> 12/01/15
<b>Lab Sample ID:</b> JC9689-3	<b>Date Received:</b> 12/02/15
<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>	I99292.D	1	12/04/15	AMA	12/03/15	M:OP45613	M:MSI3710
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	80%		26-121%		
321-60-8	2-Fluorobiphenyl	76%		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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## Report of Analysis

<b>Client Sample ID:</b> REP120115PP1	<b>Date Sampled:</b> 12/01/15
<b>Lab Sample ID:</b> JC9689-4	<b>Date Received:</b> 12/02/15
<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>	I99293.D	1	12/04/15	AMA	12/03/15	M:OP45613	M:MSI3710
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	76%		28-107%		
1718-51-0	Terphenyl-d14	87%		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  
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## **Northrop Grumman Corporation- Operable Unit 2**

### **Data Review**

BETHPAGE, NEW YORK

Volatile and Semivolatile Analyses

SDGs #JC9792 and JC9923

Analyses Performed By:  
Accutest Laboratories  
Dayton, New Jersey

Report #24833R  
December 28, 2015  
Review Level: Tier II  
Project #NY001496.1514.NAVI4



## SUMMARY

This data quality assessment summarizes the review of Sample Delivery Groups (SDGs) # JC9792 and JC9923 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
JC9792	TB120215PP1	JC9792-1	Water	12/02/2015		X				
	FB120215PP1	JC9792-2	Water	12/02/2015		X	X			
	BPOW 6-5	JC9792-3	Water	12/02/2015		X	X			
	BPOW 6-6	JC9792-4	Water	12/02/2015		X	X			
JC9923	BPOW5-3	JC9923-1	Water	12/03/2015		X	X			
	FB120315PP1	JC9923-2	Water	12/03/2015		X	X			
	TB120315PP1	JC9923-3	Water	12/03/2015		X				

Note:

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 in SDGs JC9792 or JC9923.

### 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 SDGs JC9792 and JC9923.

## **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 SDGs JC9792 or JC9923.

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

Tentatively identified compounds (TICs) were identified in SDG JC9792 in sample location FB120215PP1. The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification. TICs are qualified as estimated (JN).

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

## DATA VALIDATION CHECKLIST FOR VOCs

VOCs: 524.2	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 SDGs JC9792 or JC9923.

## **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: December 28, 2015

PEER REVIEW BY: Todd Church

DATE: December 29, 2015

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





## Report of Analysis

<b>Client Sample ID:</b> TB120215PP1		<b>Date Sampled:</b> 12/02/15
<b>Lab Sample ID:</b> JC9792-1		<b>Date Received:</b> 12/03/15
<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	1B100904.D	1	12/04/15	MD	n/a	n/a	V1B4777
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> TB120215PP1		<b>Date Sampled:</b> 12/02/15
<b>Lab Sample ID:</b> JC9792-1		<b>Date Received:</b> 12/03/15
<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	102%		78-114%
460-00-4	4-Bromofluorobenzene	103%		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> FB120215PP1		
<b>Lab Sample ID:</b> JC9792-2		<b>Date Sampled:</b> 12/02/15
<b>Matrix:</b> AQ - Field Blank Water		<b>Date Received:</b> 12/03/15
<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	1B100905.D	1	12/04/15	MD	n/a	n/a	V1B4777
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	2.1	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> FB120215PP1		<b>Date Sampled:</b> 12/02/15
<b>Lab Sample ID:</b> JC9792-2		<b>Date Received:</b> 12/03/15
<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	101%		78-114%
460-00-4	4-Bromofluorobenzene	102%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	alkene	8.24	.81	ug/l	J N
	Total TIC, Volatile		.81	ug/l	J N

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

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

4.2  
4



## Report of Analysis

<b>Client Sample ID:</b> BPOW 6-5		<b>Date Sampled:</b> 12/02/15
<b>Lab Sample ID:</b> JC9792-3		<b>Date Received:</b> 12/03/15
<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	101%		78-114%
460-00-4	4-Bromofluorobenzene	102%		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 6-6	<b>Date Sampled:</b>	12/02/15
<b>Lab Sample ID:</b>	JC9792-4	<b>Date Received:</b>	12/03/15
<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	1B100907.D	1	12/04/15	MD	n/a	n/a	V1B4777
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	0.40	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

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 6-6		<b>Date Sampled:</b> 12/02/15
<b>Lab Sample ID:</b> JC9792-4		<b>Date Received:</b> 12/03/15
<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	103%		78-114%
460-00-4	4-Bromofluorobenzene	104%		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  
4

## Report of Analysis

<b>Client Sample ID:</b> FB120215PP1		<b>Date Sampled:</b> 12/02/15
<b>Lab Sample ID:</b> JC9792-2		<b>Date Received:</b> 12/03/15
<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>	I99455.D	1	12/16/15	AMA	12/05/15	M:OP45645	M:MSI3717
Run #2							

Run #	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	66%		26-121%		
321-60-8	2-Fluorobiphenyl	60%		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

4.2  
4

## Report of Analysis

<b>Client Sample ID:</b> BPOW 6-5		<b>Date Sampled:</b> 12/02/15
<b>Lab Sample ID:</b> JC9792-3		<b>Date Received:</b> 12/03/15
<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>	I99456.D	1	12/16/15	AMA	12/05/15	M:OP45645	M:MSI3717
Run #2							

Run #	Initial Volume	Final Volume
Run #1	940 ml	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
123-91-1	1,4-Dioxane	ND	0.21	0.081	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	72%		26-121%		
321-60-8	2-Fluorobiphenyl	65%		28-107%		
1718-51-0	Terphenyl-d14	83%		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

## Report of Analysis

<b>Client Sample ID:</b> BPOW 6-6		<b>Date Sampled:</b> 12/02/15
<b>Lab Sample ID:</b> JC9792-4		<b>Date Received:</b> 12/03/15
<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>	I99457.D	1	12/16/15	AMA	12/05/15	M:OP45645	M:MSI3717
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	ND	0.22	0.083	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.

---

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  
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

Impulse

<b>FED-EX Tracking #</b> #5 <b>Accutest Quote #</b>		<b>Bottle Order Control #</b> Jc9923	
<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>631-249-7600</b> Fax #: <b>631-249-7610</b> Sample Name(s): <b>PA Perovki 516 297-6247</b>		<b>Project Information</b> Project Name: <b>AGMNYM62235 // OU2 Monitoring Wells Northrop Grumman OU2 Hydro</b> Street: _____ Billing Information (if different from Report to): Company Name: <b>Bethpage NY</b> Street Address: <b>630 Plaza Drive, Suite 600</b> City: <b>Highlands Ranch, CO</b> Zip: <b>80129</b> Client Purchase Order #: <b>NY001496.154, 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> B8270SIM14DIOX VOCs 524,2 Full list		<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		<b>Collection</b> MECH/DI/Vial # Date Time Sampled by Matrix # of bottles Number of preserved bottles HDI HNO3 HNO2 H2SO4 H2SO4 NONE ED Water MESH ENCORE	
Field ID / Point of Collection 1 BPOW 5-3 2 FB120315PP2 3 TB120315PP2		LAB USE ONLY 1 V1047	
<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"/> 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 CUMMIL+ Commercial "A" = Results Only Commercial "B" = Results + QC Summary NJ Reduced = Results + QC Summary + Partial Raw data	
<b>Comments / Special Instructions</b> INITIAL ASSESSMENT: <b>3808</b> LABEL VERIFICATION: <b>JR</b> RL reporting for metals VOCs 524,2 Full list ✓ 524,2 + 40, VMS + F113			
<b>Sample Custody must be documented below each time samples change possession, including courier delivery.</b>			
Relinquished by Sampler: 1 <b>Carlo San Giovanni</b> Relinquished by Sampler: 3 Relinquished by: 5	Date Time: 12/3/15 1945 Date Time: Date Time: Date Time:	Received By: 1 <b>Carlo San Giovanni</b> Received By: 3 Received By: 5	Relinquished By: 2 <b>James Kwag</b> Relinquished By: 4 Date Time: 12/4/15 1946 Received By: 4
Custody Seal # Ken Arcadis		Intact <input checked="" type="checkbox"/> Not Intact <input type="checkbox"/> Preserved where applicable <input type="checkbox"/> On Ice <input checked="" type="checkbox"/> Cooler Temp. 2.5°C IP	

