*Pyramid Environmental & Engineering, P.C. Project # 2014-093 Preliminary Site Assessment (PSA) – Parcel 008, Kenneth M. Taylor* 

#### PRELIMINARY SITE ASSESSMENT PARCEL 008 – KENNETH M. TAYLOR 675 CUMMINS DRIVE KENLY, JOHNSTON COUNTY, NORTH CAROLINA NC PIN: 264600-93-9514 STATE PROJECT: I-3318BB WBS ELEMENT: 34182.2.1 JUNE 27, 2014

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C-257 –Geology C-1251 – Engineering

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### Acronyms

BLS	.Below Land Surface
BTEX	.Benzene, Toluene, Ethylbenzene, & Xylenes
CADD	.Computer Aided Design and Drafting
COC	.Chain of Custody
CSA	.Comprehensive Site Assessment
DENR	.Department of Environment and Natural Resources
DRO	.Diesel Range Organics
DWM	.Division of Waste Management
EM	.Electromagnetic (as with EM-61)
EPA	.Environmental Protection Agency
GRO	.Gasoline Range Organics
GCLs	.Gross Contaminant Levels
GPR	.Ground Penetrating Radar
HASP	.Health & Safety Plan
MSCC	.Maximum Soil Contaminant Concentration
MTBE	.Methyl Tertiary Butyl Ether
μg/L	.Micrograms per Liter
mg/kg	.Milligrams per kilogram
NPDES	.National Pollutions Discharge Elimination System
NCAC	North Carolina Administrative Code
NCDOT	North Carolina Department of Transportation
OSHA	.Occupational Safety and Health Administration
OVA	.Organic Vapor Analyzer
PPM	.Parts Per Million
PID	.Photo-ionization Detector
PSA	.Preliminary Site Assessment
PVC	.Poly-vinyl Chloride
RFP	.Request for Proposal
ROW	.Right of Way
SVOCs	.Semi-volatile Organic Compounds
TW	.Temporary Well
TPH	.Total Petroleum Hydrocarbons
UVF	.Ultraviolet Fluorescence (UVF) QED Analyzer
UST	.Underground Storage Tank
US EPA	.United States Environmental Protection Agency
VOCs	.Volatile Organic Compounds

#### PRELIMINARY SITE ASSESSMENT PARCEL 008, KENNETH M. TAYLOR 675 CUMMINS DRIVE KENLY, JOHNSTON COUNTY, NORTH CAROLINA

### **EXECUTIVE SUMMARY OF RESULTS**

Pyramid Environmental & Engineering P.C. (Pyramid) has prepared this Preliminary Site Assessment (PSA) report documenting background information, field activities, assessment activities, findings, conclusions, and recommendations for Parcel 008, Kenneth M. Taylor. The purpose of this assessment was to determine the presence or absence of underground storage tanks (USTs) and impacted soils between the existing edge of pavement and the proposed ROW and/or easements, whichever distance was greater. This PSA is a part of State Project I-3318BB. The PSA was conducted with particular attention to the areas to be cut as indicated by slope stake lines and cross sections or to be excavated for the installation of drainage features. This preliminary site assessment was conducted on behalf of the North Carolina Department of Transportation (NCDOT) in accordance with Pyramid's April 23, 2014, technical proposal.

The following statements summarize the results of the PSA:

• Site History: On May 6, 2014, Pyramid emailed the Johnston County I-3318BB parcel address (675 Cummins Drive in Kenly, NC) to Mr. Jeremy Poplawski, Johnston County Incident Manager, with the Fayetteville Regional Office for the DENR UST Section, with a request to investigate any environmental incidents associated with the parcel. On May 7, 2014, Mr. Poplawski responded to the email and stated that site address did not come back with any incidents or files. He also did not find anything related to the suspected methamphetamine lab.

On May 13, 2014, Pyramid Project Manager Eric Cross performed a site visit at the property. Neither the owner nor any other personnel were present to conduct an interview. A large billboard was present on the property, and the remaining area was undeveloped land. No evidence of USTs was observed. Subsequent site visits provided evidence that it is most likely Parcel 5 (directly to the west of Parcel 8) is being used as a methamphetamine lab.

• **Geophysical Survey:** A large portion of the parcel was inaccessible due to dense/tall vegetation and forest. Two of the EM61 anomalies detected could be attributed to visible objects at the ground surface; specifically, a mailbox and a reinforced concrete drainage pipe. The remaining EM features were suspected to be associated with metallic debris, and were investigated using the GPR. The

GPR did not record any significant reflectors that would be indicative of structures such as USTs. The GPR data were consistent with areas of metallic debris or utilities. The geophysical investigation <u>did not record any evidence of metallic USTs</u> at the property.

• Limited Soil Assessment: A total of four borings were performed across the property. The DENR action levels for both TPH-GRO and TPH-DRO are 10 milligrams per kilogram (mg/kg). Soil samples were screened with a PID, and select soil samples were analyzed for DRO and GRO using a QED Analyzer. None of the soil samples analyzed exhibited DRO or GRO concentrations above 10 mg/kg.

One soil sample 8-1(4-6) was sent to the laboratory for analysis of soils using EPA Methods 8260/8270 for volatile organic compounds (VOC) and semivolatile organic compounds (SVOC) based on the suspected methamphetamine lab located at the west-adjacent parcel. The laboratory results did not detect any concentrations of VOCs or SVOCs above residential or soil-to-groundwater MSCCs in the samples that were analyzed.

• Limited Groundwater Assessment: Soil boring 8-4 was converted into a 1-inch diameter temporary monitoring well to a total depth of 12 feet BLS. The depth-to-groundwater was measured at 3.35 feet BLS. The 6200B and 625 laboratory analysis of groundwater samples did not detect any VOC or SVOC concentrations above NCAC 2L groundwater standards.

Review of the NCDOT engineering plans indicates that the NCDOT may encounter groundwater at the property during construction activities. The results of this PSA do not indicate that any contamination is present in the groundwater at the location of sampling.

• **Contaminated Soil Volumes:** No evidence of petroleum-impacted soils (DRO/GRO > 10mg/kg) was observed during this investigation. Additionally, no evidence of elevated concentrations of VOCs or SVOCs was detected by laboratory analysis of the soil samples. Therefore, no recommendations for the treatment, handling, or disposal of such materials are warranted.

### **1.0 Introduction**

Pyramid Environmental & Engineering P.C. (Pyramid) has prepared this Preliminary Site Assessment (PSA) report documenting background information, field activities, assessment activities, findings, conclusions, and recommendations for Parcel 008, Kenneth M. Taylor. The Kenneth M. Taylor property contains a commercial billboard sign and is otherwise undeveloped. The parcel is located at approximately 675 Cummins Drive, Kenly, NC. This preliminary site assessment was conducted on behalf of the North Carolina Department of Transportation (NCDOT) in accordance with Pyramid's April 23, 2014, technical proposal. This PSA is a part of State Project I-3318BB.

The purpose of this assessment was to determine the presence or absence of underground storage tanks (USTs) and impacted soils between the existing edge of pavement and the proposed ROW and/or easements, whichever distance was greater. The PSA was conducted with particular attention to the areas to be cut as indicated by slope stake lines and cross sections or to be excavated for the installation of drainage features between the existing edge of pavement and proposed ROW/easements. The location of the subject site is shown on **Figure 1**.

#### **<u>1.1 Background Information</u>**

Based on the NCDOT's April 15, 2014, *Request for Technical and Cost Proposal*, the PSA was conducted between the existing edge of pavement and the proposed ROW and/or easements, whichever distance was greater, with emphasis on the areas to be cut as indicated by slope stake lines and cross sections or to be excavated for the installation of drainage features and/or other utilities, in accordance with the CADD files provided to Pyramid by the NCDOT. The PSA included the following:

- Research the properties for past uses and possible releases.
- Conduct a preliminary geophysical site assessment and limited soil assessment across the entire parcel with emphasis on the areas to be cut as indicated by slope stake lines and cross sections or to be excavated for the installation of drainage features and/or other utilities.
- If a NCDENR Groundwater Incident has been assigned to a parcel, then a single groundwater sample will be collected (or attempted) from the parcel if groundwater is encountered in any of the soil borings on that parcel incidentally during the course of attaining the depths required for objective of soil sampling. At parcels without NCDENR assigned Groundwater Incidents, if groundwater is likely to be encountered by subsequent excavation required by construction, then Pyramid will attempt to obtain a groundwater sample from the parcel.

#### **1.2 Project Information**

Prior to field activities, a Health and Safety Plan was prepared. Prior to drilling activities, the public underground utilities were located and marked by the North Carolina One-Call Service. A private utility locator, Northstate Utility Locating Incorporated of Colfax, North Carolina was used to mark the on-site private, buried utilities.

### 2.0 Site History

The NCDOT description of Parcel 008 in the RFP provided to Pyramid on April 15, 2014, provided the following background information related to the site:

"Vacant mobile homes and a junk yard were observed at Parcels 5 & 8 during a site reconnaissance on June 9, 2011. The site(s) is located on the southern side of Cummins Drive, approximately 400 feet southeast of Bridge 114. A local area employee indicated that the property may have operated as a clandestine methamphetamine. Review of crime databases could not confirm the historical presence of a lab. According to NCDENR's UST Section Registry there are no known Facility ID's or Groundwater Incidents associated with Parcel 8."

Pyramid interviewed DENR personnel, interviewed property owners, and reviewed aerial photographs to assess past uses of the property. Pyramid reviewed historical aerial photographs obtained from the Johnston County GIS website and Google Earth dating back to 1937. The 1937, 1948, 1956, 1971, 1988, 1993, 1999, 2004, 2006, 2008, 2009 and 2012 aerial photographs are included in **Appendix A**. Historical information reviewed as part of the PSA indicated that the Kenneth M. Taylor property was first developed between 1988 and 1993. The 1993 aerial photograph shows that the western portion of the parcel had been cleared, and scattered trailers and other structures/vehicles were present that may have been associated with the apparent junkyard directly to the west at Parcel 5. The 1988 aerial shows the parcel to be undeveloped. The historical aerials indicate that the trailers and vehicles had been removed from Parcel 8 between 2006 and 2008, and no other structures or buildings were observed in the subsequent aerial photographs.

On May 6, 2014, Pyramid emailed the Johnston County I-3318BB parcel address (675 Cummins Drive in Kenly, NC) to Mr. Jeremy Poplawski, Johnston County Incident Manager, with the Fayetteville Regional Office for the DENR UST Section, with a request to investigate any environmental incidents associated with the parcel. On May 7, 2014, Mr. Poplawski responded to the email and stated that site address did not come back with any incidents or files. He also did not find anything related to the suspected methamphetamine lab.

On May 13, 2014, Pyramid Project Manager Eric Cross performed a site visit at the property. Neither the owner nor any other personnel were present to conduct an interview. A large billboard was present on the property, and the remaining area was undeveloped land. No evidence of USTs was observed. Subsequent site visits provided evidence that it is likely Parcel 5, directly to the west of Parcel 8, was reportedly being used as a methamphetamine lab (see Pyramid's Parcel 5 PSA Report).

## **3.0 Geophysical Investigation**

Pyramid's classifications of USTs for the purposes of this PSA report are based directly on the geophysical UST ratings provided to us by the NCDOT. These ratings are as follows:

	Geophysical Surveys for on NCI	Underground Stora	ge Tanks
High Confidence	Intermediate Confidence	Low Confidence	No Confidence Anomaly noted but not
Active tank - spatial location, orientation, and approximate depth determined by geophysics.	Sufficient geophysical data from both magnetic and radar surveys that is characteristic of a tank. Interpretation may be supported by physical evidence such as fill/vent pipe, metal cover plate, asphal/concrete patch, etc.	Sufficient geophysical data from either magnetic or radar surveys that is characteristic of a tank. Additional data is not sufficient enough to confirm or deny the presence of a UST.	characteristic of a UST. Should be noted in the text and may be called out in the figures at the geophysicist's discretion.

Pyramid performed electromagnetic (EM) and ground penetrating radar (GPR) surveys across the <u>accessible</u> portions of the Parcel. Two of the EM61 anomalies detected could be attributed to visible objects at the ground surface; specifically, a mailbox and a reinforced concrete drainage pipe. The remaining EM features were suspected to be associated with metallic debris, and were investigated by the GPR. The GPR did not record any significant reflectors that would be indicative of structures such as USTs. The GPR data were consistent with areas of metallic debris or utilities. The geophysical investigation <u>did not record any evidence of metallic USTs</u> at the property.

The full details of the geophysical investigation are included in the Geophysical Investigation Report as **Appendix B**.

### 4.0 Soil Sampling Activities & Results

#### 4.1 Soil Assessment Field Activities

On June 2, 2014, Pyramid mobilized to the site, drilled soil borings and collected the proposed soil samples for the PSA. Four (4) soil borings (8-1 through 8-4) were advanced on the subject property between the NCDOT proposed ROW and easements,

and edge of pavement. The soil borings were completed using a truck mounted GeoProbe drill rig. The selected locations were chosen to avoid public and private utilities while remaining in the proposed right of way and/or easement.

The soil borings were installed at or adjacent to proposed drainage features, as indicated by the NCDOT engineering plans, or generally within the proposed ROW and/or easement to obtain additional information. The locations of the borings are shown on **Figure 2**.

Soil samples were continuously collected in four-foot long disposable sleeves from each boring for geologic description, and visual examination for signs of contamination. Soil recovered from each sleeve was screened in the field using a Photo-Ionization Detector (PID) approximately every 2 feet depending on the soil recovery of each sleeve. In general, the soil sample with the highest PID reading was selected from each boring for laboratory analysis. If field screening detected an elevated reading, then additional soil samples from each boring were selectively analyzed with the QED UVF HC-1 Analyzer. The soil boring logs with the soil descriptions, visual examination, and PID screening results are included in **Appendix C**. The PID field screening results are summarized in **Table 1**. To prevent cross contamination, new disposable nitrile gloves were worn by the sampling technician during the sampling activities, and were changed between samples. No petroleum odor was detected in any of the borings during the field screening.

The soil samples selected for Total Petroleum Hydrocarbon (TPH) analyses were analyzed utilizing the QED UVF HC-1 Analyzer system from QROS-US. The NCDOT has indicated that this instrument is an acceptable method to provide total petroleum hydrocarbon (TPH) results for soil analysis for the PSA projects. Pyramid's QEDcertified technician performed the soil analyses. The soil samples selected for analysis using the QED Analyzer were analyzed for TPH as diesel range organics (DRO) and TPH as gasoline range organics (GRO). The soil samples selected for analysis using the QED were preserved in the field with methanol and were analyzed at the end of each day using the QED.

In addition to the QED analysis, select samples were collected for more comprehensive laboratory analysis using EPA Methods 8260 and 8270 for volatile and semi-volatile organic compounds, respectively. These additional analyses were performed based on the site history of the property, which suggested that other potential contaminants such as solvents may have been or are being utilized in association with a possible methamphetamine lab directly to the west of Parcel 008. Soils from the boring located nearest to Parcel 005, where the lab was suspected to be located, were selected for the additional laboratory analyses. Specifically, sample 8-1(4-6) was placed in a laboratory prepared containers and shipped to Pace Analytical in Huntersville, NC for analysis of volatile and semi-volatile organic compounds. It should be noted that a typographical error was made on the Chain of Custody form submitted to Pace Analytical, and the

sample ID number in their lab report is listed as sample 8-2(4-6). The actual sample that was analyzed was 8-1(4-6).

#### 4.2 Soil Sample Analytical Results

#### QED Results

The DENR action levels for both TPH-GRO and TPH-DRO are 10 mg/kg. Soil samples were screened with a PID, and select soil samples were analyzed for DRO and GRO using a QED Analyzer. None of the soil samples analyzed exhibited DRO or GRO concentrations above 10 mg/kg. The soil sample QED results are summarized in **Table 2**. A copy of the QED analysis report is included in **Appendix D**.

#### Laboratory Analysis for Methods 8260/8270

One soil sample 8-1(4-6) was sent to the laboratory for analysis of soils using EPA Methods 8260/8270 for volatile (VOC) and semi-volatile (SVOC) organic compounds based on the suspected methamphetamine lab located at the west-adjacent parcel. The laboratory results did not detect any concentrations of VOCs or SVOCs above residential or soil-to-groundwater MSCCs in the samples that were analyzed. The soil sample laboratory results are summarized in **Table 3**. A copy of the laboratory report and chain-of-custody is included in **Appendix E**.

#### 4.3 Temporary Monitoring Well Installation

On June 2, 2014, Pyramid converted soil boring 8-4 into a 1-inch diameter temporary monitoring well (TW). This location was chosen based on PID and QED readings, and its location at a proposed drainage structure. Soil boring 8-4(TW) was completed to a total depth of 12 feet below land surface (BLS). The temporary well was constructed with 2 feet of 1-inch diameter schedule 80 PVC casing and 10 feet 1-inch diameter of schedule 80 PVC slotted screen.

The depth-to-groundwater was measured at 3.35 feet BLS. The temporary monitoring well was sampled using a new 0.5-inch diameter disposable bailer. Upon completion of the gauging and sampling, the temporary monitoring well was properly abandoned by the drillers by removing the casing, and filling the borehole with bentonite chips and portland cement.

#### 4.4 Groundwater Analytical Results

The groundwater sample 8-4(TW) was placed in laboratory prepared containers for analysis of VOCs using EPA Method 6200B and for SVOCs using EPA Method 625. The samples were shipped to Pace Analytical in Huntersville, NC. The 6200B and 625 laboratory analysis did not detect any concentrations of VOCs or SVOCs above NCAC 2L groundwater standards in the sample. The groundwater results for sample 8-4(TW) are summarized in **Table 4**. A copy of the laboratory report and chain-of-custody is included in **Appendix E**.

### **5.0** Conclusions and Recommendations

As requested by NCDOT, Pyramid has completed a PSA at the Kenneth M. Taylor property located at 675 Cummins Drive, Kenly, NC (Parcel 008). The following is a summary of the assessment activities and results. Personnel logs for all field work are included in **Appendix F.** 

#### 5.1 Geophysical Investigation

A large portion of the parcel was inaccessible due to dense/tall vegetation and forest. Two of the EM61 anomalies detected could be attributed to visible objects at the ground surface; specifically, a mailbox and a reinforced concrete drainage pipe. The remaining EM features were suspected to be associated with metallic debris, and were investigated using the GPR. The GPR did not record any significant reflectors that would be indicative of structures such as USTs. The GPR data were consistent with areas of metallic debris or utilities. The geophysical investigation <u>did not record any evidence of metallic USTs</u> at the property.

#### 5.2 Limited Soil Assessment

#### QED Results

The DENR action levels for both TPH-GRO and TPH-DRO are 10 mg/kg. Soil samples were screened with a PID, and select soil samples were analyzed for DRO and GRO using a QED Analyzer. None of the soil samples analyzed exhibited DRO or GRO concentrations above 10 mg/kg.

#### Laboratory Results

One soil sample 8-1(4-6) was sent to the laboratory for analysis of soils using EPA Methods 8260 and 8270 for volatile (VOC) and semi-volatile (SVOC) organic compounds, respectively, based on the suspected methamphetamine lab located at the west-adjacent parcel. The laboratory results did not detect any concentrations of VOCs or SVOCs above residential or soil-to-groundwater MSCCs in the samples that were analyzed.

#### 5.3 Limited Groundwater Assessment

Soil boring 8-4 was converted into a 1-inch diameter temporary monitoring well to a total depth of 12 feet BLS. The depth-to-groundwater was measured at 3.35 feet BLS. The 6200B and 625 laboratory analysis did not detect any VOCs or SVOCs exhibiting concentrations above NCAC 2L groundwater standards in the sample.

Review of the NCDOT engineering plans indicates that the NCDOT may encounter groundwater at the property during construction activities. The results of this PSA do not indicate that any contamination is present in the groundwater at the location of sampling.

#### 5.4 Recommendations

Petroleum-Impacted Soils

No evidence of petroleum-impacted soils (DRO/GRO > 10 mg/kg) was observed during this investigation. Additionally, no evidence of elevated concentrations of VOCs or SVOCs was detected by laboratory analysis of the soil samples. Therefore, no recommendations for the treatment, handling, or disposal of such materials are warranted.

It should be noted that, if impacted soil is encountered during road construction outside of the area analyzed by this investigation, the impacted soil should be managed according to NC DENR Division of Waste Management (DWM) Guidelines and disposed of at a permitted facility.

### 6.0 Limitations

The results of this preliminary investigation are limited to the boring locations completed during this limited assessment and presented in this report. The laboratory results only reflect the current conditions at the locations sampled on the date this PSA was performed.

### 7.0 Closure

This report was prepared for, and is available solely for use by NCDOT and their designees. The contents thereof may not be used or relied upon by any other person without the express written consent and authorization of Pyramid Environmental & Engineering, P.C. (Pyramid). The observations, conclusions, and recommendations documented in this report are based on site conditions and information reviewed at the time of Pyramid's investigation. Pyramid appreciates the opportunity to provide this environmental service.

**FIGURES** 





—	PROPOSED UTILITY EASEMENT
	EXISTING ROW
	EXISTING PROPERTY BOUNDARY
<del>)</del>	PROPOSED ROW
	PROPOSED CONST. EASEMENT
	PROP. DRAINAGE UTIL. EASEMENT
	PROPOSED SS CUT LINE
	PROPOSED SS FILL LINE
	PROPOSED SS TRANSITION LINE
	PROPOSED DRAINAGE PIPING
_	





TABLES

#### TABLE 1

#### Summary of Soil Field Screening Results NCDOT Project I-3318BB 675 Cummins Drive - Parcel 008 Kenly, Johnston County, North Carolina

SOIL BORING	SAMPLE ID	DEPTH	PID
		(feet bgs)	READINGS (PPM)
	8-1(0-2)	0 to 2	1.5
	8-1(2-4)	2 to 4	2.0
8-1	8-1(4-6)	4 to 6	3.0
	8-1(6-8)	6 to 8	2.0
	8-1(8-11)	8 to 11	1.0
	8-2(0-2)	0 to 2	1.0
8-2	8-2(2-4)	2 to 4	2.0
	8-2(4-6)	4 to 6	3.0
	8-2(6-8)	6 to 8	2.0
	8-2(8-11)	8 to 11	2.0
	8-3(0-2)	0 to 2	1.0
8-3	8-3(2-4)	2 to 4	2.5
	8-3(4-6)	4 to 6	2.0
	8-3(6-8)	6 to 8	2.0
	8-4(0-2)	0 to 2	2.0
	8-4(2-4)	2 to 4	3.0
8-4	8-4(4-6)	4 to 6	2.0
	8-4(6-8)	6 to 8	2.0
	8-4(8-10)	8 to 10	13.0
	8-4(10-12)	10 to 12	17.0

bgs= below ground surface

PID= photo-ionization detector

PPM= parts-per-million

= sampled for lab analysis &/or QROS-QED analysis

OVA= Organic Vapor Analyzer

### TABLE 2

#### Summary of Soil Sample QED Analytical Results for GRO/DRO NCDOT State Project I-3318BB 675 Cummins Drive - Parcel 008 Kenly, Johnston County, North Carolina

				QROS - QED Analysis			
SAMPLE ID	DATE	DEPTH (feet)	PID (ppm)	GRO (mg/kg) (C5-C10)	DRO (mg/kg) (C10-C35)	TPH (mg/kg) (C5-C35)	
	- /- / /						
8-1(4-6)	6/2/2014	4 to 6	3.0	<0.7	0.53	0.53	
8-2(4-6)	6/2/2014	4 to 6	3.0	<0.6	0.17	0.17	
8-3(2-4)	6/2/2014	2 to 4	2.5	<0.6	0.49	0.49	
8-4(2-4)	6/2/2014	2 to 4	3.0	<0.6	0.33	0.33	
NC Initial	Action Level	<ul> <li>UST Secti</li> </ul>	on for				
503	5/5030-GRO;	3550-DRO		10	10	NA	
PID=	photo-ionizaton	detector	GRO=	Gasoline Range Organics	TPH= Total Petroleum	NA=	Not Applicable
PPM=	parts-per-million		DRO= ma/ka=	Diesel Range Organics milligrams-per-kilogram	Hydrocarbons (GRO + DRO)	"" =	No Laboratory

\* Bold values indicate concentrations above initial action levels

#### TABLE 3 Summary of Volatile/Semi-Volatile Laboratory Results of Soil Samples NCDOT Project I-3318BB 675 Cummins Drive - Parcel 008

Analytical	Analytical	SAMPLE ID NUMBER		Soil to
Parameter	Method	8-1(4-6)	Residential	Groundwater
	Sample Date:	6/3/2014	MSCC	MSCC
	Depth (feet):	4 to 6	(mg/kg)	(mg/kg)
	Location	N central		
Acetone	8260	ND	14000	24
Benzene	8260	ND	18	0.0056
Bromobenzene	8260	ND	NMSCC	NMSCC
Bromoform	8260	ND	81	0.026
2-Butanone (MEK)	8260	ND	9385	16
n-Butylbenzene	8260	ND	626	4.3
sec-Butylbenzene	8260	ND	626	3.3
Styrene	8260	ND	3128	1.5
tert-Butylbenzene	8260	ND	626	3.4
4-Chlorotoluene	8260	ND	1000	0.1
Ethylbenzene	8260	ND	1560	4.9
1,2-Dichloroethane	8260	ND	7	0.0019
Isopropyl ether (IPE)	8260	ND	156	0.37
Isopropylbenzene	8260	ND	1564	1.7
P-Isopropyltoluene	8260	ND	NMSCC	NMSCC
Naphthalene	8260	ND	313	0.16
n-Propylbenzene	8260	ND	626	1.7
Toluene	8260	ND	1200	4.3
1,2,4-Trimethylbenzene	8260	ND	782	8.5
1,3,5-Trimethylbenzene	8260	ND	782	8.3
Total Xylenes	8260	ND	3129	4.6
MTBE	8260	ND	350	0.091
2-Hexanone	8260	ND	70	0.1
Methylene chloride	8260	ND	85	0.02
All Other 8260 Parameters	8260	ND	NA	NA
Acenaphthene	8270	ND	940	8.2
bis(2-Ethylhexyl)phthalate	8270	ND	46	6.6
1-Methylnaphthalene	8270	ND	20	0.004
2-Methylnaphthalene	8270	ND	63	3.6
Naphthalene	8270	ND	313	0.16
All Other 8270 Parameters	8270	ND	NA	NA

mg/kg = parts per million (ppm). BOLD values are above MSCC levels. NS=Not Sampled for Parameter

MSCC = Maximum Soil Contaminant Conc. ND = Not Detected. J= Estimated Concentration NMSCC= No MSCC NA Not Applicable CI= Considered Immobile

#### TABLE 4

#### Summary of Groundwater Analytical Results NCDOT State Project I-3318BB 675 Cummins Drive - Parcel 008 Kenly, Johnston County, North Carolina

		SAMPLE ID	NCAC 2L			
PARAMETER	UNITS		GROUNDWATER			
		8-4(TW)	STANDARD			
EPA Method 6200B VOCs; Sample Collection Date: 6/3/14						
Benzene	ug/L	ND	1			
Chloroform	ug/L	ND	70			
Diisopropyl Ether (IPE)	ug/L	ND	70			
Ethyl Benzene	ug/L	ND	600			
Isopropylbenzene (Cumene)	ug/L	ND	70			
Naphthalene	ug/L	ND	6			
Styrene	ug/L	ND	70			
Toluene	ug/L	ND	600			
Total Xylenes	ug/L	ND	500			
n-Propylbenzene	ug/L	ND	70			
sec-Butylbenzene	ug/L	ND	70			
n-Butylbenzene	ug/L	ND	70			
tert-Butyl methyl ether (MTBE)	ug/L	ND	20			
tert-Butylbenzene	ug/L	ND	70			
1,2,4-Trimethylbenzene	ug/L	ND	400			
1,2-Dichloroethane	ug/L	ND	0.4			
1,3,5-Trimethylbenzene	ug/L	ND	400			
4-Isopropyltoluene	ug/L	ND	25			
All Other Parameters	ug/L	ND	NA			
EPA Method 625 Semi-Volatile Organic Compounds						
Acenaphthene	ug/L	ND	80			
Diethylphthalate	ug/L	ND	6000			
bis(2-Ethylhexyl)phthalate	ug/L	ND	3			
Naphthalene	ug/L	ND	6			
Phenanthrene	ug/L	ND	200			
Phenol	ug/L	ND	30			
Pyrene	ug/L	ND	200			
All Other Parameters	ug/L	ND	NA			

ug/L= micrograms-per-liter

ND= Not Detected at or above adjusted reporting limit.