5.1  
5



## Report of Analysis

<b>Client Sample ID:</b>	BPOW5-3	<b>Date Sampled:</b>	12/03/15
<b>Lab Sample ID:</b>	JC9923-1	<b>Date Received:</b>	12/04/15
<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	1B100962.D	1	12/08/15	MD	n/a	n/a	V1B4779
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> BPOW5-3		<b>Date Sampled:</b> 12/03/15
<b>Lab Sample ID:</b> JC9923-1		<b>Date Received:</b> 12/04/15
<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	98%		78-114%
460-00-4	4-Bromofluorobenzene	97%		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> FB120315PP1		<b>Date Sampled:</b> 12/03/15
<b>Lab Sample ID:</b> JC9923-2		<b>Date Received:</b> 12/04/15
<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	98%		78-114%
460-00-4	4-Bromofluorobenzene	98%		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> TB120315PP1		<b>Date Sampled:</b> 12/03/15
<b>Lab Sample ID:</b> JC9923-3		<b>Date Received:</b> 12/04/15
<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	100%		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> BPOW5-3		<b>Date Sampled:</b> 12/03/15
<b>Lab Sample ID:</b> JC9923-1		<b>Date Received:</b> 12/04/15
<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>	I99391.D	1	12/14/15	AMA	12/07/15	M:OP45658	M:MSI3715
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	0.39	0.21	0.080	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	77%		26-121%		
321-60-8	2-Fluorobiphenyl	68%		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> FB120315PP1	<b>Date Sampled:</b> 12/03/15
<b>Lab Sample ID:</b> JC9923-2	<b>Date Received:</b> 12/04/15
<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>	I99392.D	1	12/14/15	AMA	12/07/15	M:OP45658	M:MSI3715
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	71%		26-121%		
321-60-8	2-Fluorobiphenyl	64%		28-107%		
1718-51-0	Terphenyl-d14	83%		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

4.2  
4

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

### **Data Review**

BETHPAGE, NEW YORK

Volatile and Semivolatile Analyses

SDGs #JC8939, JC9090 and JC9091

Analyses Performed By:  
Accutest Laboratories  
Dayton, New Jersey

Report #24831R  
December 22, 2015  
Review Level: Tier II  
Project #NY001496.1514.NAVI4

## SUMMARY

This data quality assessment summarizes the review of Sample Delivery Groups (SDGs) # JC8939, JC9090 and JC9091 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
JC8939	RE 117D1	JC8939-1	Water	11/18/2015		X	X			
	RE 117D2	JC8939-2	Water	11/18/2015		X	X			
	FB111815PP1	JC8939-3	Water	11/18/2015		X	X			
	TB111815PP1	JC8939-4	Water	11/18/2015		X				
JC9090	BPOW5-7	JC9090-1	Water	11/20/2015		X	X			
	FB112015PP1	JC9090-2	Water	11/20/2015		X	X			
	TB112015PP2	JC9090-3	Water	11/20/2015		X				
JC9091	RE119D1	JC9091-1	Water	11/20/2015		X	X			
	TB112015PP1	JC9091-2	Water	11/20/2015		X				

Notes:

1. Semivolatile (1,4-Dioxane) analysis was performed at Accutest Laboratories, Marlborough, MA (Accutest-NE) facility.
2. SDG JC8939: Matrix spike/matrix spike duplicate analysis was performed on sample location RE 117D2 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 8260C, 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..
SW-846 8260C			

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 not detected above the MDL in the associated blanks; therefore detected sample results were not associated with blank contamination in SDGs JC8939 and JC9091.

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
SDG JC9090:			
BPOW5-7	Acetone	Detected sample results >RL and <BAL	"UB" at the RL
	TIC: Unknown (RT7.60)	Detected sample results less than 5 times blank result	R

RL Reporting limit

## 3. Surrogates/System Monitoring Compounds

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

All surrogate recoveries were within control limits.

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

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

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

A MS/MSD analysis was not performed on a sample location associated with SDGs JC9090 or JC9091.

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

Sample Locations	Compound
<u>SDG JC8939:</u>	
RE 117D2	All compounds , except Freon 113

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 in SDGs JC8939, JC9090 and JC9091.

#### 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 SDGs JC8939, JC9090 or JC9091.

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

Tentatively identified compounds (TICs) were identified in SDG JC9090 in sample locations: BPOW5-7 and FB112015PP1. The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification. TICs are qualified as estimated (JN).

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

## DATA VALIDATION CHECKLIST FOR VOCs

VOCs: 524.2 and 8260C	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 in SDG JC8939.

A MS/MSD analysis was not performed on a sample location associated with SDGs JC9090 and JC9091.

## **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 for SDGs JC8939, JC9090 and JC9091.

## **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: December 22, 2015

PEER REVIEW BY: Todd Church

DATE: December 29, 2015

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

GW  
WFB  
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>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>631-249-7600</b> Fax #: <b>631-249-7610</b> Sampler(s) Name(s): <b>Pat Prerato 516-642-5166</b>		<b>Project Information</b> Project Name: <b>AGMNYM62235 // OU2 Monitoring 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: <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>		FED-EX Tracking # <b>#5</b> Accutest Quote # <b>JC8939</b> Bottle Order Control # Accutest Job # <b>JC8939</b>	
Requested Analysis (see TEST CODE sheet)		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 - Wipes 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"/> 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		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 CUMMUL+		Comments / Special Instructions RL reporting for metals Please use RE 11702 as a QA/QC MS/MSO sample	
Sample Custody must be documented below each time samples change possession, including courier delivery.					
Requisitioned by Sampler: Date Time: <b>11/18/15 1900</b>		Received By: <b>[Signature]</b> Date Time: <b>11/18/15 17:50</b>		Requisitioned By: Date Time: <b>11/18/15 16:11</b>	
Requisitioned by Sampler: Date Time:		Received By: Date Time:		Requisitioned By: Date Time:	
Requisitioned by Sampler: Date Time:		Received By: Date Time:		Requisitioned By: Date Time:	
Requisitioned by Sampler: Date Time:		Received By: Date Time:		Requisitioned By: 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>3.20</b>	

5.1  
5

LAB USE ONLY  
V900

INITIAL ASSESSMENT **3A Done**  
LABEL VERIFICATION **76**



FED-EX Tracking # <b>#5</b>		Bottle Order Control #	
Accutest Quote #		Accutest Job # <b>TC9939</b>	
<b>Client / Reporting Information</b>		<b>Project Information</b>	
Company Name <b>Arcadis</b>		Project Name: <b>AGMNYM62235 // OU2 Monitoring Wells Northrop Grumman OU2 Hydro</b>	
Street Address <b>2 Huntington Quad, Suite 1S10</b>		Billing Information (if different from Report to) Company Name <b>Arcadis, U.S., Inc. Attn: Accts Payable</b>	
City State Zip <b>Melville NY 11747</b>		Street Address <b>630 Plaza Drive, Suite 600</b>	
Project Contact <b>Soma Das, soma.das@arcadis-us.com</b>		City State Zip <b>Highlands Ranch, CO 80129</b>	
Phone # <b>631-249-7600</b>		Work Authorization #: NY001496_2015	
Sample(s) Name(s) <b>RE 117D1, RE 117D2, FB 111815PP2</b>		Project Manager <b>Carlo San Giovanni</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 T/A date 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 PMP) / Date:		RL-reporting for metals <b>Please use RE 117D2 as a QA/QC MS/MSD sample</b>	
Sample Custody must be documented below each time samples change possession, including courier delivery.			
Relinquished by: <b>[Signature]</b>	Date/Time: <b>11/18/15 1830</b>	Received By: <b>FDEX</b>	Date/Time: <b>09:30</b>
Relinquished by: <b>[Signature]</b>	Date/Time:	Received By:	Date/Time: <b>11/19/15</b>
Relinquished by:	Date/Time:	Received By:	Date/Time:
Relinquished by:	Date/Time:	Received By:	Date/Time:
Custody Seal #		On Ice / Cooler Temp. <b>15, 20, 25</b>	

5.1  
5

**JC8939: Chain of Custody**

**Page 3 of 4**

## Report of Analysis

<b>Client Sample ID:</b> RE 117D1		<b>Date Sampled:</b> 11/18/15
<b>Lab Sample ID:</b> JC8939-1		<b>Date Received:</b> 11/19/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8260C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	2D150433.D	1	11/24/15	BK	n/a	n/a	V2D6321
Run #2							