NA= Not Applicable

Bold values above 2L

### APPENDIX A





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Study Area

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### APPENDIX B



# **GEOPHYSICAL SURVEY**

### PARCEL 008 – KENNETH M. TAYLOR 675 CUMMINS DRIVE, KENLY, NC NCDOT PROJECT I-3318BB (WBS 34182.2.1)

#### KENLY, JOHNSTON COUNTY, NC

JUNE 19, 2014

Report prepared for:

Mr. Gordon Box GeoEnvironmental Project Manager Geotechnical Engineering Unit 1020 Birch Ridge Drive Raleigh, NC 27610

Prepared by:

Eric C. Cross, P.G. NC License #2181

Reviewed by: \_

Douz Canavello

viewed by: \_\_\_\_\_

Douglas A. Canavello, P.G. NC License #1066

503 INDUSTRIAL AVENUE, GREENSBORO, NC 27406 P: 336.335.3174 F: 336.691.0648 C257: GEOLOGY C1251: ENGINEERING
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- Figure 2 Parcel 008 EM61 Differential Results Contour Map
- Figure 3 Parcel 008 Overlay of EM61 Contour Map On Engineering Plans
- Figure 4 Parcel 008 GPR Transect Locations
- Figure 5 Parcel 008 GPR Transect Images

# **EXECUTIVE SUMMARY**

**Project Description:** Pyramid Environmental conducted a geophysical investigation for the North Carolina Department of Transportation (NCDOT), at the Kenneth M. Taylor property, Parcel 008, 675 Cummins Drive, Kenly, Johnston County, NC. The survey was part of an NCDOT Right-of-Way (ROW) investigation (NCDOT Project I-3318BB). The geophysical survey boundaries at the project site were designed to include the portions of the property between the existing edge of pavement and the proposed ROW and easements, whichever distance was greater. The geophysical investigation consisted of electromagnetic (EM) induction-metal detection and ground penetrating radar (GPR) surveys.

**Geophysical Results:** Two of the EM61 anomalies detected could be attributed to visible objects at the ground surface; specifically, a mailbox and a reinforced concrete drainage pipe. The remaining EM features were suspected to be associated with metallic debris, and were investigated by the GPR. The GPR did not record any significant reflectors that would be indicative of structures such as USTs. The GPR data were consistent with areas of metallic debris or utilities. The geophysical investigation did not record any evidence of metallic USTs at the property.

## INTRODUCTION

Pyramid Environmental conducted a geophysical investigation for the North Carolina Department of Transportation (NCDOT), at the Kenneth M. Taylor property, Parcel 008, 675 Cummins Drive, Kenly, Johnston County, NC. The survey was part of an NCDOT Right-of-Way (ROW) investigation (NCDOT Project I-3318BB). The geophysical survey boundaries at the project site were designed to include the portions of the property between the existing edge of pavement and the proposed ROW and easements, whichever distance was greater. The survey grid spanned approximately 440 feet from west to east and a maximum of approximately 120 feet from north to south. Conducted on May 22 and 23, 2014, the geophysical investigation was performed to determine if unknown, metallic underground storage tanks (USTs) were present beneath the survey area.

The site contained vacant lots with billboard signs, an asphalt road, open grassy areas and zones of dense/tall vegetation. It should be noted that significant portions of the parcel that were within the proposed ROW and/or easements were not accessible by the geophysical equipment due to the vegetation. Surveys were performed in all accessible areas. Aerial photographs showing the survey area boundaries and ground-level photographs are shown in **Figure 1**.

## FIELD METHODOLOGY

The geophysical investigation consisted of electromagnetic (EM) induction-metal detection and ground penetrating radar (GPR) surveys. The EM survey was performed on May 22, 2014, using a Geonics EM61 metal detection instrument integrated with a Trimble AG-114 GPS antennae. The integrated GPS system allows the location of the instrument to be recorded in real-time during data collection, resulting in an EM data set that geo-referenced and can be overlain on aerial photographs and CADD drawings. A boundary grid was established around the perimeter of the site and at select interior locations with marks every 10 feet to maintain orientation of the instrument throughout the survey and assure complete coverage of the area.

According to the instrument specifications, the EM61 can detect a metal drum down to a maximum depth of approximately 8 feet. Smaller objects (1-foot or less in size) can be detected to a maximum depth of 4 to 5 feet. The EM61 data were digitally collected at approximately 0.8 foot intervals generally along north-south trending or east-west trending, parallel survey lines spaced five feet apart. The data were downloaded to a computer and reviewed in the field and office using the Geonics NAV61 and Surfer for Windows Version 11.0 software programs.

GPR data were acquired across select EM differential anomalies on May 23, 2014, using a Geophysical Survey Systems, Inc. (GSSI) UtilityScan DF unit with a dual frequency 300/800 MHz antenna. Data were collected generally from east to west and north to south across the property. The GPR data were viewed in real time using a vertical scan of 512 samples, at a rate of 36 scans per foot. GPR data were viewed down to a maximum depth of approximately 4-5 feet, based on an estimated two-way travel time of 8 nanoseconds per foot. GPR Transects across specific anomalies were saved to the hard drive of the DF unit for post-processing and figure generation.

# **DISCUSSION OF RESULTS**

A contour plot of the EM61 differential results obtained across survey area at the property is presented in **Figure 2**. The differential results are obtained from the difference between the top and bottom coils of the EM61 instrument. The differential results focus on the larger metal objects such as drum and UST-size objects and ignore the smaller insignificant metal objects.

**Discussion of EM Anomalies**: The EM response on the west side of the survey area near the entry drive to the adjacent parcel was associated with a mailbox. The EM linear response on the east side of the survey area corresponded to the location of a reinforced concrete drainage pipe crossing under the road. The remaining features in the center of the survey could not be attributed to visible objects at the ground surface, and were investigated further with the GPR.

**Figure 3** provides an overlay of the EM61 contour map on the NCDOT engineering plans for the site to provide a reference of proposed ROW and construction features with the geophysical data.

**Discussion of GPR Survey: Figure 4** presents the locations of the GPR transects performed at the property, and **Figure 5** presents the transect images. GPR Transects 1-6 were performed at various locations across the scattered EM anomalies in the center of the profile. Collectively, the 6 transects did not record any significant reflectors that would be indicative of structures such as USTs. Some minor soil disturbances and discontinuous reflectors were observed that are consistent with buried debris, and some isolated high amplitude reflectors were suggestive of objects or utilities. It should be noted that historical aerial photographs obtained as part of the full Preliminary Site Assessment (PSA) show unidentifiable objects/structures at the location of these EM features in the past. It is likely that these structures have been covered by vegetation/soil and are the source of the EM responses.

The geophysical investigation did not <u>record any evidence of metallic USTs</u> at the property within the survey area limits. It should be re-stated that a significant portion of the parcel was inaccessible due to dense/tall vegetation.

# **SUMMARY & CONCLUSIONS**

Our evaluation of the EM61 and GPR data collected across Parcel 008 in Kenly, North Carolina, provides the following summary and conclusions:

- The EM61 and GPR surveys provided reliable results for the detection of metallic USTs within the accessible portions of the geophysical survey area.
- Two of the EM61 anomalies detected could be attributed to visible objects at the ground surface; specifically, a mailbox and a reinforced concrete drainage pipe.
- The remaining EM features were suspected to be associated with metallic debris, and were investigated by the GPR.
- The GPR did not record any significant reflectors that would be indicative of structures such as USTs. The GPR data were consistent with areas of metallic debris or utilities.
- The geophysical investigation did not <u>record any evidence of metallic USTs</u> at the property.

# LIMITATIONS

Geophysical surveys have been performed and this report prepared for the NCDOT in accordance with generally accepted guidelines for EM61 and GPR surveys. It is generally recognized that the results of the EM61 and GPR surveys are non-unique and may not represent actual subsurface conditions. The EM61 and GPR results obtained for this project have not conclusively determined the definitive presence or absence of metallic USTs, but that the evidence collected is sufficient to result in the conclusions made in this report. Additionally, it should be understood that areas containing extensive vegetation, reinforced concrete, or other restrictions to the accessibility of the geophysical instruments could not be fully investigated.





Approximate Boundaries of Geophysical Survey Area



View of West Portion of Survey Area (Facing Approximately East)



View of East Portion of Survey Area (Facing Approximately West)

TITLE PARCEL 008: EM61 GEOPHYSICAL SURVEY PATH AND SITE PHOTOGRAPHS					
PROJECT NCDOT PROJECT I-3318BB (34182.2.1) KENLY, JOHNSTON COUNTY, NC					
	503 INDUSTRIAL AVENUE GREENSBORO, NC 27460 (336) 335-3174 (p) (336) 691-0648 (f) License # C1251 Eng. / License # C257 Geology				
DATE	6/17/2014		CLIENT NCDOT		
PYRAMID PROJECT #:	2014-093		FIGURE 1		



# **EM61 Differential Results**



# NO EVIDENCE OF METALLIC USTs OBSERVED

The contour plot shows the differential results of the EM61 instrument in millivolts (mV). The differential response focuses on larger, buried metallic objects such as drums and USTs and ignores smaller miscellaneous buried, metal debris. The EM61 data were collected on May 22, 2014, using a Geonics EM61 instrument. Ground penetrating radar (GPR) data were collected on May 23, 2014, using a GSSI UtilityScan DF unit and a dual frequency 300/800 MHz antenna.





TITLE

# PARCEL 008: EM61 DIFFERENTIAL RESULTS CONTOUR MAP

PROJECT

NCDOT PROJECT I-3318BB (34182.2.1)
KENLY, JOHNSTON COUNTY, NC

PYRAMID ENVIRONMENTAL & ENGINEERING, P.C	503 INDUSTRIAL AVENUE GREENSBORO, NC 27460 (336) 335-3174 (p) (336) 691-0648 (f) License # C1251 Eng. / License # C257 Geology
DATE 6/19/2014	CLIENT NCDOT
PYRAMID 2014-093	FIGURE 2





Geophysical Survey Area Overlain on NCDOT Engineering Plans (areas not included in survey are the result of dense/tall vegetation)

TITLE	TITLE PARCEL 008: GEOPHYSICAL SURVEY AREA OVERLAIN ON NCDOT CADD				
PROJECT	PROJECT NCDOT PROJECT I-3318BB (34182.2.1) KENLY, JOHNSTON COUNTY, NC				
PYR	503 INDUSTRIAL AVENUE GREENSBORO, NC 27460 (336) 335-3174 (p) (336) 691-0648 (f) License # C1251 Eng. / License # C257 Geology				
DATE	6/17/2014		CLIENT NCDOT		
PYRAMID PROJECT #:	2014-093		FIGURE 3		

# **GPR Transect Locations**









**GPR TRANSECT 1** 



GPR TRANSECT 2

GPR TRANSECT 3



GPR TRANSECT 4



**GPR TRANSECT 5** 



**GPR TRANSECT 6** 



TITLE PARCEL 008: GPR TRANSECT IMAGES				
PROJECT	NCDOT PRO KENLY, J	DJEC' OHNS	T I-3318BB (34182.2.1) STON COUNTY, NC	
503 INDUSTRIAL AVENUE GREENSBORO, NC 27460 (336) 335-3174 (p) (336) 691-0648 (f) License # C1251 Eng. / License # C257 Geold				
DATE	6/19/2014		CLIENT NCDOT	
PYRAMID PROJECT #:	2014-093		FIGURE 5	

# APPENDIX C

# FIELD DRILLING RECORD

PROJECT NAME: PROJECT NUMBER:	NC DOT I-3318BB Parcel 008 Kenneth M. Taylor 2014-093	C DOT I-3318BB Parcel 008 Kenneth M. Taylor 2014-093 BORING/WELL NO:	
SITE LOCATION:	675 Cummins Drive, Kenly, Johnston County, NC	BORING/WELL LOCATION:	W end of parcel
START DATE:	6/2/14	COMPLETED:	6/2/14
GEOLOGIST:	T. Leatherman	DRILLER:	Solutions, IES
DRILL METHOD:	Geoprobe	SAMPLE METHOD:	Macro-core
BORING DIA:	2-inch	CASING DIA:	None
TOTAL DEPTH:	11 feet	CASING DEPTH:	None

	VISUAL MANUAL SOIL CLASSIFICATION	OVA RESULTS
DEPTH	COLOR, TEXTURE, STRUCTURE, CONSISTENCY, ODOR, ETC.	PERCENT RECOVERY
(ft.)		BLOW COUNTS

	Depths correspond to changes in soil type	
0-2'	clayey sandy silt (ML); brown, moist, no odor	PID=8-1(0-2): 1.5 PPM
2-4'	clayey sandy silt (ML); brown, moist, no odor	PID=8-1(2-4): 2 PPM
4-6'	sandy silty clay (CL); brown to tan, moist, no odor	PID=8-1(4-6): 3 PPM
6-8'	sandy silty clay (CL); brown to tan, moist, no odor	PID=8-1(6-8): 2 PPM
8-11'	clayey silty sand (SM) to clayey sand (SC); brown to tan, rock frags	PID=8-1(8-11): 1 PPM
	refusal at 11', groundwater encountered at 5.3 feet	

RISER LENGTH (ft)	DEPTH (ft)	DIAMETE	R (in)	MATERIAL	
SCREEN LENGTH (ft)	DEPTH (ft)	DIAMETE	R (in)	MATERIAL	
DEPTH TO TOP OF SAND		BAGS OF	SAND		
DEPTH TO TOP SEAL	B	ENTONITE USED		BAGS OF CEMENT	USED

# FIELD DRILLING RECORD

PROJECT NAME: PROJECT NUMBER:	NC DOT I-3318BB Parcel 008 Kenneth M. Taylor 2014-093	BORING/WELL NO:	8-2
SITE LOCATION:	675 Cummins Drive, Kenly, Johnston County, NC	BORING/WELL LOCATION:	W-central near road
START DATE:	6/2/14	COMPLETED:	6/2/14
GEOLOGIST:	T. Leatherman	DRILLER:	Solutions, IES
DRILL METHOD:	Geoprobe	SAMPLE METHOD:	Macro-core
BORING DIA:	2-inch	CASING DIA:	None
TOTAL DEPTH:	11 feet	CASING DEPTH:	None

	VISUAL MANUAL SOIL CLASSIFICATION	OVA RESULTS
DEPTH	COLOR, TEXTURE, STRUCTURE, CONSISTENCY, ODOR, ETC.	PERCENT RECOVERY
(ft.)		BLOW COUNTS

	Depths correspond to changes in soil type	
0-2'	clayey sandy silt (ML); brown, moist, no odor	PID=8-2(0-2): 1 PPM
2-4'	clayey sandy silt (ML); brown, moist, no odor	PID=8-2(2-4): 2 PPM
4-6'	clayey sandy silt (ML); brown, very moist, no odor	PID=8-2(4-6): 3 PPM
6-8'	clayey sandy silt (ML); brown, very moist, no odor	PID=8-2(6-8): 2 PPM
8-11'	clayey sandy silt (ML) & silty sand (SM); brown, rocks, saturated, no odor	PID=8-2(8-11): 2 PPM
	refusal at 11', groundwater encountered at 3.7 feet	

RISER LENGTH (ft)	DEPTH (ft)		DIAMETER (in)		MATERIAL	
SCREEN LENGTH (ft)	DEPTH (ft)		DIAMETER (in)		MATERIAL	
DEPTH TO TOP OF SAND			BAGS OF SAND	<u> </u>		
DEPTH TO TOP SEAL		BENTONIT	E USED		BAGS OF CEME	NT USED

# FIELD DRILLING RECORD

PROJECT NAME: PROJECT NUMBER:	NC DOT I-3318BB Parcel 008 Kenneth M. Taylor 2014-093	BORING/WELL NO:	8-3
SITE LOCATION:	675 Cummins Drive, Kenly, Johnston County, NC	BORING/WELL LOCATION:	Center parcel near road
START DATE:	6/2/14	COMPLETED:	6/2/14
GEOLOGIST:	T. Leatherman	DRILLER:	Solutions, IES
DRILL METHOD:	Geoprobe	SAMPLE METHOD:	Macro-core
BORING DIA:	2-inch	CASING DIA:	None
TOTAL DEPTH:	8 feet	CASING DEPTH:	None

	VISUAL MANUAL SOIL CLASSIFICATION	OVA RESULTS
DEPTH	COLOR, TEXTURE, STRUCTURE, CONSISTENCY, ODOR, ETC.	PERCENT RECOVERY
(ft.)		BLOW COUNTS

		1
	Depths correspond to changes in soil type	
0-2'	clayey sandy silt (ML); brown to gray, moist, no odor	PID=8-3(0-2): 1 PPM
2-4'	sandy silty clay (CL); brown to tan, moist, no odor	PID=8-3(2-4): 2.5 PPM
4-6'	sandy silty clay (CL); brown to gray, saturated, no odor	PID=8-3(4-6): 2 PPM
6-8'	clayey sand (SC); brown to gray, saturated, no odor	PID=8-3(6-8): 2 PPM
	refusal at 8', groundwater encountered at 7.6 feet	

RISER LENGTH (ft)	DEPTH (ft)	DIAMETER (in)	MATERIAL
SCREEN LENGTH (ft)	DEPTH (ft)	DIAMETER (in)	MATERIAL
DEPTH TO TOP OF SAND		BAGS OF SAND	
DEPTH TO TOP SEAL	BEN	TONITE USED	BAGS OF CEMENT USED

# FIELD DRILLING RECORD

PROJECT NAME: PROJECT NUMBER:	NC DOT I-3318BB Parcel 008 Kenneth M. Taylor 2014-093	BORING/WELL NO:	8-4
SITE LOCATION:	675 Cummins Drive, Kenly, Johnston County, NC	BORING/WELL LOCATION:	East parcel near road
START DATE:	6/2/14	COMPLETED:	6/2/14
GEOLOGIST:	T. Leatherman	DRILLER:	Solutions, IES
DRILL METHOD:	Geoprobe	SAMPLE METHOD:	Macro-core
BORING DIA:	2-inch	CASING DIA:	1-inch
TOTAL DEPTH:	12 feet	CASING DEPTH:	1-inch

	VISUAL MANUAL SOIL CLASSIFICATION	OVA RESULTS
DEPTH	COLOR, TEXTURE, STRUCTURE, CONSISTENCY, ODOR, ETC.	PERCENT RECOVERY
(ft.)		BLOW COUNTS

	Depths correspond to changes in soil type	
0-2'	clayey sandy silt (ML); brown, moist, no odor	PID=8-4(0-2): 2 PPM
2-4'	sandy silty clay (CL); brown to light gray, moist, no odor	PID=8-4(2-4): 3 PPM
4-6'	sandy silty clay (CL); tan to It gray, very moist, no odor	PID=8-4(4-6): 2 PPM
6-8'	sandy silty clay (CL); tan to It gray, very moist, no odor	PID=8-4(6-8): 2 PPM
8-10'	sandy silty clay (CL); brown to lt gray, saturated, no odor	PID=8-4(8-10): 13 PPM
10-12'	clayey silty sand (SM); brown, rock frags, saturated, no odor	PID=8-4(10-12): 17 PPM
	refusal at 12', Temp well set to 12 feet, groundwater at 3.35 feet	

RISER LENGTH (ft) 2	DEPTH (ft) 0-2	DIAMETER (in) 1	MATERIAL <u>PVC</u> .
SCREEN LENGTH (ft) 10	DEPTH (ft) 2-12	DIAMETER (in) 1	MATERIAL <u>PVC</u> .
DEPTH TO TOP OF SAND		BAGS OF SAND	
DEPTH TO TOP SEAL	BENTON	ITE USED	BAGS OF CEMENT USED

# APPENDIX D

Q	ED											ſ	<u>QROS</u>
				Hydroca	arbon An	alysis R	esults						
Client: Address:	NCDOT - Johnston County I-3318 Off Old River Road in Kenly, NC; I	BB Parcel 8							Sa Sample Sampl	mples f es extra es ana	aken acted lysed		8-1 thru 8-4
Contact: Proiect:	NCDOT - Johnston County I-3318	BB. Parce	<del>3</del>   8							Оре	erator		Ryan Kramer
110,000.		<b>DD</b> , 1 010	510										
Matrix	Sample ID	Dilution used	BTEX (C6 - C9)	GRO (C5 - C10)	DRO (C10 - C35)	TPH (C5 - C35)	Total Aromatics (C10-C35)	16 EPA PAHs	BaP		Ratios		HC Fingerprint Match
										% light	% mid	% heavy	
S	8-1 (4-6)	13.0	<0.7	<0.7	0.53	0.53	0.41	0.05	<0.013	47.3	35.7	17	Road Tar 91.2%
S	8-2 (4-6)	12.0	<0.6	<0.6	0.17	0.17	<0.12	<0.01	<0.012	0	7.7	92.3	Deg.Fuel Residue (FCM) (P) 9.4%
S	8-3 (2-4)	12.0	<0.6	<0.6	0.49	0.49	0.3	0.08	<0.012	69.8	4.4	25.7	Waste Oil (FCM) (P) 14.9%
S	8-4 (2-4)	13.0	<0.6	<0.6	0.33	0.33	<0.13	<0.01	<0.013	0	10.3	89.7	Deg.Fuel Residue (FCM) (P) 14.4%
	Initial Ca	alibrator	QC check	OK					Final FC	CM QC (	Check	OK	90.4%
Results gen Fingerprints (SBS) or (LE	erated by a QED HC-1 analyser. Concent provide a tentative hydrocarbon identification 3S) = Site Specific or Library Background St	tration value on. The abbi ubtraction a	s in mg/kg f eviations ar oplied to res	or soil sample e:- FCM = R ult : (PFM) =	es and mg/L f esults calcula Poor Fingerpi	or water samp ted using Fun rint Match : (T	bles. Soil valu damental Calib ) = Turbid : (P)	ues are not pration Mod = Particula	corrected fo le : % = cont ate present	or moistur fidence fo	e or sto or sampl	ne conte e fingerj	ent print match to library

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# APPENDIX E



Pace Analytical Services, Inc. 9800 Kincey Ave. Suite 100 Huntersville, NC 28078 (704)875-9092

June 12, 2014

Chemical Testing Engineer NCDOT Materials & Tests Unit 1801 Blue Ridge Road Raleigh, NC 27607

RE: Project: 2014-093 Johnston 34182.1.2 Pace Project No.: 92204085

Dear Chemical Engineer:

Enclosed are the analytical results for sample(s) received by the laboratory on June 04, 2014. The results relate only to the samples included in this report. Results reported herein conform to the most current TNI standards and the laboratory's Quality Assurance Manual, where applicable, unless otherwise noted in the body of the report.

Analyses were performed at the Pace Analytical Services location indicated on the sample analyte page for analysis unless otherwise footnoted.

If you have any questions concerning this report, please feel free to contact me.

Sincerely,

for Brudley

Jon D Bradley jon.bradley@pacelabs.com Project Manager

Enclosures

cc: Tim Leatherman, Pyramid



# **REPORT OF LABORATORY ANALYSIS**

This report shall not be reproduced, except in full, without the written consent of Pace Analytical Services, Inc..



Pace Analytical Services, Inc. 9800 Kincey Ave. Suite 100 Huntersville, NC 28078 (704)875-9092

## CERTIFICATIONS

Project: 2014-093 Johnston 34182.1.2

Pace Project No.: 92204085

#### **Charlotte Certification IDs**

9800 Kincey Ave. Ste 100, Huntersville, NC 28078 North Carolina Drinking Water Certification #: 37706 North Carolina Field Services Certification #: 5342 North Carolina Wastewater Certification #: 12 South Carolina Certification #: 99006001

Florida/NELAP Certification #: E87627 Kentucky UST Certification #: 84 West Virginia Certification #: 357 Virginia/VELAP Certification #: 460221



# SAMPLE ANALYTE COUNT

 Project:
 2014-093 Johnston 34182.1.2

 Pace Project No.:
 92204085

Lab ID	Sample ID	Method	Analysts	Analytes Reported	Laboratory
92204085001	8-2 (4-6)	EPA 8270	RES	74	PASI-C
		EPA 8260	DLK	70	PASI-C
		ASTM D2974-87	ZAK	1	PASI-C
92204085002	8-4 (TW)	EPA 625	BPJ	58	PASI-C
		SM 6200B	CAH	63	PASI-C



Project: 2014-093 Johnston 34182.1.2

Pace Project No.: 92204085

## Method: EPA 625

Description:625 MSSVClient:NCDOT East CentralDate:June 12, 2014

### General Information:

1 sample was analyzed for EPA 625. All samples were received in acceptable condition with any exceptions noted below.

#### Hold Time:

The samples were analyzed within the method required hold times with any exceptions noted below.

### Sample Preparation:

The samples were prepared in accordance with EPA 625 with any exceptions noted below.

### Initial Calibrations (including MS Tune as applicable):

All criteria were within method requirements with any exceptions noted below.

### Continuing Calibration:

All criteria were within method requirements with any exceptions noted below.

## Internal Standards:

All internal standards were within QC limits with any exceptions noted below.

### Surrogates:

All surrogates were within QC limits with any exceptions noted below.

## Method Blank:

All analytes were below the report limit in the method blank, where applicable, with any exceptions noted below.

### Laboratory Control Spike:

All laboratory control spike compounds were within QC limits with any exceptions noted below.

### Matrix Spikes:

All percent recoveries and relative percent differences (RPDs) were within acceptance criteria with any exceptions noted below.

### QC Batch: OEXT/28189

A matrix spike and/or matrix spike duplicate (MS/MSD) were performed on the following sample(s): 92204282003

- R1: RPD value was outside control limits.
- MSD (Lab ID: 1218107)
  - 2,4-Dinitrophenol

## Additional Comments:



Project: 2014-093 Johnston 34182.1.2

Pace Project No.: 92204085

Method:EPA 8270Description:8270 MSSV MicrowaveClient:NCDOT East CentralDate:June 12, 2014

### General Information:

1 sample was analyzed for EPA 8270. All samples were received in acceptable condition with any exceptions noted below.

#### Hold Time:

The samples were analyzed within the method required hold times with any exceptions noted below.

### Sample Preparation:

The samples were prepared in accordance with EPA 3546 with any exceptions noted below.

#### Initial Calibrations (including MS Tune as applicable):

All criteria were within method requirements with any exceptions noted below.

#### Continuing Calibration:

All criteria were within method requirements with any exceptions noted below.

### Internal Standards:

All internal standards were within QC limits with any exceptions noted below.

#### Surrogates:

All surrogates were within QC limits with any exceptions noted below.

#### Method Blank:

All analytes were below the report limit in the method blank, where applicable, with any exceptions noted below.

#### Laboratory Control Spike:

All laboratory control spike compounds were within QC limits with any exceptions noted below.

#### Matrix Spikes:

All percent recoveries and relative percent differences (RPDs) were within acceptance criteria with any exceptions noted below.

#### Additional Comments:



Project: 2014-093 Johnston 34182.1.2

Pace Project No.: 92204085

### Method: SM 6200B

Description:6200B MSVClient:NCDOT East CentralDate:June 12, 2014

## General Information:

1 sample was analyzed for SM 6200B. All samples were received in acceptable condition with any exceptions noted below.

#### Hold Time:

The samples were analyzed within the method required hold times with any exceptions noted below.

## Initial Calibrations (including MS Tune as applicable):

All criteria were within method requirements with any exceptions noted below.

#### Continuing Calibration:

All criteria were within method requirements with any exceptions noted below.

#### Internal Standards:

All internal standards were within QC limits with any exceptions noted below.

### Surrogates:

All surrogates were within QC limits with any exceptions noted below.

### Method Blank:

All analytes were below the report limit in the method blank, where applicable, with any exceptions noted below.

### Laboratory Control Spike:

All laboratory control spike compounds were within QC limits with any exceptions noted below.

## Matrix Spikes:

All percent recoveries and relative percent differences (RPDs) were within acceptance criteria with any exceptions noted below.

### QC Batch: MSV/27102

A matrix spike and/or matrix spike duplicate (MS/MSD) were performed on the following sample(s): 92204081001

M0: Matrix spike recovery and/or matrix spike duplicate recovery was outside laboratory control limits.

- MSD (Lab ID: 1215547)
  - Vinyl chloride

### Additional Comments:



Project: 2014-093 Johnston 34182.1.2

Pace Project No.: 92204085

#### Method: EPA 8260

Description:8260/5035A Volatile OrganicsClient:NCDOT East CentralDate:June 12, 2014

## General Information:

1 sample was analyzed for EPA 8260. All samples were received in acceptable condition with any exceptions noted below.

#### Hold Time:

The samples were analyzed within the method required hold times with any exceptions noted below.

**Initial Calibrations (including MS Tune as applicable):** All criteria were within method requirements with any exceptions noted below.

#### Continuing Calibration:

All criteria were within method requirements with any exceptions noted below.

#### **Internal Standards:**

All internal standards were within QC limits with any exceptions noted below.

#### Surrogates:

All surrogates were within QC limits with any exceptions noted below.

#### Method Blank:

All analytes were below the report limit in the method blank, where applicable, with any exceptions noted below.

#### Laboratory Control Spike:

All laboratory control spike compounds were within QC limits with any exceptions noted below.

## Matrix Spikes:

All percent recoveries and relative percent differences (RPDs) were within acceptance criteria with any exceptions noted below.

### **Duplicate Sample:**

All duplicate sample results were within method acceptance criteria with any exceptions noted below.

### Additional Comments:

This data package has been reviewed for quality and completeness and is approved for release.