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

## VOA OU2 GW List

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	10	3.3	ug/l	
71-43-2	Benzene	ND	0.50	0.24	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.23	ug/l	
75-25-2	Bromoform	ND	1.0	0.23	ug/l	
74-83-9	Bromomethane	ND	2.0	0.42	ug/l	
78-93-3	2-Butanone (MEK)	ND	10	5.6	ug/l	
75-15-0	Carbon disulfide	ND	2.0	0.25	ug/l	
56-23-5	Carbon tetrachloride	ND	1.0	0.22	ug/l	
108-90-7	Chlorobenzene	ND	1.0	0.19	ug/l	
75-00-3	Chloroethane	ND	1.0	0.34	ug/l	
67-66-3	Chloroform	ND	1.0	0.19	ug/l	
74-87-3	Chloromethane	ND	1.0	0.41	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.15	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.17	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.18	ug/l	
75-35-4	1,1-Dichloroethene	ND	1.0	0.51	ug/l	
156-59-2	cis-1,2-Dichloroethene	ND	1.0	0.27	ug/l	
156-60-5	trans-1,2-Dichloroethene	ND	1.0	0.65	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.39	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.21	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.19	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.27	ug/l	
76-13-1	Freon 113	ND	5.0	0.52	ug/l	
591-78-6	2-Hexanone	ND	5.0	1.7	ug/l	
108-10-1	4-Methyl-2-pentanone(MIBK)	ND	5.0	1.0	ug/l	
75-09-2	Methylene chloride	ND	2.0	0.73	ug/l	
100-42-5	Styrene	ND	1.0	0.27	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	1.0	0.21	ug/l	
127-18-4	Tetrachloroethene	ND	1.0	0.40	ug/l	
108-88-3	Toluene	0.75	1.0	0.16	ug/l	J
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.25	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.21	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> RE 117D1		<b>Date Sampled:</b> 11/18/15
<b>Lab Sample ID:</b> JC8939-1		<b>Date Received:</b> 11/19/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8260C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**VOA OU2 GW List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethene	9.4	1.0	0.22	ug/l	
75-01-4	Vinyl chloride	ND	1.0	0.15	ug/l	
	m,p-Xylene	ND	1.0	0.38	ug/l	
95-47-6	o-Xylene	ND	1.0	0.17	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	102%		76-120%
17060-07-0	1,2-Dichloroethane-D4	100%		73-122%
2037-26-5	Toluene-D8	100%		84-119%
460-00-4	4-Bromofluorobenzene	102%		78-117%

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> RE 117D2		<b>Date Sampled:</b> 11/18/15
<b>Lab Sample ID:</b> JC8939-2		<b>Date Received:</b> 11/19/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8260C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	2D150333.D	1	11/21/15	BK	n/a	n/a	V2D6317
Run #2							

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

## VOA OU2 GW List

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	10	3.3	ug/l	J
71-43-2	Benzene	ND	0.50	0.24	ug/l	J
75-27-4	Bromodichloromethane	ND	1.0	0.23	ug/l	J
75-25-2	Bromoform	ND	1.0	0.23	ug/l	J
74-83-9	Bromomethane	ND	2.0	0.42	ug/l	J
78-93-3	2-Butanone (MEK)	ND	10	5.6	ug/l	J
75-15-0	Carbon disulfide	ND	2.0	0.25	ug/l	J
56-23-5	Carbon tetrachloride	ND	1.0	0.22	ug/l	J
108-90-7	Chlorobenzene	ND	1.0	0.19	ug/l	J
75-00-3	Chloroethane	ND	1.0	0.34	ug/l	J
67-66-3	Chloroform	ND	1.0	0.19	ug/l	J
74-87-3	Chloromethane	ND	1.0	0.41	ug/l	J
124-48-1	Dibromochloromethane	ND	1.0	0.15	ug/l	J
75-34-3	1,1-Dichloroethane	ND	1.0	0.17	ug/l	J
107-06-2	1,2-Dichloroethane	ND	1.0	0.18	ug/l	J
75-35-4	1,1-Dichloroethene	ND	1.0	0.51	ug/l	J
156-59-2	cis-1,2-Dichloroethene	ND	1.0	0.27	ug/l	J
156-60-5	trans-1,2-Dichloroethene	ND	1.0	0.65	ug/l	J
78-87-5	1,2-Dichloropropane	ND	1.0	0.39	ug/l	J
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.21	ug/l	J
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.19	ug/l	J
100-41-4	Ethylbenzene	ND	1.0	0.27	ug/l	J
76-13-1	Freon 113	ND	5.0	0.52	ug/l	J
591-78-6	2-Hexanone	ND	5.0	1.7	ug/l	J
108-10-1	4-Methyl-2-pentanone(MIBK)	ND	5.0	1.0	ug/l	J
75-09-2	Methylene chloride	ND	2.0	0.73	ug/l	J
100-42-5	Styrene	ND	1.0	0.27	ug/l	J
79-34-5	1,1,2,2-Tetrachloroethane	ND	1.0	0.21	ug/l	J
127-18-4	Tetrachloroethene	ND	1.0	0.40	ug/l	J
108-88-3	Toluene	0.98	1.0	0.16	ug/l	J
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.25	ug/l	J
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.21	ug/l	J

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> RE 117D2		<b>Date Sampled:</b> 11/18/15
<b>Lab Sample ID:</b> JC8939-2		<b>Date Received:</b> 11/19/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8260C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**VOA OU2 GW List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethene	ND	1.0	0.22	ug/l	J
75-01-4	Vinyl chloride	ND	1.0	0.15	ug/l	J
	m,p-Xylene	ND	1.0	0.38	ug/l	J
95-47-6	o-Xylene	ND	1.0	0.17	ug/l	J

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	100%		76-120%
17060-07-0	1,2-Dichloroethane-D4	98%		73-122%
2037-26-5	Toluene-D8	100%		84-119%
460-00-4	4-Bromofluorobenzene	100%		78-117%

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> FB111815PP1		
<b>Lab Sample ID:</b> JC8939-3		<b>Date Sampled:</b> 11/18/15
<b>Matrix:</b> AQ - Field Blank Water		<b>Date Received:</b> 11/19/15
<b>Method:</b> SW846 8260C		<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	2D150331.D	1	11/21/15	BK	n/a	n/a	V2D6317
Run #2							

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

## VOA OU2 GW List

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	10	3.3	ug/l	
71-43-2	Benzene	ND	0.50	0.24	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.23	ug/l	
75-25-2	Bromoform	ND	1.0	0.23	ug/l	
74-83-9	Bromomethane	ND	2.0	0.42	ug/l	
78-93-3	2-Butanone (MEK)	ND	10	5.6	ug/l	
75-15-0	Carbon disulfide	ND	2.0	0.25	ug/l	
56-23-5	Carbon tetrachloride	ND	1.0	0.22	ug/l	
108-90-7	Chlorobenzene	ND	1.0	0.19	ug/l	
75-00-3	Chloroethane	ND	1.0	0.34	ug/l	
67-66-3	Chloroform	ND	1.0	0.19	ug/l	
74-87-3	Chloromethane	ND	1.0	0.41	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.15	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.17	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.18	ug/l	
75-35-4	1,1-Dichloroethene	ND	1.0	0.51	ug/l	
156-59-2	cis-1,2-Dichloroethene	ND	1.0	0.27	ug/l	
156-60-5	trans-1,2-Dichloroethene	ND	1.0	0.65	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.39	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.21	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.19	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.27	ug/l	
76-13-1	Freon 113	ND	5.0	0.52	ug/l	
591-78-6	2-Hexanone	ND	5.0	1.7	ug/l	
108-10-1	4-Methyl-2-pentanone(MIBK)	ND	5.0	1.0	ug/l	
75-09-2	Methylene chloride	ND	2.0	0.73	ug/l	
100-42-5	Styrene	ND	1.0	0.27	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	1.0	0.21	ug/l	
127-18-4	Tetrachloroethene	ND	1.0	0.40	ug/l	
108-88-3	Toluene	ND	1.0	0.16	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.25	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.21	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> FB111815PP1		<b>Date Sampled:</b> 11/18/15
<b>Lab Sample ID:</b> JC8939-3		<b>Date Received:</b> 11/19/15
<b>Matrix:</b> AQ - Field Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8260C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**VOA OU2 GW List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethene	ND	1.0	0.22	ug/l	
75-01-4	Vinyl chloride	ND	1.0	0.15	ug/l	
	m,p-Xylene	ND	1.0	0.38	ug/l	
95-47-6	o-Xylene	ND	1.0	0.17	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	99%		76-120%
17060-07-0	1,2-Dichloroethane-D4	97%		73-122%
2037-26-5	Toluene-D8	99%		84-119%
460-00-4	4-Bromofluorobenzene	101%		78-117%

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> TB111815PP1		<b>Date Sampled:</b> 11/18/15
<b>Lab Sample ID:</b> JC8939-4		<b>Date Received:</b> 11/19/15
<b>Matrix:</b> AQ - Trip Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8260C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**VOA OU2 GW List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethene	ND	1.0	0.22	ug/l	
75-01-4	Vinyl chloride	ND	1.0	0.15	ug/l	
	m,p-Xylene	ND	1.0	0.38	ug/l	
95-47-6	o-Xylene	ND	1.0	0.17	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	100%		76-120%
17060-07-0	1,2-Dichloroethane-D4	97%		73-122%
2037-26-5	Toluene-D8	99%		84-119%
460-00-4	4-Bromofluorobenzene	101%		78-117%

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  
4

## Report of Analysis

<b>Client Sample ID:</b> RE 117D1	<b>Date Sampled:</b> 11/18/15
<b>Lab Sample ID:</b> JC8939-1	<b>Date Received:</b> 11/19/15
<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>	I99203.D	1	11/27/15	AMA	11/22/15	M:OP45518	M:MSI3704
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	77%		26-121%		
321-60-8	2-Fluorobiphenyl	64%		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