Project: 2014-093 Johnston 34182.1.2

Pace Project No.: 92204085

Sample: 8-2 (4-6)	Lab ID: 922	04085001	Collected: 06/02/	14 16:0	0 Received: 06	04/14 17:30 N	latrix: Solid	
Results reported on a "dry-weigl	ht" basis							
Parameters	Results	Units	Report Limit	DF	Prepared	Analyzed	CAS No.	Qual
8270 MSSV Microwave	Analytical Meth	nod: EPA 827	0 Preparation Met	hod: EF	PA 3546			
Acenaphthene	ND ug	/kg	422	1	06/05/14 13:07	06/10/14 12:26	83-32-9	
Acenaphthylene	ND ug	/ka	422	1	06/05/14 13:07	06/10/14 12:26	208-96-8	
Aniline	ND ug	/kg	422	1	06/05/14 13:07	06/10/14 12:26	62-53-3	
Anthracene	ND ug	/kg	422	1	06/05/14 13:07	06/10/14 12:26	120-12-7	
Benzo(a)anthracene	ND ug	/kg	422	1	06/05/14 13:07	06/10/14 12:26	56-55-3	
Benzo(a)pyrene	ND ug	/ka	422	1	06/05/14 13:07	06/10/14 12:26	50-32-8	
Benzo(b)fluoranthene	ND ug	/ka	422	1	06/05/14 13:07	06/10/14 12:26	205-99-2	
Benzo(a,h,i)pervlene	ND ug	/ka	422	1	06/05/14 13:07	06/10/14 12:26	191-24-2	
Benzo(k)fluoranthene	ND ug	/ka	422	1	06/05/14 13:07	06/10/14 12:26	207-08-9	
Benzoic Acid	ND ug	/ka	2110	1	06/05/14 13:07	06/10/14 12:26	65-85-0	
Benzvl alcohol	ND ug	/ka	844	1	06/05/14 13:07	06/10/14 12:26	100-51-6	
4-Bromophenylphenyl ether	ND ug	/ka	422	1	06/05/14 13:07	06/10/14 12:26	101-55-3	
Butylbenzylphthalate	ND ug	/ka	422	1	06/05/14 13:07	06/10/14 12:26	85-68-7	
4-Chloro-3-methylphenol	ND ug	/ka	844	1	06/05/14 13:07	06/10/14 12:26	59-50-7	
4-Chloroaniline	ND ug	/ka	2110	1	06/05/14 13:07	06/10/14 12:26	106-47-8	
bis(2-Chloroethoxy)methane	ND ug	/ka	422	1	06/05/14 13:07	06/10/14 12:26	111-91-1	
bis(2-Chloroethyl) ether	ND ug	/ka	422	1	06/05/14 13:07	06/10/14 12:26	111-44-4	
bis(2-Chloroisopropyl) ether	ND ug	/ka	422	1	06/05/14 13:07	06/10/14 12:26	108-60-1	
2-Chloronaphthalene	ND ug	/ka	422	1	06/05/14 13:07	06/10/14 12:26	91-58-7	
2-Chlorophenol	ND ug	/kg	422	1	06/05/14 13:07	06/10/14 12:26	95-57-8	
4-Chlorophenylphenyl ether	ND ug	/ka	422	1	06/05/14 13:07	06/10/14 12:26	7005-72-3	
Chrysene	ND ug	/ka	422	1	06/05/14 13:07	06/10/14 12:26	218-01-9	
Dibenz(a h)anthracene	ND ug	/ka	422	1	06/05/14 13:07	06/10/14 12:26	53-70-3	
Dibenzofuran	ND ug	/kg	422	1	06/05/14 13:07	06/10/14 12:26	132-64-9	
1.2-Dichlorobenzene	ND ug	/kg	422	1	06/05/14 13:07	06/10/14 12:26	95-50-1	
1.3-Dichlorobenzene	ND ug	/ka	422	1	06/05/14 13:07	06/10/14 12:26	541-73-1	
1 4-Dichlorobenzene	ND ug	/ka	422	1	06/05/14 13:07	06/10/14 12:26	106-46-7	
3 3'-Dichlorobenzidine	ND ug	/kg	2110	1	06/05/14 13:07	06/10/14 12:26	91-94-1	
2 4-Dichlorophenol	ND ug	/kg	422	1	06/05/14 13:07	06/10/14 12:26	120-83-2	
Diethylphthalate	ND ug	/kg	422	1	06/05/14 13:07	06/10/14 12:26	84-66-2	
2 4-Dimethylphenol	ND ug	/ka	422	1	06/05/14 13:07	06/10/14 12:26	105-67-9	
Dimethylphthalate	ND ug	/kg	422	1	06/05/14 13:07	06/10/14 12:26	131-11-3	
Di-n-butylphthalate	ND ug	/kg	422	1	06/05/14 13:07	06/10/14 12:26	84-74-2	
4 6-Dinitro-2-methylphenol	ND ug	/kg	844	1	06/05/14 13:07	06/10/14 12:26	534-52-1	
2 4-Dinitrophenol	ND ug	/kg	2110	1	06/05/14 13:07	06/10/14 12:26	51-28-5	
2 4-Dinitrotoluene	ND ug	/kg	422	1	06/05/14 13:07	06/10/14 12:26	121-14-2	
2 6-Dinitrotoluene	ND ug	/ka	422	1	06/05/14 13:07	06/10/14 12:26	606-20-2	
Di-n-octylphthalate	ND ug	/kg	422	1	06/05/14 13:07	06/10/14 12:26	117-84-0	
bis(2-Ethylbeyyl)phthalate	ND ug	/kg	422	1	06/05/14 13:07	06/10/14 12:26	117-81-7	
Fluoranthene	ND ug	/kg	422	1	06/05/14 13:07	06/10/14 12:26	206-44-0	
Fluorene	ND ug	/kg	422	1	06/05/14 13:07	06/10/14 12:26	86-73-7	
Hexachloro-1 3-butadiene		/ka	422	1	06/05/14 13:07	06/10/14 12:20	87-68-3	
Hexachlorobenzene		/ka		1	06/05/14 13:07	06/10/14 12:20	118-74-1	
Hexachlorocyclopentadiene		/ka		1	06/05/14 13:07	06/10/14 12:20	77-47-4	
Hexachloroethane		/ka	422	1	06/05/14 13:07	06/10/14 12:20	67-72-1	
Indeno(1.2.3-cd)pyrene	ND ug	/ka	422	1	06/05/14 13:07	06/10/14 12:20	193-39-5	
	ND ug	/ NG	422		00/00/14 10.07	00/10/14 12.20	100-00-0	

## **REPORT OF LABORATORY ANALYSIS**

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Project: 2014-093 Johnston 34182.1.2

Pace Project No.: 92204085

Sample: 8-2 (4-6)	Lab ID: 9220	04085001	Collected: 06/02/	4 16:00	0 Received: 06	/04/14 17:30 N	latrix: Solid	
Results reported on a "dry-weight	t" basis							
Parameters	Results	Units	Report Limit	DF	Prepared	Analyzed	CAS No.	Qual
8270 MSSV Microwave	Analytical Meth	od: EPA 82	70 Preparation Met	hod: EF	PA 3546			
Isophorone	ND ug/	/kg	422	1	06/05/14 13:07	06/10/14 12:26	78-59-1	
1-Methylnaphthalene	ND ug	/kg	422	1	06/05/14 13:07	06/10/14 12:26	90-12-0	
2-Methylnaphthalene	ND ug	/kg	422	1	06/05/14 13:07	06/10/14 12:26	91-57-6	
2-Methylphenol(o-Cresol)	ND ug	/kg	422	1	06/05/14 13:07	06/10/14 12:26	95-48-7	
3&4-Methylphenol(m&p Cresol)	ND ug	/kg	422	1	06/05/14 13:07	06/10/14 12:26		
Naphthalene	ND ug	/kg	422	1	06/05/14 13:07	06/10/14 12:26	91-20-3	
2-Nitroaniline	ND ug	/kg	2110	1	06/05/14 13:07	06/10/14 12:26	88-74-4	
3-Nitroaniline	ND ug	/kg	2110	1	06/05/14 13:07	06/10/14 12:26	99-09-2	
4-Nitroaniline	ND ug	/kg	844	1	06/05/14 13:07	06/10/14 12:26	100-01-6	
Nitrobenzene	ND ug	/kg	422	1	06/05/14 13:07	06/10/14 12:26	98-95-3	
2-Nitrophenol	ND ug	/kg	422	1	06/05/14 13:07	06/10/14 12:26	88-75-5	
4-Nitrophenol	ND ug	/ka	2110	1	06/05/14 13:07	06/10/14 12:26	100-02-7	
N-Nitrosodimethylamine	ND ug	/ka	422	1	06/05/14 13:07	06/10/14 12:26	62-75-9	
N-Nitroso-di-n-propylamine	ND ug	/ka	422	1	06/05/14 13:07	06/10/14 12:26	621-64-7	
N-Nitrosodiphenvlamine	ND ug	/ka	422	1	06/05/14 13:07	06/10/14 12:26	86-30-6	
Pentachlorophenol	ND ug	/ka	2110	1	06/05/14 13:07	06/10/14 12:26	87-86-5	
Phenanthrene	ND ug	/ka	422	1	06/05/14 13:07	06/10/14 12:26	85-01-8	
Phenol	ND ug	/ka	422	1	06/05/14 13:07	06/10/14 12:26	108-95-2	
Pyrene	ND ug	/ka	422	1	06/05/14 13:07	06/10/14 12:26	129-00-0	
1.2.4-Trichlorobenzene	ND ug	/ka	422	1	06/05/14 13:07	06/10/14 12:26	120-82-1	
2.4.5-Trichlorophenol	ND ug	/ka	422	1	06/05/14 13:07	06/10/14 12:26	95-95-4	
2.4.6-Trichlorophenol	ND ug	/ka	422	1	06/05/14 13:07	06/10/14 12:26	88-06-2	
Surrogates						00,10,1112120	00 00 1	
Nitrobenzene-d5 (S)	43 %		23-110	1	06/05/14 13:07	06/10/14 12:26	4165-60-0	
2-Fluorobiphenyl (S)	47 %		30-110	1	06/05/14 13:07	06/10/14 12:26	321-60-8	
Terphenyl-d14 (S)	79 %		28-110	1	06/05/14 13:07	06/10/14 12:26	1718-51-0	
Phenol-d6 (S)	44 %		22-110	1	06/05/14 13:07	06/10/14 12:26	13127-88-3	
2-Fluorophenol (S)	45 %		13-110	1	06/05/14 13:07	06/10/14 12:26	367-12-4	
2,4,6-Tribromophenol (S)	59 %		27-110	1	06/05/14 13:07	06/10/14 12:26	118-79-6	
8260/5035A Volatile Organics	Analytical Meth	od: EPA 820	60					
Anatana	ND	//	440			00/00/4447-50	07.04.4	
Acetone	ND ug/	/Kg //	110	1		06/09/14 17:52	67-64-1	
Benzene	ND ug/	кg	5.5	1		06/09/14 17:52	71-43-2	
Bromobenzene	ND ug/	′кg //	5.5	1		06/09/14 17:52	108-86-1	
Bromocniorometnane	ND ug/	′кg //	5.5	1		06/09/14 17:52	74-97-5	
Bromodicniorometnane	ND ug/	′кg	5.5	1		06/09/14 17:52	75-27-4	
Bromotorm	ND ug/	′kg	5.5	1		06/09/14 17:52	75-25-2	
Bromomethane	ND ug/	′kg	11.0	1		06/09/14 17:52	74-83-9	
2-Butanone (MEK)	ND ug/	′kg	110	1		06/09/14 17:52	78-93-3	
n-Butylbenzene	ND ug/	′kg	5.5	1		06/09/14 17:52	104-51-8	
sec-Butylbenzene	ND ug/	'kg	5.5	1		06/09/14 17:52	135-98-8	
tert-Butylbenzene	ND ug/	'kg	5.5	1		06/09/14 17:52	98-06-6	
Carbon tetrachloride	ND ug/	'kg	5.5	1		06/09/14 17:52	56-23-5	
Chlorobenzene	ND ug/	′kg	5.5	1		06/09/14 17:52	108-90-7	
Chloroethane	ND ug/	′kg	11.0	1		06/09/14 17:52	75-00-3	
Chloroform	ND ug/	′kg	5.5	1		06/09/14 17:52	67-66-3	

# **REPORT OF LABORATORY ANALYSIS**

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Project: 2014-093 Johnston 34182.1.2

Pace Project No.: 92204085

Sample: 8-2 (4-6)	Lab ID: 9220	04085001 Collect	cted: 06/02/	14 16:00	Received: 06	5/04/14 17:30 N	latrix: Solid	
Results reported on a "dry-weigh	t" basis							
Parameters	Results	Units R	Report Limit	DF	Prepared	Analyzed	CAS No.	Qual
8260/5035A Volatile Organics	Analytical Meth	nod: EPA 8260						
Chloromethane	ND ug	/kg	11.0	1		06/09/14 17:52	74-87-3	
2-Chlorotoluene	ND ug	/kg	5.5	1		06/09/14 17:52	95-49-8	
4-Chlorotoluene	ND ug	/kg	5.5	1		06/09/14 17:52	106-43-4	
1,2-Dibromo-3-chloropropane	ND ug	/kg	5.5	1		06/09/14 17:52	96-12-8	
Dibromochloromethane	ND ug	/kg	5.5	1		06/09/14 17:52	124-48-1	
1,2-Dibromoethane (EDB)	ND ug	/kg	5.5	1		06/09/14 17:52	106-93-4	
Dibromomethane	ND ug	/kg	5.5	1		06/09/14 17:52	74-95-3	
1,2-Dichlorobenzene	ND ug	/kg	5.5	1		06/09/14 17:52	95-50-1	
1,3-Dichlorobenzene	ND ug	/kg	5.5	1		06/09/14 17:52	541-73-1	
1,4-Dichlorobenzene	ND ug	/kg	5.5	1		06/09/14 17:52	106-46-7	
Dichlorodifluoromethane	ND ug	/kg	11.0	1		06/09/14 17:52	75-71-8	
1,1-Dichloroethane	ND ug	/kg	5.5	1		06/09/14 17:52	75-34-3	
1,2-Dichloroethane	ND ug	/kg	5.5	1		06/09/14 17:52	107-06-2	
1,1-Dichloroethene	ND ug	/kg	5.5	1		06/09/14 17:52	75-35-4	
cis-1,2-Dichloroethene	ND ug	/kg	5.5	1		06/09/14 17:52	156-59-2	
trans-1,2-Dichloroethene	ND ug	/kg	5.5	1		06/09/14 17:52	156-60-5	
1,2-Dichloropropane	ND ug	/kg	5.5	1		06/09/14 17:52	78-87-5	
1,3-Dichloropropane	ND ug	/kg	5.5	1		06/09/14 17:52	142-28-9	
2,2-Dichloropropane	ND ug	/kg	5.5	1		06/09/14 17:52	594-20-7	
1,1-Dichloropropene	ND ug	/kg	5.5	1		06/09/14 17:52	563-58-6	
cis-1,3-Dichloropropene	ND ug	/kg	5.5	1		06/09/14 17:52	10061-01-5	
trans-1,3-Dichloropropene	ND ug	/kg	5.5	1		06/09/14 17:52	10061-02-6	
Diisopropyl ether	ND ug	/kg	5.5	1		06/09/14 17:52	108-20-3	
Ethylbenzene	ND ug	/kg	5.5	1		06/09/14 17:52	100-41-4	
Hexachloro-1,3-butadiene	ND ug	/kg	5.5	1		06/09/14 17:52	87-68-3	
2-Hexanone	ND ug	/kg	55.0	1		06/09/14 17:52	591-78-6	
Isopropylbenzene (Cumene)	ND ug	/kg	5.5	1		06/09/14 17:52	98-82-8	
p-Isopropyltoluene	ND ug	/kg	5.5	1		06/09/14 17:52	99-87-6	
Methylene Chloride	ND ug	/kg	22.0	1		06/09/14 17:52	75-09-2	
4-Methyl-2-pentanone (MIBK)	ND ug	/kg	55.0	1		06/09/14 17:52	108-10-1	
Methyl-tert-butyl ether	ND ug	/kg	5.5	1		06/09/14 17:52	1634-04-4	
Naphthalene	ND ug	/kg	5.5	1		06/09/14 17:52	91-20-3	
n-Propylbenzene	ND ug	/kg	5.5	1		06/09/14 17:52	103-65-1	
Styrene	ND ug	/kg	5.5	1		06/09/14 17:52	100-42-5	
1,1,1,2-Tetrachloroethane	ND ug	/kg	5.5	1		06/09/14 17:52	630-20-6	
1,1,2,2-Tetrachloroethane	ND ug	/kg	5.5	1		06/09/14 17:52	79-34-5	
Tetrachloroethene	ND ug	/kg	5.5	1		06/09/14 17:52	127-18-4	
Toluene	ND ug	/kg	5.5	1		06/09/14 17:52	108-88-3	
1,2,3-Trichlorobenzene	ND ug	/kg	5.5	1		06/09/14 17:52	87-61-6	
1,2,4-Trichlorobenzene	ND ug	/kg	5.5	1		06/09/14 17:52	120-82-1	
1,1,1-Trichloroethane	ND ug	/kg	5.5	1		06/09/14 17:52	71-55-6	
1,1,2-Trichloroethane	ND ug	/kg	5.5	1		06/09/14 17:52	79-00-5	
Trichloroethene	ND ug	/kg	5.5	1		06/09/14 17:52	79-01-6	
Trichlorofluoromethane	ND ug	/kg	5.5	1		06/09/14 17:52	75-69-4	
1,2,3-Trichloropropane	ND ug	/kg	5.5	1		06/09/14 17:52	96-18-4	
1,2,4-Trimethylbenzene	ND ug	/kg	5.5	1		06/09/14 17:52	95-63-6	



Project: 2014-093 Johnston 34182.1.2

Pace Project No.: 92204085

Sample: 8-2 (4-6)	Lab ID: 9220	04085001	Collected:	06/02/1	4 16:00	Received: 0	6/04/14 17:30 N	latrix: Solid				
Results reported on a "dry-weight	Results reported on a "dry-weight" basis											
Parameters	Results	Units	Report	Limit	DF	Prepared	Analyzed	CAS No.	Qual			
8260/5035A Volatile Organics	Analytical Meth	od: EPA 82	260									
1,3,5-Trimethylbenzene	ND ug/	/kg		5.5	1		06/09/14 17:52	108-67-8				
Vinyl acetate	ND ug/	/kg		55.0	1		06/09/14 17:52	108-05-4				
Vinyl chloride	ND ug/	/kg		11.0	1		06/09/14 17:52	75-01-4				
Xylene (Total)	ND ug/	/kg		11.0	1		06/09/14 17:52	1330-20-7				
m&p-Xylene	ND ug/	/kg		11.0	1		06/09/14 17:52	179601-23-1				
o-Xylene	ND ug/	/kg		5.5	1		06/09/14 17:52	95-47-6				
Surrogates												
Toluene-d8 (S)	101 %		7	0-130	1		06/09/14 17:52	2037-26-5				
4-Bromofluorobenzene (S)	102 %		7	0-130	1		06/09/14 17:52	460-00-4				
1,2-Dichloroethane-d4 (S)	113 %		7	0-132	1		06/09/14 17:52	17060-07-0				
Percent Moisture	Analytical Meth	od: ASTM	D2974-87									
Percent Moisture	<b>21.8</b> %			0.10	1		06/12/14 16:10					



Project: 2014-093 Johnston 34182.1.2

Pace Project No.: 92204085

Sample: 8-4 (TW)	Lab ID: 92	204085002	Collected:	06/03/1	4 08:05	Received: 06	i/04/14 17:30 N	latrix: Water	
Parameters	Results	Units	Repor	t Limit	DF	Prepared	Analyzed	CAS No.	Qual
625 MSSV	Analytical Me	thod: EPA 62	25 Preparati	on Metho	od: EPA	625			
Acenaphthene	ND u	g/L		5.6	1	06/10/14 16:53	06/11/14 22:27	83-32-9	
Acenaphthylene	ND u	g/L		5.6	1	06/10/14 16:53	06/11/14 22:27	208-96-8	
Anthracene	ND u	g/L		5.6	1	06/10/14 16:53	06/11/14 22:27	120-12-7	
Benzo(a)anthracene	ND u	g/L		5.6	1	06/10/14 16:53	06/11/14 22:27	56-55-3	
Benzo(a)pyrene	ND u	g/L		5.6	1	06/10/14 16:53	06/11/14 22:27	50-32-8	
Benzo(b)fluoranthene	ND u	g/L		5.6	1	06/10/14 16:53	06/11/14 22:27	205-99-2	
Benzo(g,h,i)perylene	ND u	g/L		5.6	1	06/10/14 16:53	06/11/14 22:27	191-24-2	
Benzo(k)fluoranthene	ND u	g/L		5.6	1	06/10/14 16:53	06/11/14 22:27	207-08-9	
4-Bromophenylphenyl ether	ND u	g/L		5.6	1	06/10/14 16:53	06/11/14 22:27	101-55-3	
Butylbenzylphthalate	ND u	g/L		5.6	1	06/10/14 16:53	06/11/14 22:27	85-68-7	
4-Chloro-3-methylphenol	ND u	g/L		5.6	1	06/10/14 16:53	06/11/14 22:27	59-50-7	
bis(2-Chloroethoxy)methane	ND u	g/L		11.1	1	06/10/14 16:53	06/11/14 22:27	111-91-1	
bis(2-Chloroethyl) ether	ND u	a/L		5.6	1	06/10/14 16:53	06/11/14 22:27	111-44-4	
bis(2-Chloroisopropyl) ether	ND u	a/L		5.6	1	06/10/14 16:53	06/11/14 22:27	108-60-1	
2-Chloronaphthalene	ND u	9' – a/l		5.6	1	06/10/14 16:53	06/11/14 22:27	91-58-7	
2-Chlorophenol	ND u	g/ = a/l		5.6	1	06/10/14 16:53	06/11/14 22:27	95-57-8	
4-Chlorophenylphenyl ether	ND u	g/L a/l		5.6	1	06/10/14 16:53	06/11/14 22:27	7005-72-3	
Chrysene		g/⊑ a/l		5.6	1	06/10/14 16:53	06/11/14 22:27	218-01-9	
Dibenz(a b)anthracene		g/L a/l		5.6	1	06/10/14 16:53	06/11/14 22:27	53-70-3	
3 3'-Dichlorobenzidine		g/∟ α/I		27.8	1	06/10/14 16:53	06/11/14 22:27	91-94-1	
2 4-Dichlorophenol		g/∟ α/I		5.6	1	06/10/14 16:53	06/11/14 22:27	120-83-2	
Diethylohthalate	ND u	g/L g/l		5.6	1	06/10/14 16:53	06/11/14 22:27	84-66-2	
2 4-Dimethylphenol		g/L g/l		11 1	1	06/10/14 16:53	06/11/14 22:27	105-67-0	
Dimethylohthalate		g/∟ α/I		56	1	06/10/14 16:53	06/11/14 22:27	131-11-3	
		g/∟ g/l		5.0	1	06/10/14 16:53	06/11/14 22:27	94 74 2	
4 6 Dipitro 2 mothylphonol		g/∟ α/I		2.0	1	06/10/14 10:53	06/11/14 22.27	04-74-2 524 52 4	
2.4 Dinitrophonol		g/∟ α/I		22.2 55.6	1	06/10/14 16:53	06/11/14 22.27	51 29 5	
2,4-Dinitrophenoi		g/∟ α/I		55.0	1	06/10/14 16:53	06/11/14 22.27	01-20-0 101 14 0	
2,4-Dinitrotoluene		g/∟ α/I		5.0	1	06/10/14 16:53	06/11/14 22.27	121-14-2	
2,6-Dimitrotoluene		g/∟ ~/!		5.0 5.0	1	06/10/14 16:53	06/11/14 22:27	000-20-2	
bis (2. Ethyllh surd) ab th a late	ND u	g/L		0.C	1	06/10/14 16:53	06/11/14 22:27	117-64-0	
		g/L ~/l		5.0 5.0	1	06/10/14 16:53	06/11/14 22:27	117-01-7	
Fluorance		g/∟ ≈/l		5.0	1	06/10/14 10:55	06/11/14 22.27	200-44-0	
Fluorene		g/∟ ~/!		5.0 5.0	1	06/10/14 16:53	06/11/14 22:27	00-73-7	
Hexachioro-1,3-buladiene	ND u	g/L		0.C	1	06/10/14 16:53	06/11/14 22:27	07-00-3	
Hexachiorobenzene	ND U	g/L		0.0	1	06/10/14 16:53	06/11/14 22:27	118-74-1	
Hexachlorocyclopentadiene	ND U	g/L		11.1	1	06/10/14 16:53	06/11/14 22:27	77-47-4	
	ND U	g/L		5.6	1	06/10/14 16:53	06/11/14 22:27	67-72-1	
Indeno(1,2,3-cd)pyrene	ND U	g/L		5.6	1	06/10/14 16:53	06/11/14 22:27	193-39-5	
Isophorone	ND u	g/L		11.1	1	06/10/14 16:53	06/11/14 22:27	78-59-1	
Naphthalene	ND u	g/L		5.6	1	06/10/14 16:53	06/11/14 22:27	91-20-3	
	ND u	g/L		5.6	1	06/10/14 16:53	06/11/14 22:27	98-95-3	
	ND u	g/∟		5.6	1	06/10/14 16:53	06/11/14 22:27	88-75-5	
4-INITrophenol	ND u	g/L		55.6	1	06/10/14 16:53	06/11/14 22:27	100-02-7	
N-Nitrosodimethylamine	ND u	g/L		5.6	1	06/10/14 16:53	06/11/14 22:27	62-75-9	
N-Nıtroso-di-n-propylamine	ND u	g/L		5.6	1	06/10/14 16:53	06/11/14 22:27	621-64-7	
N-Nitrosodiphenylamine	ND u	g/L		11.1	1	06/10/14 16:53	06/11/14 22:27	86-30-6	
Pentachlorophenol	ND u	g/L		11.1	1	06/10/14 16:53	06/11/14 22:27	87-86-5	

## **REPORT OF LABORATORY ANALYSIS**

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Project: 2014-093 Johnston 34182.1.2