4.1  
4

## Report of Analysis

<b>Client Sample ID:</b> RE 117D2		<b>Date Sampled:</b> 11/18/15
<b>Lab Sample ID:</b> JC8939-2		<b>Date Received:</b> 11/19/15
<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>	I99186.D	1	11/26/15	AMA	11/22/15	M:OP45518	M:MSI3703
Run #2							

Run #	Initial Volume	Final Volume
Run #1	1000 ml	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
123-91-1	1,4-Dioxane	ND	0.20	0.076	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	75%		26-121%		
321-60-8	2-Fluorobiphenyl	66%		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> FB111815PP1	<b>Date Sampled:</b> 11/18/15
<b>Lab Sample ID:</b> JC8939-3	<b>Date Received:</b> 11/19/15
<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>	I99204.D	1	11/27/15	AMA	11/22/15	M:OP45518	M:MSI3704
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	ND	0.21	0.079	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	79%		26-121%		
321-60-8	2-Fluorobiphenyl	68%		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

GW  
FB  
WTB

NJ

**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 # 44 Bottle Order Control #  
Accutest Quote # 44 Accutest Job # JC9090

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>												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															
City State Zip <b>Melville NY 11747</b>		City State <b>Bethpage NY</b>												LAB USE ONLY  INITIAL ASSESSMENT <u>3B CB</u> LABEL VERIFICATION <u>05</u>			
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 # Fax # <b>631-249-7600 631-249-7610</b>		Project # <b>1614</b>															
Sample(s) Name(s) <b>Pat Krzotki 516 287-0247</b>		Client Purchase Order # <b>NY001496.514:NAV13</b>															
		Work Authorization #: NY001496_2015															
		Project Manager <b>Carlo San Giovanni</b>															
		Attention: <b>Soma Das</b>															
		City State Zip <b>Highlands Ranch, CO 80129</b>															
		Collection															
Account Sample #	Field ID / Point of Collection	MEOH/DI Vial #	Date	Time	Sampled by	Metric	# of bottles	HCl	NH <sub>4</sub> Cl	NH <sub>3</sub> Cl	NH <sub>2</sub> SO <sub>4</sub>	NH <sub>2</sub> NO <sub>2</sub>	DI Water	MEDIA	ENCLOSURE		
1	BP0W 5-7		11/20/15	1528	KP	GW	3										
2	FB112015PP2		11/20/15	1528	KP	FB	2										
3	TB112015PP2		11/20/15	0940	-	TB	2										

VOCs 524.2 Full List  
B82705M14 P10X

5.1  
5

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"/> 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 CUMML+ Commercial "A" = Results Only Commercial "B" = Results + QC Summary NJ Reduced = Results + QC Summary + Partial Raw data	
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:
1 <i>Carlo San Giovanni</i>	11/20/15 16:00	1 Robert chambers	11/20/15	2 Robert chambers	11/20/15
Relinquished by Sampler:	Date Time:	Received By:	Date Time:	Relinquished By:	Date Time:
3		3		4	
Relinquished by:	Date Time:	Received By:	Date Time:	Relinquished By:	Date Time:
		5			
Custody Seal # <u>361</u>		<input type="checkbox"/> Intact <input type="checkbox"/> Not Intact		Preserved where applicable <input type="checkbox"/> On Ice <input checked="" type="checkbox"/> Cooler Temp. <u>2.0°C</u>	

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>Satellite (s) Name(s)</b> Vest Product 51687 6477		<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> Arcadis, U.S., Inc. Attn: Accts Payable <b>Project #</b> NY001496.1614.NAVI3 <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>Attention:</b> Soma Das		<b>FED-EX Tracking #</b> #4 <b>Accutest Quote #</b> <b>Accutest Job #</b> JC9090		<b>Matrix Codes</b> DW - Drinking Water GW - Ground Water WW - Water SW - Surface Water SO - Soil SL - Sludge SEP - 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 (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+		<b>Comments / Special Instructions</b> VMS 524.2 Full list V524SL4 + 40 VMS + F113	
Sample Custody must be documented below each time sample change possession, including courier delivery.							
<b>Relinquished by Sampler</b> 1 _____ <b>Date Time:</b> 11/20/15 1730	<b>Received By:</b> 1 _____ <b>Date Time:</b> _____	<b>Relinquished by Sampler</b> 2 _____ <b>Date Time:</b> _____	<b>Received By:</b> 2 _____ <b>Date Time:</b> _____	<b>Relinquished by Sampler</b> 3 _____ <b>Date Time:</b> _____	<b>Received By:</b> 3 _____ <b>Date Time:</b> _____	<b>Relinquished by Sampler</b> 4 _____ <b>Date Time:</b> _____	<b>Received By:</b> 4 _____ <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 type="checkbox"/> Inset <input type="checkbox"/> Not Inset <b>Preserved where applicable:</b> <input type="checkbox"/> On Ice <input type="checkbox"/> Cooler Temp.		_____ _____	

5.1  
5

17B



## Report of Analysis

<b>Client Sample ID:</b> BPOW5-7		<b>Date Sampled:</b> 11/20/15
<b>Lab Sample ID:</b> JC9090-1		<b>Date Received:</b> 11/20/15
<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	102%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	unknown	7.60	.65	ug/l	J R
	Total TIC, Volatile		.65	ug/l	J

ND = Not detected      MDL = Method Detection Limit  
 RL = Reporting Limit

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> FB112015PP1		
<b>Lab Sample ID:</b> JC9090-2		<b>Date Sampled:</b> 11/20/15
<b>Matrix:</b> AQ - Field Blank Water		<b>Date Received:</b> 11/20/15
<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	1B100759.D	1	11/24/15	MD	n/a	n/a	V1B4770
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	0.20	0.50	0.047	ug/l	J
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> FB112015PP1		<b>Date Sampled:</b> 11/20/15
<b>Lab Sample ID:</b> JC9090-2		<b>Date Received:</b> 11/20/15
<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	102%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	unknown	7.60	.59	ug/l	J N
	Total TIC, Volatile		.59	ug/l	J N

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

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



## Report of Analysis

<b>Client Sample ID:</b> TB112015PP2		<b>Date Sampled:</b> 11/20/15
<b>Lab Sample ID:</b> JC9090-3		<b>Date Received:</b> 11/20/15
<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	103%		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

4.3  
4

## Report of Analysis

<b>Client Sample ID:</b> BPOW5-7		<b>Date Sampled:</b> 11/20/15
<b>Lab Sample ID:</b> JC9090-1		<b>Date Received:</b> 11/20/15
<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>	I99205.D	1	11/27/15	AMA	11/22/15	M:OP45518	M:MSI3704
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	69%		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> FB112015PP1	<b>Date Sampled:</b> 11/20/15
<b>Lab Sample ID:</b> JC9090-2	<b>Date Received:</b> 11/20/15
<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>	I99206.D	1	11/27/15	AMA	11/22/15	M:OP45518	M:MSI3704
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	ND	0.22	0.083	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	82%		26-121%		
321-60-8	2-Fluorobiphenyl	74%		28-107%		
1718-51-0	Terphenyl-d14	87%		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> RE119D1		<b>Date Sampled:</b> 11/20/15
<b>Lab Sample ID:</b> JC9091-1		<b>Date Received:</b> 11/20/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8260C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	2B137063.D	1	12/01/15	EH	n/a	n/a	V2B6106
Run #2							

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

## VOA OU2 GW List

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	10	3.3	ug/l	
71-43-2	Benzene	ND	0.50	0.24	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.23	ug/l	
75-25-2	Bromoform	ND	1.0	0.23	ug/l	
74-83-9	Bromomethane	ND	2.0	0.42	ug/l	
78-93-3	2-Butanone (MEK)	ND	10	5.6	ug/l	
75-15-0	Carbon disulfide	ND	2.0	0.25	ug/l	
56-23-5	Carbon tetrachloride	ND	1.0	0.22	ug/l	
108-90-7	Chlorobenzene	ND	1.0	0.19	ug/l	
75-00-3	Chloroethane	ND	1.0	0.34	ug/l	
67-66-3	Chloroform	ND	1.0	0.19	ug/l	
74-87-3	Chloromethane	ND	1.0	0.41	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.15	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.17	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.18	ug/l	
75-35-4	1,1-Dichloroethene	ND	1.0	0.51	ug/l	
156-59-2	cis-1,2-Dichloroethene	ND	1.0	0.27	ug/l	
156-60-5	trans-1,2-Dichloroethene	ND	1.0	0.65	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.39	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.21	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.19	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.27	ug/l	
76-13-1	Freon 113	ND	5.0	0.52	ug/l	
591-78-6	2-Hexanone	ND	5.0	1.7	ug/l	
108-10-1	4-Methyl-2-pentanone(MIBK)	ND	5.0	1.0	ug/l	
75-09-2	Methylene chloride	ND	2.0	0.73	ug/l	
100-42-5	Styrene	ND	1.0	0.27	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	1.0	0.21	ug/l	
127-18-4	Tetrachloroethene	ND	1.0	0.40	ug/l	
108-88-3	Toluene	0.72	1.0	0.16	ug/l	J
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.25	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.21	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> RE119D1		<b>Date Sampled:</b> 11/20/15
<b>Lab Sample ID:</b> JC9091-1		<b>Date Received:</b> 11/20/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8260C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**VOA OU2 GW List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethene	ND	1.0	0.22	ug/l	
75-01-4	Vinyl chloride	ND	1.0	0.15	ug/l	
	m,p-Xylene	ND	1.0	0.38	ug/l	
95-47-6	o-Xylene	ND	1.0	0.17	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	105%		76-120%
17060-07-0	1,2-Dichloroethane-D4	115%		73-122%
2037-26-5	Toluene-D8	100%		84-119%
460-00-4	4-Bromofluorobenzene	100%		78-117%

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  
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## Report of Analysis

<b>Client Sample ID:</b> TB112015PP1		<b>Date Sampled:</b> 11/20/15
<b>Lab Sample ID:</b> JC9091-2		<b>Date Received:</b> 11/20/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8260C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	2B137000.D	1	11/28/15	EH	n/a	n/a	V2B6102
Run #2							

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

## VOA OU2 GW List

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	10	3.3	ug/l	
71-43-2	Benzene	ND	0.50	0.24	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.23	ug/l	
75-25-2	Bromoform	ND	1.0	0.23	ug/l	
74-83-9	Bromomethane	ND	2.0	0.42	ug/l	
78-93-3	2-Butanone (MEK)	ND	10	5.6	ug/l	
75-15-0	Carbon disulfide	ND	2.0	0.25	ug/l	
56-23-5	Carbon tetrachloride	ND	1.0	0.22	ug/l	
108-90-7	Chlorobenzene	ND	1.0	0.19	ug/l	
75-00-3	Chloroethane	ND	1.0	0.34	ug/l	
67-66-3	Chloroform	ND	1.0	0.19	ug/l	
74-87-3	Chloromethane	ND	1.0	0.41	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.15	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.17	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.18	ug/l	
75-35-4	1,1-Dichloroethene	ND	1.0	0.51	ug/l	
156-59-2	cis-1,2-Dichloroethene	ND	1.0	0.27	ug/l	
156-60-5	trans-1,2-Dichloroethene	ND	1.0	0.65	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.39	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.21	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.19	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.27	ug/l	
76-13-1	Freon 113	ND	5.0	0.52	ug/l	
591-78-6	2-Hexanone	ND	5.0	1.7	ug/l	
108-10-1	4-Methyl-2-pentanone(MIBK)	ND	5.0	1.0	ug/l	
75-09-2	Methylene chloride	ND	2.0	0.73	ug/l	
100-42-5	Styrene	ND	1.0	0.27	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	1.0	0.21	ug/l	
127-18-4	Tetrachloroethene	ND	1.0	0.40	ug/l	
108-88-3	Toluene	ND	1.0	0.16	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.25	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.21	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> TB112015PP1		<b>Date Sampled:</b> 11/20/15
<b>Lab Sample ID:</b> JC9091-2		<b>Date Received:</b> 11/20/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8260C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**VOA OU2 GW List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethene	ND	1.0	0.22	ug/l	
75-01-4	Vinyl chloride	ND	1.0	0.15	ug/l	
	m,p-Xylene	ND	1.0	0.38	ug/l	
95-47-6	o-Xylene	ND	1.0	0.17	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	102%		76-120%
17060-07-0	1,2-Dichloroethane-D4	103%		73-122%
2037-26-5	Toluene-D8	100%		84-119%
460-00-4	4-Bromofluorobenzene	102%		78-117%

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> RE119D1		<b>Date Sampled:</b> 11/20/15
<b>Lab Sample ID:</b> JC9091-1		<b>Date Received:</b> 11/20/15
<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>	I99207.D	1	11/27/15	AMA	11/22/15	M:OP45518	M:MSI3704
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	74%		26-121%		
321-60-8	2-Fluorobiphenyl	67%		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

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

### **Data Review**

BETHPAGE, NEW YORK

Volatile and Semivolatile Analyses

SDGs #JC9220, JC9566 and JC9689

Analyses Performed By:  
Accutest Laboratories  
Dayton, New Jersey

Report #24832R  
December 28, 2015  
Review Level: Tier II  
Project #NY001496.1514.NAVI4

## SUMMARY

This data quality assessment summarizes the review of Sample Delivery Groups (SDGs) # JC9220, JC9566 and JC9689 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
JC9220	FB112315PP1	JC9220-1	Water	11/23/2015		X	X			
	TB112315PP1	JC9220-2	Water	11/23/2015		X				
	RE118D1	JC9220-3	Water	11/23/2015		X	X			
JC9566	FB113015PP1	JC9566-1	Water	11/30/2015		X	X			
	BPOW 6-1	JC9566-2	Water	11/30/2015		X	X			
	BPOW 6-2	JC9566-3	Water	11/30/2015		X	X			
	TB113015PP1	JC9566-4	Water	11/30/2015		X				
JC9689	FB120115PP1	JC9689-1	Water	12/01/2015		X	X			
	BPOW 6-3	JC9689-2	Water	12/01/2015		X	X			
	BPOW 6-4	JC9689-3	Water	12/01/2015		X	X			
	REP120115PP1	JC9689-4	Water	12/01/2015	BPOW 6-4	X	X			
	TB120115PP1	JC9689-5	Water	12/01/2015		X				

Note:

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 8260C, 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..
SW-846 8260C			

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 not detected above the MDL in the associated blanks; therefore detected sample results were not associated with blank contamination in SDG JC9220.

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 in SDG JC9566.

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
SDG JC9689:			
BPOW 6-3	Acetone	Detected sample results >RL and <BAL	"UB" at the RL

RL Reporting limit

## 3. Surrogates/System Monitoring Compounds

All samples to be analyzed for organic compounds are spiked with surrogate compounds prior to sample preparation to evaluate overall laboratory performance and efficiency of the analytical technique. VOC

analysis requires that all surrogates associated with the analysis exhibit recoveries within the laboratory-established acceptance limits.

All surrogate recoveries were within control limits.

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

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

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

A MS/MSD analysis was not performed on a sample location associated with SDGs JC9220, JC9566 or JC9689.

#### 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 SDGs JC9220 or JC9566.

Results for duplicate samples are summarized in the following table.

Sample ID/Duplicate ID	Compound	Sample Result	Duplicate Result	RPD
SDG JC9689:				
BPOW 6-4/ REP120115PP1	All compounds	U	U	AC

AC Acceptable

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

#### 7. Laboratory Duplicate Analysis

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

concentrations are less than or equal to 5 times the RL, a control limit of two times the RL is applied for water matrices.

A laboratory duplicate sample was not performed on a sample location associated with SDGs JC9220, JC9566 or JC9689.

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

Tentatively identified compounds (TICs) were not identified in SDG JC9220. TICs were identified in SDG JC9566 in sample locations FB113015PP1; and, in SDG JC9689 in sample location FB12015PP1. The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification. TICs are qualified as estimated (JN).

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

## DATA VALIDATION CHECKLIST FOR VOCs

VOCs: 524.2 and 8260C	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		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 SDGs JC9220, JC9566 or JC9689.

## 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 SDG JC9220 and JC9566

Results for duplicate samples are summarized in the following table.