Pace Project No.: 92204085

Parameters         Results         Units         Report Limit         DF         Prepared         Analyzed         CAS No.         Qual           625 MSSV         Analytical Method: EPA 625         Freparation Method: EPA 625         Freparation Method: EPA 626         Freparation Method: EPA 626 <td< th=""><th>Sample: 8-4 (TW)</th><th>Lab ID: 92204</th><th>4085002</th><th>Collected: 06/03/</th><th>14 08:05</th><th>Received: 06</th><th>/04/14 17:30 N</th><th>latrix: Water</th><th></th></td<>	Sample: 8-4 (TW)	Lab ID: 92204	4085002	Collected: 06/03/	14 08:05	Received: 06	/04/14 17:30 N	latrix: Water	
SSSV         Analytical Method: EPA 625         Preparation Method: EPA 625           Phenanihrene         ND ug/L         5.6         1         06/10/14 16:53         06/11/14 22:27         128-0-0           Prene         ND ug/L         5.6         1         06/10/14 16:53         06/11/14 22:27         128-0-0           12.4 - Trichlorobenzene         ND ug/L         5.6         1         06/10/14 16:53         06/11/14 22:27         128-0-0           Act-Trichlorobenzene         ND ug/L         5.6         1         06/10/14 16:53         06/11/14 22:27         128-0-0           Act-Trichlorobenzene         ND ug/L         5.10         06/10/14 16:53         06/11/14 22:27         128-0-0           Act-Trichlorobenzene         61 %         10-120         1         06/10/14 16:53         06/11/14 22:27         127-0-0           Act-Trichlorobenzene         9 %         10-120         1         06/10/14 16:53         06/11/14 22:27         127-24-2           Act-Trichlorobenzene         ND ug/L         0.00         1         06/05/14 16:42         71-42-2           Act-Trichlorobenzene         ND ug/L         0.50         1         06/05/14 16:42         74-97-5           Bromodachizomethane         ND ug/L         0.50         1	Parameters	Results	Units	Report Limit	DF	Prepared	Analyzed	CAS No.	Qual
Phenal         ND ug1         5.6         1         08/10/14 16:33         06/11/14 22:27         12:3           Prene         ND ug1         5.6         1         06/10/14 16:33         06/11/14 22:27         12:8-0-0           1:2.4-Trichlorobenzene         ND ug1         5.6         1         06/10/14 16:33         06/11/14 22:27         12:8-0-0           A:4-Trichlorobenzene         ND ug1         5.6         1         06/10/14 16:33         06/11/14 22:7         12:8-0-0           Surrogets         ND         1         06/10/14 16:33         06/11/14 22:7         21:8-0-0           Surrogets         1         06/10/14 16:33         06/11/14 22:7         12:16-0           Prenot-d6 (S)         39 %         10-120         1         06/10/14 16:33         06/11/14 2:27         12:7:48-10           Prenot-d6 (S)         39 %         10-120         1         06/10/14 16:33         06/11/14 2:27         12:7:48-10           Prenot-d6 (S)         39 %         10-120         1         06/10/14 16:33         06/11/14 2:27         12:7:48-10           Prenot-d6 (S)         39 %         10-120         1         06/00/14 16:42         12:7:43-12           Z-Horophenol (S)         100 %         ND ug1         0.60	625 MSSV	Analytical Metho	od: EPA 62	25 Preparation Meth	od: EPA	625			
Phenol         ND ug/L         5.6         1         06/10/14 1633         06/11/14 2227         108-95-2           1.2.4-Trichlorobenzene         ND ug/L         5.6         1         06/10/14 1633         06/11/14 2227         120-00-0           2.4.6-Trichlorobenzene         ND ug/L         1.1         1         06/10/14 1653         06/11/14 2227         188-06-2           Surrogates         NTOSEnzene-dS (S)         83         %         10-120         1         06/10/14 1653         06/11/14 2227         1165-00-3           TerphenyL-d1 (S)         69         %         11-131         1         06/10/14 1653         06/11/14 2227         171-85-10           Phenol-d6 (S)         29         10.3         01/11/14 2227         171-85-10         1           2-Floorobiphenol (S)         39         %         10-120         1         06/10/14 1653         06/11/14 2227         127-83-3           2-Floorobiphenol (S)         39         %         10-120         1         06/05/14 1642         71-83-2           2-Floorobiphenol (S)         109         %         1         06/05/14 1642         71-83-2           2-GOED MSV         Analytical Method: SM 6200B         1         06/05/14 1642         74-95-2	Phenanthrene	ND ug/l	_	5.6	1	06/10/14 16:53	06/11/14 22:27	85-01-8	
Pyrene         ND ug/L         5.6         1         06/10/14 1623         06/11/14 2227         129-0-0           1.2.4.Trichlorophenol         ND ug/L         1.1         1         06/10/14 1653         06/11/14 2227         88-06-2           Surrogets         -         -         06/10/14 1653         06/11/14 2227         38-06-2           Z-Fluorobipheny (S)         61 %         15-120         1         06/10/14 1653         06/11/14 2227         321-60-3           Z-Fluorobipheny (S)         69 %         1131         1         06/10/14 1653         06/11/14 2227         71-85-10           Z-Fluorobipheny (S)         39 %         10-120         1         06/10/14 1653         06/11/14 2227         71-85-10           Z-Horobneno (S)         39 %         10-120         1         06/10/14 1623         06/11/14 2227         71-85-10           Z-Horobneno (S)         39 %         10-137         1         06/10/14 1642         71-43-2           Bromochiconomethane         ND ug/L         0.50         1         06/05/14 1642         71-43-2           Bromochiconomethane         ND ug/L         0.50         1         06/05/14 1642         74-9-5           Bromochiconomethane         ND ug/L         0.50         1 <td>Phenol</td> <td>ND ug/L</td> <td>_</td> <td>5.6</td> <td>1</td> <td>06/10/14 16:53</td> <td>06/11/14 22:27</td> <td>108-95-2</td> <td></td>	Phenol	ND ug/L	_	5.6	1	06/10/14 16:53	06/11/14 22:27	108-95-2	
1.2.4.Trichlonobenzene       ND ug/L       5.6       1       06/10/14 16:53       06/11/14 22:27       129-82-1         Surrogates       ND ug/L       11.1       1       06/10/14 16:53       06/11/14 22:27       1465-60-0         Surrogates       88.3 %       10-120       1       06/10/14 16:53       06/11/14 22:27       1718-51-0         TerphenyL-d1 (S)       69 %       11-131       1       06/10/14 16:53       06/11/14 22:27       1718-51-0         Phenol-d6 (S)       29 %       10-120       1       06/10/14 16:53       06/11/14 22:27       1372-83-3         2-Flocrophenol (S)       39 %       10-120       1       06/10/14 16:53       06/11/14 22:27       1372-83-3         2-Flocrophenol (S)       39 %       10-120       1       06/10/14 16:23       06/11/14 22:27       1372-83-3         2-A6-Tribromophenol (S)       100 %       10-130       1       06/05/14 16:42       74-74-3         Bromochoromethane       ND ug/L       0.50       1       06/05/14 16:42       74-75-5         Bromochoromethane       ND ug/L       0.50       1       06/05/14 16:42       74-75-5         Bromochoromethane       ND ug/L       0.50       1       06/05/14 16:42       74-75-5 <t< td=""><td>Pyrene</td><td>ND ug/L</td><td>_</td><td>5.6</td><td>1</td><td>06/10/14 16:53</td><td>06/11/14 22:27</td><td>129-00-0</td><td></td></t<>	Pyrene	ND ug/L	_	5.6	1	06/10/14 16:53	06/11/14 22:27	129-00-0	
2.4.6.Tickloirophend         ND ugL         11.         1         06/10/14 16:53         06/11/14 22:27         1465:60-0           Nitrobarzanc-d5 (S)         63         %         15-120         1         06/10/14 16:53         06/11/14 22:27         1718-51-0           Tephenyl-14 (S)         69 %         11-131         1         06/10/14 18:53         06/11/14 22:27         1718-51-0           Tephenyl-14 (S)         99 %         10-120         1         06/10/14 18:53         06/11/14 22:27         13127-88-3           2-Hourophenol (S)         39 %         10-120         1         06/10/14 18:53         06/11/14 22:27         13127-88-3           2-Hourophenol (S)         39 %         10-130         1         06/05/14 16:42         71-43-2           2-Romobenzane         ND ugL         0.50         1         06/05/14 16:42         71-43-2           Bromochiloromethane         ND ugL         0.50         1         06/05/14 16:42         74-97-5           Bromochiloromethane         ND ugL         0.50         1         06/05/14 16:42         72-74           Bromochiloromethane         ND ugL         0.50         1         06/05/14 16:42         73-95           Bromochiloromethane         ND ugL         0.50	1,2,4-Trichlorobenzene	ND ug/L	_	5.6	1	06/10/14 16:53	06/11/14 22:27	120-82-1	
Surrogates         Surrogates           Surrogates         06/10/14 16:53         66/11/14 22:27         321-60-8           Terphenyl-d14 (S)         66 %         11-131         06/10/14 16:53         06/11/14 22:27         321-60-8           Terphenyl-d16 (S)         29 %         10-120         1         06/10/14 16:53         06/11/14 22:27         3127-38-3           2-Fluorophenol (S)         39 %         10-120         1         06/10/14 16:53         06/11/14 22:27         37-12-4           2.4.6.Tribromophenol (S)         100 %         100 %         06/10/14 16:53         06/11/14 22:27         37-12-4           2.4.6.Tribromophenol (S)         100 %         06/10         1         06/05/14 16:42         71-43-2           Bromochloromethane         ND ug/L         0.50         1         06/05/14 16:42         74-3-2           Bromochloromethane         ND ug/L         0.50         1         06/05/14 16:4	2,4,6-Trichlorophenol	ND ug/L	_	11.1	1	06/10/14 16:53	06/11/14 22:27	88-06-2	
Nincbezened5 (S)         B3 %         10-120         1         06/10/14 16:S3         06/11/14 22:27         1465-60-0           2-Fluoroblphenyl (S)         69 %         11-131         1         06/10/14 16:S3         06/11/14 22:27         1718-51-0           2-Fluoroblphenyl (S)         39 %         10-120         1         06/10/14 16:S3         06/11/14 22:27         13127-88-3           2-Fluoroblenol (S)         39 %         10-120         1         06/10/14 16:S3         06/11/14 22:27         13127-88-3           2-Fluoroblenol (S)         39 %         10-130         1         06/05/14 16:42         71.43-2           Benzene         ND ug/L         0.50         1         06/05/14 16:42         71.43-2           Bromochloromethane         ND ug/L         0.50         1         06/05/14 16:42         74.87-5           Bromochloromethane         ND ug/L         0.50         1         06/05/14 16:42         74.83-3           Bromochloromethane         ND ug/L         0.50         1         06/05/14 16:42         74.83-3           Bromochloromethane         ND ug/L         0.50         1         06/05/14 16:42         74.83-3           Bromochloromethane         ND ug/L         0.50         1         06/05/14 16:42 <td>Surrogates</td> <td>C C</td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td>	Surrogates	C C							
2-Fluorobjehenyl (S)         61 %         15-120         1         06/10/14 16:53         06/11/14 22:27         7321-80-8           Prephenyl-df (S)         29 %         10-120         1         06/10/14 16:53         06/11/14 22:27         7371-78-8-3           2-Fluorophenol (S)         39 %         10-120         1         06/10/14 16:53         06/11/14 22:27         7371-24           2-Fluorophenol (S)         100 %         10-137         1         06/10/14 16:32         06/11/14 22:27         7371-24           2-AG-Thrbromophenol (S)         100 %         10-137         1         06/05/14 16:42         71-43-2           Bromochoromethane         ND ug/L         0.50         1         06/05/14 16:42         71-43-2           Bromochoromethane         ND ug/L         0.50         1         06/05/14 16:42         72-74           Bromochoromethane         ND ug/L         0.50         1         06/05/14 16:42         75-25-2           Bromochoromethane         ND ug/L         0.50         1         06/05/14 16:42         75-25-2           Bromochoromethane         ND ug/L         0.50         1         06/05/14 16:42         75-25-2           Bromochoromethane         ND ug/L         0.50         1         06/05/14 16:4	Nitrobenzene-d5 (S)	83 %		10-120	1	06/10/14 16:53	06/11/14 22:27	4165-60-0	
TerphenyLift 4 (S)         69 %         11-131         1         06/10/14 16:53         06/11/14 22:27         7118-51-0           PhenoLdG (S)         29 %         10-120         1         06/10/14 16:53         06/11/14 22:27         1817-88-3           2-Fluorophenol (S)         100 %         10-120         1         06/10/14 16:53         06/11/14 22:27         1817-88-3           2,4.5 Tribromophenol (S)         100 %         10-137         1         06/05/14 16:42         71-43-2           Bromobersene         ND ug/L         0.50         1         06/05/14 16:42         71-43-2           Bromochhoromethane         ND ug/L         0.50         1         06/05/14 16:42         74-97-5           Bromochhoromethane         ND ug/L         0.50         1         06/05/14 16:42         74-97-5           Bromochhoromethane         ND ug/L         0.50         1         06/05/14 16:42         74-33-9           n-Butylbenzene         ND ug/L         0.50         1         06/05/14 16:42         74-97-5           Bromochhoromethane         ND ug/L         0.50         1         06/05/14 16:42         75-27-4           Bromochhoromethane         ND ug/L         0.50         1         06/05/14 16:42         75-93-3 </td <td>2-Fluorobiphenyl (S)</td> <td>61 %</td> <td></td> <td>15-120</td> <td>1</td> <td>06/10/14 16:53</td> <td>06/11/14 22:27</td> <td>321-60-8</td> <td></td>	2-Fluorobiphenyl (S)	61 %		15-120	1	06/10/14 16:53	06/11/14 22:27	321-60-8	
Phenol.dB (S)         29 %         10-120         1         06/10/14 16:53         06/11/14 22:27         13127-88-3           2-Fluorophenol (S)         100 %         10-130         1         06/10/14 16:53         06/11/14 22:27         1367-12:4           2-A6-Tribromophenol (S)         100 %         10-137         1         06/05/14 16:42         71-43-2           6200B MSV         Analytical Method: SM 6200B         1         06/05/14 16:42         71-43-2           Benzene         ND ug/L         0.50         1         06/05/14 16:42         74-37-5           Bromochoromethane         ND ug/L         0.50         1         06/05/14 16:42         74-97-5           Bromochoromethane         ND ug/L         0.50         1         06/05/14 16:42         74-97-5           Bromochoromethane         ND ug/L         0.50         1         06/05/14 16:42         74-97-5           Bromochoromethane         ND ug/L         0.50         1         06/05/14 16:42         74-93-9           Bromochoromethane         ND ug/L         0.50         1         06/05/14 16:42         74-93-9           Bromochoromethane         ND ug/L         0.50         1         06/05/14 16:42         18-98-8           Chiorobenzene	Terphenyl-d14 (S)	69 %		11-131	1	06/10/14 16:53	06/11/14 22:27	1718-51-0	
2-Fluorophenol (S)         39 %         10-120         1         06/10/14 16:53         06/11/14 22:27         367-12-4           2,4,6-Tribromophenol (S)         100 %         10         06/10/14 16:53         06/11/14 22:27         118-79-6           6200B MSV         Analytical Method: SM 6200B         5         5         5         5           Benzene         ND ug/L         0.50         1         06/05/14 16:42         71-43-2           Bromochirormethane         ND ug/L         0.50         1         06/05/14 16:42         74-97-5           Bromochirormethane         ND ug/L         0.50         1         06/05/14 16:42         75-25-2           Bromochirormethane         ND ug/L         0.50         1         06/05/14 16:42         98-06-6           Cathor tetrachioride         N	Phenol-d6 (S)	29 %		10-120	1	06/10/14 16:53	06/11/14 22:27	13127-88-3	
2.4.6-Tribromophenol (S)         100 %         10-137         1         06/10/141 16:33         06/11/14 22:27         118-79-6           6200B MSV         Analytical Method: SM 6200B               Benzene         ND ug/L         0.50         1         06/05/14 16:42         71-43-2           Bromobenzene         ND ug/L         0.50         1         06/05/14 16:42         74-97-5           Bromodinichloromethane         ND ug/L         0.50         1         06/05/14 16:42         75-27-4           Bromothane         ND ug/L         0.50         1         06/05/14 16:42         75-27-4           Bromothane         ND ug/L         0.50         1         06/05/14 16:42         74-97-5           Bromothane         ND ug/L         0.50         1         06/05/14 16:42         74-98-9           Bromothane         ND ug/L         0.50         1         06/05/14 16:42         74-83-9           ScaPutylbenzene         ND ug/L         0.50         1         06/05/14 16:42         74-83-9           Chrotophane         ND ug/L         0.50         1         06/05/14 16:42         75-0-3           Chlorobenzene         ND ug/L         0.50         1 <th< td=""><td>2-Fluorophenol (S)</td><td>39 %</td><td></td><td>10-120</td><td>1</td><td>06/10/14 16:53</td><td>06/11/14 22:27</td><td>367-12-4</td><td></td></th<>	2-Fluorophenol (S)	39 %		10-120	1	06/10/14 16:53	06/11/14 22:27	367-12-4	