Sample ID/Duplicate ID	Compound	Sample Result	Duplicate Result	RPD
SDG JC9689:				
BPOW 6-4/ REP120115PP1	1,4-Dioxane	0.22 U	0.21 U	AC

AC Acceptable

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

## 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		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: December 28, 2015

PEER REVIEW BY: Todd Church

DATE: December 29, 2015

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





# Report of Analysis

<b>Client Sample ID:</b> FB112315PP1	
<b>Lab Sample ID:</b> JC9220-1	<b>Date Sampled:</b> 11/23/15
<b>Matrix:</b> AQ - Field Blank Water	<b>Date Received:</b> 11/24/15
<b>Method:</b> SW846 8260C	<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	1A156464.D	1	11/30/15	VC	n/a	n/a	V1A6717
Run #2							

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

### VOA OU2 GW List

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	10	3.3	ug/l	
71-43-2	Benzene	ND	0.50	0.24	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.23	ug/l	
75-25-2	Bromoform	ND	1.0	0.23	ug/l	
74-83-9	Bromomethane	ND	2.0	0.42	ug/l	
78-93-3	2-Butanone (MEK)	ND	10	5.6	ug/l	
75-15-0	Carbon disulfide	ND	2.0	0.25	ug/l	
56-23-5	Carbon tetrachloride	ND	1.0	0.22	ug/l	
108-90-7	Chlorobenzene	ND	1.0	0.19	ug/l	
75-00-3	Chloroethane	ND	1.0	0.34	ug/l	
67-66-3	Chloroform	ND	1.0	0.19	ug/l	
74-87-3	Chloromethane	ND	1.0	0.41	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.15	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.17	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.18	ug/l	
75-35-4	1,1-Dichloroethene	ND	1.0	0.51	ug/l	
156-59-2	cis-1,2-Dichloroethene	ND	1.0	0.27	ug/l	
156-60-5	trans-1,2-Dichloroethene	ND	1.0	0.65	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.39	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.21	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.19	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.27	ug/l	
76-13-1	Freon 113	ND	5.0	0.52	ug/l	
591-78-6	2-Hexanone	ND	5.0	1.7	ug/l	
108-10-1	4-Methyl-2-pentanone(MIBK)	ND	5.0	1.0	ug/l	
75-09-2	Methylene chloride	ND	2.0	0.73	ug/l	
100-42-5	Styrene	ND	1.0	0.27	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	1.0	0.21	ug/l	
127-18-4	Tetrachloroethene	ND	1.0	0.40	ug/l	
108-88-3	Toluene	ND	1.0	0.16	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.25	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.21	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> FB112315PP1 <b>Lab Sample ID:</b> JC9220-1 <b>Matrix:</b> AQ - Field Blank Water <b>Method:</b> SW846 8260C <b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY	<b>Date Sampled:</b> 11/23/15 <b>Date Received:</b> 11/24/15 <b>Percent Solids:</b> n/a
--	---

**VOA OU2 GW List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethene	ND	1.0	0.22	ug/l	
75-01-4	Vinyl chloride	ND	1.0	0.15	ug/l	
	m,p-Xylene	ND	1.0	0.38	ug/l	
95-47-6	o-Xylene	ND	1.0	0.17	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	100%		76-120%
17060-07-0	1,2-Dichloroethane-D4	97%		73-122%
2037-26-5	Toluene-D8	99%		84-119%
460-00-4	4-Bromofluorobenzene	100%		78-117%

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> TB112315PP1		<b>Date Sampled:</b> 11/23/15
<b>Lab Sample ID:</b> JC9220-2		<b>Date Received:</b> 11/24/15
<b>Matrix:</b> AQ - Trip Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8260C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1A156465.D	1	11/30/15	VC	n/a	n/a	V1A6717
Run #2							

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

## VOA OU2 GW List

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	10	3.3	ug/l	
71-43-2	Benzene	ND	0.50	0.24	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.23	ug/l	
75-25-2	Bromoform	ND	1.0	0.23	ug/l	
74-83-9	Bromomethane	ND	2.0	0.42	ug/l	
78-93-3	2-Butanone (MEK)	ND	10	5.6	ug/l	
75-15-0	Carbon disulfide	ND	2.0	0.25	ug/l	
56-23-5	Carbon tetrachloride	ND	1.0	0.22	ug/l	
108-90-7	Chlorobenzene	ND	1.0	0.19	ug/l	
75-00-3	Chloroethane	ND	1.0	0.34	ug/l	
67-66-3	Chloroform	ND	1.0	0.19	ug/l	
74-87-3	Chloromethane	ND	1.0	0.41	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.15	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.17	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.18	ug/l	
75-35-4	1,1-Dichloroethene	ND	1.0	0.51	ug/l	
156-59-2	cis-1,2-Dichloroethene	ND	1.0	0.27	ug/l	
156-60-5	trans-1,2-Dichloroethene	ND	1.0	0.65	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.39	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.21	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.19	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.27	ug/l	
76-13-1	Freon 113	ND	5.0	0.52	ug/l	
591-78-6	2-Hexanone	ND	5.0	1.7	ug/l	
108-10-1	4-Methyl-2-pentanone(MIBK)	ND	5.0	1.0	ug/l	
75-09-2	Methylene chloride	ND	2.0	0.73	ug/l	
100-42-5	Styrene	ND	1.0	0.27	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	1.0	0.21	ug/l	
127-18-4	Tetrachloroethene	ND	1.0	0.40	ug/l	
108-88-3	Toluene	ND	1.0	0.16	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.25	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.21	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> TB112315PP1		<b>Date Sampled:</b> 11/23/15
<b>Lab Sample ID:</b> JC9220-2		<b>Date Received:</b> 11/24/15
<b>Matrix:</b> AQ - Trip Blank Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8260C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**VOA OU2 GW List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethene	ND	1.0	0.22	ug/l	
75-01-4	Vinyl chloride	ND	1.0	0.15	ug/l	
	m,p-Xylene	ND	1.0	0.38	ug/l	
95-47-6	o-Xylene	ND	1.0	0.17	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	100%		76-120%
17060-07-0	1,2-Dichloroethane-D4	97%		73-122%
2037-26-5	Toluene-D8	100%		84-119%
460-00-4	4-Bromofluorobenzene	99%		78-117%

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> RE 118D1		<b>Date Sampled:</b> 11/23/15
<b>Lab Sample ID:</b> JC9220-3		<b>Date Received:</b> 11/24/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8260C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1A156488.D	1	12/01/15	VC	n/a	n/a	V1A6718
Run #2							

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

### VOA OU2 GW List

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	10	3.3	ug/l	
71-43-2	Benzene	ND	0.50	0.24	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.23	ug/l	
75-25-2	Bromoform	ND	1.0	0.23	ug/l	
74-83-9	Bromomethane	ND	2.0	0.42	ug/l	
78-93-3	2-Butanone (MEK)	ND	10	5.6	ug/l	
75-15-0	Carbon disulfide	ND	2.0	0.25	ug/l	
56-23-5	Carbon tetrachloride	ND	1.0	0.22	ug/l	
108-90-7	Chlorobenzene	ND	1.0	0.19	ug/l	
75-00-3	Chloroethane	ND	1.0	0.34	ug/l	
67-66-3	Chloroform	ND	1.0	0.19	ug/l	
74-87-3	Chloromethane	ND	1.0	0.41	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.15	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.17	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.18	ug/l	
75-35-4	1,1-Dichloroethene	ND	1.0	0.51	ug/l	
156-59-2	cis-1,2-Dichloroethene	ND	1.0	0.27	ug/l	
156-60-5	trans-1,2-Dichloroethene	ND	1.0	0.65	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.39	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.21	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.19	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.27	ug/l	
76-13-1	Freon 113	ND	5.0	0.52	ug/l	
591-78-6	2-Hexanone	ND	5.0	1.7	ug/l	
108-10-1	4-Methyl-2-pentanone(MIBK)	ND	5.0	1.0	ug/l	
75-09-2	Methylene chloride	ND	2.0	0.73	ug/l	
100-42-5	Styrene	ND	1.0	0.27	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	1.0	0.21	ug/l	
127-18-4	Tetrachloroethene	ND	1.0	0.40	ug/l	
108-88-3	Toluene	0.57	1.0	0.16	ug/l	J
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.25	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.21	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> RE 118D1		<b>Date Sampled:</b> 11/23/15
<b>Lab Sample ID:</b> JC9220-3		<b>Date Received:</b> 11/24/15
<b>Matrix:</b> AQ - Ground Water		<b>Percent Solids:</b> n/a
<b>Method:</b> SW846 8260C		
<b>Project:</b> Northrop Grumman, OU2 Hydro, Bethpage, NY		

**VOA OU2 GW List**

CAS No.	Compound	Result	RL	MDL	Units	Q
79-01-6	Trichloroethene	ND	1.0	0.22	ug/l	
75-01-4	Vinyl chloride	ND	1.0	0.15	ug/l	
	m,p-Xylene	ND	1.0	0.38	ug/l	
95-47-6	o-Xylene	ND	1.0	0.17	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	98%		76-120%
17060-07-0	1,2-Dichloroethane-D4	97%		73-122%
2037-26-5	Toluene-D8	100%		84-119%
460-00-4	4-Bromofluorobenzene	101%		78-117%

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> FB112315PP1		<b>Date Sampled:</b> 11/23/15
<b>Lab Sample ID:</b> JC9220-1		<b>Date Received:</b> 11/24/15
<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>	I99247.D	1	12/01/15	AMA	11/24/15	M:OP45543	M:MSI3707
Run #2							

	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	ND	0.21	0.079	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	88%		26-121%		
321-60-8	2-Fluorobiphenyl	75%		28-107%		
1718-51-0	Terphenyl-d14	86%		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

4.1  
4

## Report of Analysis

<b>Client Sample ID:</b> RE 118D1		<b>Date Sampled:</b> 11/23/15
<b>Lab Sample ID:</b> JC9220-3		<b>Date Received:</b> 11/24/15
<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>	I99248.D	1	12/01/15	AMA	11/24/15	M:OP45543	M:MSI3707
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	ND	0.22	0.083	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	67%		26-121%		
321-60-8	2-Fluorobiphenyl	60%		28-107%		
1718-51-0	Terphenyl-d14	87%		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

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 # **#5** Bottle Order Control #  
Accutest Quote # **5C9566** Accutest Job #

Client / Reporting Information		Project Information		Requested Analysis (see TEST CODE sheet)												Matrix Codes
Company Name <b>Arcadis</b>		Project Name <b>AGMNYM62235 // OU2 Monitoring Wells</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>Northrop Grumman OU2 Hydro</b>														
City State Zip <b>Melville NY 11747</b>		Billing Information (if different from Report to) Company Name <b>Bethpage NY</b>														LAB USE ONLY  V988
Project Contact <b>Soma Das, soma.das@arcadis-us.com</b>		Street Address <b>Arcadis, U.S., Inc. Attn: Accts Payable</b>														
Phone # <b>631-249-7600</b>		Client Purchase Order # <b>NY001946.1514.NAVIS</b>														
Samples Name(s) <b>11/30/15</b>		Work Authorization # <b>NY001496_2015</b>														
Phone # <b>516 237-6247</b>		Project Manager <b>Carlo San Giovanni</b>														
City State Zip <b>Highlands Ranch, CO 80129</b>		Attention: <b>Soma Das</b>														

Matrix	# of bottles	HC	HNOC	HRBC	HNOC	HRBC	ED Water	MICH	ENDORE
FB	3	3							
GH	3	3							
TB	2	2							

VC62602NG6666WY740  
 B8270S1M1 4DIOX  
 VDCS 524.2 Full list

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  
 Emergency & Rush T/A data available VIA Lablink

Approved By (Accutest PM): / Date: INITIAL ASSESSMENT - NL3A  
LABEL VERIFICATION - COB

Data Deliverable Information  
 Commercial "A" (Level 1)  
 Commercial "B" (Level 2)  
 FULLT (Level 3+4)  
 NJ Reduced  
 Commercial "C"  
 NYASP Category A  
 NYASP Category B  
 State Forms  
 EDD Format  
 Other LUMMLC+

Commercial "A" = Results Only  
 Commercial "B" = Results + QC Summary  
 NJ Reduced = Results + QC Summary + Partial Raw data

Comments / Special Instructions  
VDCS 524.2 Full list  
V5245L4+40 VMS + F113

Sample Custody must be documented below each time samples change possession, including courier delivery.

Relinquished by: <u>[Signature]</u>	Date Time: <u>11/30/15 2:30</u>	Received By: <u>Robert Chambers</u>	Date Time: <u>12-1-15 10:15</u>	Relinquished By: <u>Robert Chambers</u>	Date Time: <u>12-1-15</u>	Received By: <u>[Signature]</u>
Relinquished by:	Date Time:	Received By:	Date Time:	Relinquished By:	Date Time:	Received By:
Relinquished by:	Date Time:	Received By:	Date Time:	Relinquished By:	Date Time:	Received By:

Custody Seal # 364  
 Intact  
 Not Intact  
 Preserved where applicable  
 On Ice  Cooler Temp: 2.0°C

5.1  
5

Client / Reporting Information		Project Information				Requested Analysis (see TEST CODE sheet)										Matrix Codes							
Company Name <b>Arcadis</b>		Project Name: <b>AGMNYM62235 // OU2 Monitoring Wells Northrop Grumman OU2 Hydro</b>				Requested Analysis (see TEST CODE sheet)  Matrix Codes: DW - Drinking Water GW - Ground Water WW - Water SW - Surface Water SO - Soil SL - Sludge SED - Sediment CI - Oil LIQ - Other Liquid AIR - Air SOL - Other Solid WP - Wipe FB - Field Blank EB - Equipment Blank RB - Rinses Blank TB - Trip Blank										LAB USE ONLY							
Street Address <b>2 Huntington Quad, Suite 1S10</b>		City <b>Melville NY</b>		State <b>NY</b>														Zip <b>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>N4014961514.1AV13</b>		Street Address <b>630 Plaza Drive, Suite 600</b>														City <b>Highlands Ranch, CO</b>		State <b>CO</b>		Zip <b>80129</b>	
Phone # <b>631-249-7600</b>		Fax # <b>631-249-7610</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>	
Sampler(s) Name(s) <b>Pat Prusti 547-6247</b>		Collection		Number of preserved bottles																			
Field ID / Point of Collection		MECHNOV #		Date		Time		Sampled by		Matrix		# of bottles		PCI NICK HINCH HURCH NDRE DI Water MESH ENDUSE									
-1 <b>FB 113015002</b>				<b>11/30/15</b>		<b>1310</b>		<b>AP</b>		<b>FB</b>		<b>2</b>											
-2 <b>Box G-1</b>				<b>11/30/15</b>		<b>1635</b>		<b>AP</b>		<b>GW</b>		<b>2</b>											
-3 <b>Box G-2</b>				<b>11/30/15</b>		<b>1635</b>		<b>AP</b>		<b>GW</b>		<b>2</b>											
Turnaround Time (Business days)		Approved By (Accutest PI#): / Date:				Data Deliverable Information										Comments / Special Instructions							
<input type="checkbox"/> SKI, 15 Business Days <input checked="" type="checkbox"/> SKI, 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						<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 LUMMU+		RL reporting for metals  _____ _____ _____					
Emergency & Rush T/A data available via Lablink																							
Sample Custody must be documented below each time samples change possession, including courier delivery.																							
1 Relinquished By: <b>Pat Prusti</b>		Date Time: <b>11/30/15 1915</b>		Received By: <b>1</b>		Relinquished By: <b>Pat</b>		Date Time: <b>12/1/15 1935</b>		Received By: <b>2</b>		Relinquished By: <b>3</b>		Date Time: <b>12/1/15 1935</b>		Received By: <b>4</b>							
3 Relinquished By: <b>Pat Prusti</b>		Date Time: <b>11/30/15 1915</b>		Received By: <b>3</b>		Relinquished By: <b>4</b>		Date Time: <b>12/1/15 1935</b>		Received By: <b>4</b>		Relinquished By: <b>5</b>		Date Time: <b>12/1/15 1935</b>		Received By: <b>5</b>							
5 Relinquished By: <b>Pat Prusti</b>		Date Time: <b>11/30/15 1915</b>		Received By: <b>5</b>		Custody Seal #		<input type="checkbox"/> Intact <input type="checkbox"/> Not Intact		Preserved where applicable		On Ice <input checked="" type="checkbox"/>		Cooler Temp: <b>0.9°C</b>									

5.1  
5

17D

## Report of Analysis

<b>Client Sample ID:</b> FB113015PP1		
<b>Lab Sample ID:</b> JC9566-1		<b>Date Sampled:</b> 11/30/15
<b>Matrix:</b> AQ - Field Blank Water		<b>Date Received:</b> 12/01/15
<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	1B100892.D	1	12/03/15	MD	n/a	n/a	V1B4776
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	2.2	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> FB113015PP1		<b>Date Sampled:</b> 11/30/15
<b>Lab Sample ID:</b> JC9566-1		<b>Date Received:</b> 12/01/15
<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	104%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	unknown	7.61	.65	ug/l	J N
	Total TIC, Volatile		.65	ug/l	J N

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

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

4.1  
4



## Report of Analysis

<b>Client Sample ID:</b>	BPOW 6-1	<b>Date Sampled:</b>	11/30/15
<b>Lab Sample ID:</b>	JC9566-2	<b>Date Received:</b>	12/01/15
<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		

Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B100893.D	1	12/03/15	MD	n/a	n/a	V1B4776
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 6-1		<b>Date Sampled:</b> 11/30/15
<b>Lab Sample ID:</b> JC9566-2		<b>Date Received:</b> 12/01/15
<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	101%		78-114%
460-00-4	4-Bromofluorobenzene	102%		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> 11/30/15
<b>Lab Sample ID:</b> JC9566-3		<b>Date Received:</b> 12/01/15
<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	103%		78-114%
460-00-4	4-Bromofluorobenzene	104%		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> TB113015PP1		<b>Date Sampled:</b> 11/30/15
<b>Lab Sample ID:</b> JC9566-4		<b>Date Received:</b> 12/01/15
<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		

Run #1	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1B100895.D	1	12/03/15	MD	n/a	n/a	V1B4776
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> TB113015PP1		<b>Date Sampled:</b> 11/30/15
<b>Lab Sample ID:</b> JC9566-4		<b>Date Received:</b> 12/01/15
<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	102%		78-114%
460-00-4	4-Bromofluorobenzene	104%		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  
4

## Report of Analysis

<b>Client Sample ID:</b> FB113015PP1	<b>Date Sampled:</b> 11/30/15
<b>Lab Sample ID:</b> JC9566-1	<b>Date Received:</b> 12/01/15
<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>	I99338.D	1	12/10/15	AMA	12/02/15	M:OP45603	M:MSI3713
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	67%		26-121%		
321-60-8	2-Fluorobiphenyl	61%		28-107%		
1718-51-0	Terphenyl-d14	75%		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

4.1  
4

## Report of Analysis

<b>Client Sample ID:</b> BPOW 6-1	<b>Date Sampled:</b> 11/30/15
<b>Lab Sample ID:</b> JC9566-2	<b>Date Received:</b> 12/01/15
<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>	I99339.D	1	12/10/15	AMA	12/02/15	M:OP45603	M:MSI3713
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	72%		26-121%		
321-60-8	2-Fluorobiphenyl	64%		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

4.2  
4



## Report of Analysis

<b>Client Sample ID:</b> BPOW 6-2	<b>Date Sampled:</b> 11/30/15
<b>Lab Sample ID:</b> JC9566-3	<b>Date Received:</b> 12/01/15
<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>	I99340.D	1	12/10/15	AMA	12/02/15	M:OP45603	M:MSI3713
Run #2							

Run #	Initial Volume	Final Volume
Run #1	890 ml	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
123-91-1	1,4-Dioxane	ND	0.22	0.085	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	80%		26-121%		
321-60-8	2-Fluorobiphenyl	71%		28-107%		
1718-51-0	Terphenyl-d14	76%		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





**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> 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>631-249-7600 631-249-7610</b>		Project Name: <b>AGMNYM62235 // OU2 Monitoring 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> Work Authorization #: <b>NY001496_2015</b> Project Manager <b>Carlo San Giovanni</b>		Bottle Order Control # <b>#5</b> Accutest Quote # <b>JC9689</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 WIP - Wipe FB-Field Blank EB-Equipment Blank RB-Rinsate Blank TB-Trip Blank			
Samples (Name(s)) <b>FB 12.01.15 PPA</b> <b>BPW G-3</b> <b>BPW G-4</b> <b>Rep 12.01.15 PPA</b>		Attention: <b>Soma Das</b>		B8270 SIM14DIOX ✓ ✓ ✓		LAB USE ONLY			
Account Sample #	Field ID / Point of Collection	MEOHDI Val #	Date			Time	Sampled by	Matrix	# of bottles
1	FB 12.01.15 PPA		12/1/15			1050	CP	FB	2
2	BPW G-3		12/1/15			1350	AKH	GW	2
3	BPW G-4		12/1/15			1349	CP	GW	2
4	Rep 12.01.15 PPA		12/1/15		CP	GW	2		
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 TIA data available VIA Lablink		<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 LUM/MUC+		RL reporting for metals			
Sample Custody must be documented below each time samples change possession, including courier delivery.									
Relinquished by:	Date Time:	Received By:	Date Time:	Relinquished by:	Date Time:	Received By:	Date Time:		
1	12/1/15 1810	1		2		2			
Relinquished by:	Date Time:	Received By:	Date Time:	Relinquished by:	Date Time:	Received By:	Date Time:		
3	12-2-15 1045	3		4		4			
Relinquished by:	Date Time:	Received By:	Date Time:	Custody Seal #	<input type="checkbox"/> Intact <input type="checkbox"/> Not Intact	Preserved where applicable	On Ice <input type="checkbox"/> Cooler Temp. <input type="checkbox"/>		
5		5					IRGWI		

5.1  
5



## Report of Analysis

<b>Client Sample ID:</b> FB120115PP1		<b>Date Sampled:</b> 12/01/15
<b>Lab Sample ID:</b> JC9689-1		<b>Date Received:</b> 12/02/15
<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	1B100883.D	1	12/03/15	MD	n/a	n/a	V1B4776
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	2.7	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> FB120115PP1		<b>Date Sampled:</b> 12/01/15
<b>Lab Sample ID:</b> JC9689-1		<b>Date Received:</b> 12/02/15
<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	104%		78-114%
460-00-4	4-Bromofluorobenzene	105%		77-115%

CAS No.	Tentatively Identified Compounds	R. T.	Est. Conc.	Units	Q
	alkene	8.25	2	ug/l	J N
	Total TIC, Volatile		2	ug/l	J N

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

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

4.1  
4



## Report of Analysis

<b>Client Sample ID:</b> BPOW 6-3		<b>Date Sampled:</b> 12/01/15
<b>Lab Sample ID:</b> JC9689-2		<b>Date Received:</b> 12/02/15
<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	102%		78-114%
460-00-4	4-Bromofluorobenzene	103%		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 6-4		<b>Date Sampled:</b> 12/01/15
<b>Lab Sample ID:</b> JC9689-3		<b>Date Received:</b> 12/02/15
<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	102%		78-114%
460-00-4	4-Bromofluorobenzene	103%		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> REP120115PP1		<b>Date Sampled:</b> 12/01/15
<b>Lab Sample ID:</b> JC9689-4		<b>Date Received:</b> 12/02/15
<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	103%		78-114%
460-00-4	4-Bromofluorobenzene	103%		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

4.4  
4



## Report of Analysis

<b>Client Sample ID:</b> TB120115PP1		<b>Date Sampled:</b> 12/01/15
<b>Lab Sample ID:</b> JC9689-5		<b>Date Received:</b> 12/02/15
<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	102%		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

4.5  
4

## Report of Analysis

<b>Client Sample ID:</b> FB120115PP1		<b>Date Sampled:</b> 12/01/15
<b>Lab Sample ID:</b> JC9689-1		<b>Date Received:</b> 12/02/15
<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>	I99290.D	1	12/04/15	AMA	12/03/15	M:OP45613	M:MSI3710
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	74%		26-121%		
321-60-8	2-Fluorobiphenyl	74%		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

4.1  
4

## Report of Analysis

<b>Client Sample ID:</b> BPOW 6-3		<b>Date Sampled:</b> 12/01/15
<b>Lab Sample ID:</b> JC9689-2		<b>Date Received:</b> 12/02/15
<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>	I99291.D	1	12/04/15	AMA	12/03/15	M:OP45613	M:MSI3710
Run #2							

Run #	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	83%		26-121%		
321-60-8	2-Fluorobiphenyl	78%		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 6-4		<b>Date Sampled:</b> 12/01/15
<b>Lab Sample ID:</b> JC9689-3		<b>Date Received:</b> 12/02/15
<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>	I99292.D	1	12/04/15	AMA	12/03/15	M:OP45613	M:MSI3710
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	80%		26-121%		
321-60-8	2-Fluorobiphenyl	76%		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  
4



## Report of Analysis

<b>Client Sample ID:</b> REP120115PP1	<b>Date Sampled:</b> 12/01/15
<b>Lab Sample ID:</b> JC9689-4	<b>Date Received:</b> 12/02/15
<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>	I99293.D	1	12/04/15	AMA	12/03/15	M:OP45613	M:MSI3710
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	76%		28-107%		
1718-51-0	Terphenyl-d14	87%		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

**Appendix D**

**ARCADIS Separate and Ongoing OU2 Monitoring of Navy Wells**

Well	Well Owner	1st Q	2nd Q	3rd Q	4th Q	VOC Analysis Method
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**Outpost wells**

BPOW1-1	Navy		X		X	524.2
BPOW1-2	Navy		X		X	524.2
BPOW1-3	Navy		X		X	524.2
BPOW1-4	Navy		X		X	524.2
BPOW1-5	Navy		X		X	524.2
BPOW1-6	Navy		X		X	524.2
BPOW2-1	Navy		X		X	524.2
BPOW2-2	Navy		X		X	524.2
BPOW2-3	Navy		X		X	524.2
BPOW3-1	Navy		X		X	524.2
BPOW3-2	Navy		X		X	524.2
BPOW3-3	Navy		X		X	524.2
BPOW3-4	Navy		X		X	524.2

**Semi-annual and annual**

TT102D	Navy		X		X	8260C
TT102D2	Navy		X		X	8260C
FW-03	Navy		X			8260C
GM-15D	Navy		X		X	8260C
GM-15D2	Navy		X		X	8260C
GM-17D	Navy		X		X	8260C
GM-17I	Navy		X		X	8260C
GM-18D	Navy		X		X	8260C
GM-21D	Navy		X			8260C
GM-39DA	Navy		X		X	8260C
GM-39DB	Navy		X		X	8260C
GM-73D	Navy		X		X	8260C
GM-73D2	Navy		X		X	8260C
GM-74D	Navy		X		X	8260C
GM-74I	Navy		X		X	8260C
GM-75D2	Navy		X		X	8260C
GM-78I	Navy		X			8260C
GM-78S	Navy		X			8260C
GM-79D	Navy		X		X	8260C
GM-79I	Navy		X		X	8260C
HN-24I	Navy		X			8260C
HN-40I	Navy		X			8260C
HN-40S	Navy		X			8260C
HN-42I	Navy		X			8260C
HN-42S	Navy		X			8260C

Q: Quarter

VOC: volatile organic compound