Benzene         ND ug/L         0.50         1         06/05/14 16:42         71-43-2           Bromobenzene         ND ug/L         0.50         1         06/05/14 16:42         70-8-8-1           Bromochloromethane         ND ug/L         0.50         1         06/05/14 16:42         70-8-8-1           Bromochloromethane         ND ug/L         0.50         1         06/05/14 16:42         75-27-4           Bromodrom         ND ug/L         0.50         1         06/05/14 16:42         74-83-9           n-Butylbenzene         ND ug/L         0.50         1         06/05/14 16:42         74-83-9           charbon tetrachloride         ND ug/L         0.50         1         06/05/14 16:42         74-83-9           charbon tetrachloride         ND ug/L         0.50         1         06/05/14 16:42         74-83-3           chlorobenzene         ND ug/L         0.50         1         06/05/14 16:42 <td>2,4,6-Tribromophenol (S)</td> <td>100 %</td> <td></td> <td>10-137</td> <td>1</td> <td>06/10/14 16:53</td> <td>06/11/14 22:27</td> <td>118-79-6</td> <td></td>	2,4,6-Tribromophenol (S)	100 %		10-137	1	06/10/14 16:53	06/11/14 22:27	118-79-6	
Benzene         ND ug/L         0.50         1         06/05/14 16:42         71-43-2           Bromobenzene         ND ug/L         0.50         1         06/05/14 16:42         70-8-8-1           Bromodichloromethane         ND ug/L         0.50         1         06/05/14 16:42         74-87-5           Bromodichloromethane         ND ug/L         0.50         1         06/05/14 16:42         75-27-4           Bromodichloromethane         ND ug/L         0.50         1         06/05/14 16:42         75-27-2           Bromodichloromethane         ND ug/L         0.50         1         06/05/14 16:42         74-83-9           n-Butylbenzene         ND ug/L         0.50         1         06/05/14 16:42         164-81           sec-Butylbenzene         ND ug/L         0.50         1         06/05/14 16:42         164-80-7           Chiorobinzene         ND ug/L         0.50         1         06/05/14 16:42         164-80-7           Chiorobinzene         ND ug/L         0.50         1         06/05/14 16:42         164-83-8           Chiorobinzene         ND ug/L         0.50         1         06/05/14 16:42         164-83           2-Chiorobinzene         ND ug/L         0.50         1	6200B MSV	Analytical Metho	od: SM 62	00B					
Bromobenzene         ND ug/L         0.50         1         06/05/14 16:42         10:8-8-1           Bromochloromethane         ND ug/L         0.50         1         06/05/14 16:42         74-97-5           Bromochloromethane         ND ug/L         0.50         1         06/05/14 16:42         75-27-4           Bromotharomethane         ND ug/L         0.50         1         06/05/14 16:42         75-27-2           Bromotharomethane         ND ug/L         0.50         1         06/05/14 16:42         73-27-2           Bromotharomethane         ND ug/L         0.50         1         06/05/14 16:42         73-93-8           Bromotharomethane         ND ug/L         0.50         1         06/05/14 16:42         135-98-8           Earl-Butylbenzene         ND ug/L         0.50         1         06/05/14 16:42         108-90-7           Chlorobenzene         ND ug/L         0.50         1         06/05/14 16:42         108-90-7           Chlorobenzene         ND ug/L         0.50         1         06/05/14 16:42         108-90-7           Chlorobethane         ND ug/L         0.50         1         06/05/14 16:42         108-90-7           Chlorobethane         ND ug/L         0.50         1	Benzene	ND ug/l		0.50	1		06/05/14 16:42	71-43-2	
Biomochioromethane         ND ug/L         0.50         1         06/05/14         16.4.2         74-97-5           Bromochioromethane         ND ug/L         0.50         1         06/05/14         16.4.2         74-97-5           Bromochioromethane         ND ug/L         0.50         1         06/05/14         16.4.2         74-87-9           Bromochioromethane         ND ug/L         0.50         1         06/05/14         16.4.2         74-83-9           n-Butylbenzene         ND ug/L         0.50         1         06/05/14         16.4.2         74-83-9           ac-Butylbenzene         ND ug/L         0.50         1         06/05/14         16.4.2         74-83-9           chlorobenzene         ND ug/L         0.50         1         06/05/14         16.4.2         74-83-9           chlorobenzene         ND ug/L         0.50         1         06/05/14         16.4.2         74-87-3           chlorobenzene         ND ug/L         0.50         1         06/05/14         16.4.2         74-87-3           chlorobenzene         ND ug/L         0.50         1         06/05/14         16.4.2         96-92-8           chlorobenzene         ND ug/L         0.50         1	Bromobenzene	ND ug/L	-	0.50	1		06/05/14 16:42	108-86-1	
Biomodichloromethane         ND ug/L         0.50         1         06/05/14 16:42         75-27-4           Bromodichloromethane         ND ug/L         0.50         1         06/05/14 16:42         75-27-4           Bromodichloromethane         ND ug/L         0.50         1         06/05/14 16:42         75-25-2           Bromodichloromethane         ND ug/L         0.50         1         06/05/14 16:42         74-83-9           Bromodichloromethane         ND ug/L         0.50         1         06/05/14 16:42         135-98-8           etart-Butylbenzene         ND ug/L         0.50         1         06/05/14 16:42         58-08-6           Carbon tetrachloride         ND ug/L         0.50         1         06/05/14 16:42         58-06-6           Carbon tetrachloride         ND ug/L         0.50         1         06/05/14 16:42         57-66-3           Chlorobenzene         ND ug/L         0.50         1         06/05/14 16:42         67-66-3           Chlorobenzene         ND ug/L         0.50         1         06/05/14 16:42         95-49-8           4-Chlorobluene         ND ug/L         0.50         1         06/05/14 16:42         95-49-8           1_2-Dibromo-3-chloropropane         ND ug/L	Bromochloromethane	ND ug/l	-	0.50	1		06/05/14 16:42	74-97-5	
Display         Display <thdisplay< th=""> <th< td=""><td>Bromodichloromethane</td><td>ND ug/L</td><td>-</td><td>0.50</td><td>1</td><td></td><td>06/05/14 16:42</td><td>75-27-4</td><td></td></th<></thdisplay<>	Bromodichloromethane	ND ug/L	-	0.50	1		06/05/14 16:42	75-27-4	
Bromomethane         ND ug/L         0.50         1         06/05/14 16:42         74.82-9           n-Butylbenzene         ND ug/L         0.50         1         06/05/14 16:42         104.83-9           sec-Butylbenzene         ND ug/L         0.50         1         06/05/14 16:42         104.83-9           sec-Butylbenzene         ND ug/L         0.50         1         06/05/14 16:42         108-96-7           Carbon tetrachloride         ND ug/L         0.50         1         06/05/14 16:42         108-90-7           Chlorobenzene         ND ug/L         0.50         1         06/05/14 16:42         76-66-3           Chlorotom         ND ug/L         1.0         1         06/05/14 16:42         76-86-3           Chlorotom         ND ug/L         0.50         1         06/05/14 16:42         74-87-3           2-Chlorotoluene         ND ug/L         0.50         1         06/05/14 16:42         74-87-3           2-Chlorotoluene         ND ug/L         0.50         1         06/05/14 16:42         74-87-3           1_2-Dibromochlane (EDB)         ND ug/L         0.50         1         06/05/14 16:42         124-85-1           Dibromomethane         ND ug/L         0.50         1 <td< td=""><td>Bromoform</td><td>ND ug/L</td><td>-</td><td>0.50</td><td>1</td><td></td><td>06/05/14 16:42</td><td>75-25-2</td><td></td></td<>	Bromoform	ND ug/L	-	0.50	1		06/05/14 16:42	75-25-2	
Display         Display <t< td=""><td>Bromomethane</td><td>ND ug/L</td><td>-</td><td>5.0</td><td>1</td><td></td><td>06/05/14 16:42</td><td>74-83-9</td><td></td></t<>	Bromomethane	ND ug/L	-	5.0	1		06/05/14 16:42	74-83-9	
In Burghendering         ND ug/L         0.50         1         06/05/14         16.12         19.5           tert-Butylbenzene         ND ug/L         0.50         1         06/05/14         16.42         98-06-6           Carbon tetrachioride         ND ug/L         0.50         1         06/05/14         16.42         56-23-5           Chlorobenzene         ND ug/L         0.50         1         06/05/14         16.42         75-00-3           Chlorobenzene         ND ug/L         0.50         1         06/05/14         16.42         76-0-3           Chlorobenzene         ND ug/L         0.50         1         06/05/14         16.42         74-87-3           2-Chlorotoluene         ND ug/L         0.50         1         06/05/14         16.42         74-87-3           2-Chlorotoluene         ND ug/L         0.50         1         06/05/14         16.42         96-42-8           2-Chlorotoluene         ND ug/L         0.50         1         06/05/14         16.42         16-43-4           1,2-Dibromo-3-chloropropane         ND ug/L         0.50         1         06/05/14         16.42         16-43-4           1,2-Dibromoethane (EDB)         ND ug/L         0.50         1 <td>n-Butylbenzene</td> <td>ND ug/l</td> <td>-</td> <td>0.50</td> <td>1</td> <td></td> <td>06/05/14 16:42</td> <td>104-51-8</td> <td></td>	n-Butylbenzene	ND ug/l	-	0.50	1		06/05/14 16:42	104-51-8	
Side Day Bolt Zoho         ND ug/L         0.50         1         06/05/14 16.42         98-06-6           Carbon tetrachloride         ND ug/L         0.50         1         06/05/14 16.42         98-06-6           Carbon tetrachloride         ND ug/L         0.50         1         06/05/14 16.42         75-00-3           Chlorobenzene         ND ug/L         0.50         1         06/05/14 16.42         75-00-3           Chlorobenzene         ND ug/L         0.50         1         06/05/14 16.42         75-00-3           Chloroberzene         ND ug/L         0.50         1         06/05/14 16.42         95-49-8           Chlorobremethane         ND ug/L         0.50         1         06/05/14 16.42         95-49-8           4-Chlorotoluene         ND ug/L         0.50         1         06/05/14 16.42         95-49-8           4-Chlorotoluene         ND ug/L         0.50         1         06/05/14 16.42         95-92-8           Dibromochloromethane         ND ug/L         0.50         1         06/05/14 16.42         106-93-4           1.2-Dichlorobenzene         ND ug/L         0.50         1         06/05/14 16.42         148-9-3           1.2-Dichlorobenzene         ND ug/L         0.50	sec-Butylbenzene	ND ug/L	-	0.50	1		06/05/14 16:42	135-98-8	
Introduction         Ind by L         0.50         1         06/05/14 16:42         50:50           Carbon tetrachloride         ND ug/L         0.50         1         06/05/14 16:42         50:50:50           Chlorobenzene         ND ug/L         0.50         1         06/05/14 16:42         50:50:50           Chlorobenzene         ND ug/L         0.50         1         06/05/14 16:42         57:60-3           Chlorobenzene         ND ug/L         1.0         1         06/05/14 16:42         78:73           2-Chlorotoluene         ND ug/L         0.50         1         06/05/14 16:42         95:49:8           4-Chlorotoluene         ND ug/L         0.50         1         06/05/14 16:42         16:42:10:4:4:14           1,2-Dibromo-3-chloropropane         ND ug/L         0.50         1         06/05/14 16:42         14:4:4:14           1,2-Dibromoethane (EDB)         ND ug/L         0.50         1         06/05/14 16:42         14:4:73:1           1,2-Dichlorobenzene         ND ug/L         0.50         1         06/05/14 16:42         54:1-73:1           1,3-Dichlorobenzene         ND ug/L         0.50         1         06/05/14 16:42         54:1-73:1           1,4-Dichlorobenzene         ND ug/L <t< td=""><td>tert-Butylbenzene</td><td>ND ug/L</td><td>-</td><td>0.50</td><td>1</td><td></td><td>06/05/14 16:42</td><td>98-06-6</td><td></td></t<>	tert-Butylbenzene	ND ug/L	-	0.50	1		06/05/14 16:42	98-06-6	
Diborn Chromethane         ND ug/L         0.55         1         06/05/14 16:42         08-90-7           Chlorobertane         ND ug/L         1.0         1         06/05/14 16:42         75-00-3           Chlorobertane         ND ug/L         0.50         1         06/05/14 16:42         75-00-3           Chlorobertane         ND ug/L         0.50         1         06/05/14 16:42         75-66-3           Chlorobertane         ND ug/L         0.50         1         06/05/14 16:42         74-87-3           2-Chlorobluene         ND ug/L         0.50         1         06/05/14 16:42         95-49-8           4-Chlorobluene         ND ug/L         0.50         1         06/05/14 16:42         96-12-8           Dibromochloromethane         ND ug/L         0.50         1         06/05/14 16:42         106-93-4           1,2-Dibromoethane (EDB)         ND ug/L         0.50         1         06/05/14 16:42         14-4-81           1,3-Dichlorobenzene         ND ug/L         0.50         1         06/05/14 16:42         24-48-1           1,4-Dichlorobenzene         ND ug/L         0.50         1         06/05/14 16:42         74-95-3           1,2-Dichlorobenzene         ND ug/L         0.50	Carbon tetrachloride	ND ug/L	-	0.50	1		06/05/14 16:42	56-23-5	
Onlocation         ND         ug/L         1.0         1         06/05/14         16.42         75-00-3           Chlorotorm         ND         ug/L         0.50         1         06/05/14         16.42         75-00-3           Chlorotoluene         ND         ug/L         1.0         1         06/05/14         16.42         74-87-3           2-Chlorotoluene         ND         ug/L         0.50         1         06/05/14         16.42         95-49-8           4-Chlorotoluene         ND         ug/L         0.50         1         06/05/14         16.42         96-42-8           Dibromo-3-chloropropane         ND         ug/L         1.0         1         06/05/14         16.42         96-42-8           Dibromochloromethane         ND         ug/L         0.50         1         06/05/14         16.42         96-93-4           1,2-Dibromosethane (EDB)         ND         ug/L         0.50         1         06/05/14         16.42         96-50-1           1,3-Dichlorobenzene         ND         ug/L         0.50         1         06/05/14         16.42         541-73-1           1,4-Dichlorobenzene         ND         ug/L         0.50         1         06/05/14 </td <td>Chlorobenzene</td> <td>ND ug/L</td> <td>-</td> <td>0.50</td> <td>1</td> <td></td> <td>06/05/14 16:42</td> <td>108-90-7</td> <td></td>	Chlorobenzene	ND ug/L	-	0.50	1		06/05/14 16:42	108-90-7	
Onlocidation         ND         ug/L         0.50         1         06/05/14         16:42         67-66-3           Chloroform         ND         ug/L         1.0         1         06/05/14         16:42         67-66-3           Chloroform         ND         ug/L         0.50         1         06/05/14         16:42         74-87-3           2-Chlorotoluene         ND         ug/L         0.50         1         06/05/14         16:42         96-49-8           4-Chlorotoluene         ND         ug/L         0.50         1         06/05/14         16:42         96-12-8           Dibromochloromethane         ND         ug/L         0.50         1         06/05/14         16:42         106-93-4           1,2-Dibromo-3-chloropropane         ND         ug/L         0.50         1         06/05/14         16:42         106-93-4           1,2-Dibromoethane (EDB)         ND         ug/L         0.50         1         06/05/14         16:42         106-46-7           1,4-Dichlorobenzene         ND         ug/L         0.50         1         06/05/14         16:42         106-46-7           Dichlorodfiluoromethane         ND         ug/L         0.50         1 <td< td=""><td>Chloroethane</td><td>ND ug/l</td><td>-</td><td>1.0</td><td>1</td><td></td><td>06/05/14 16:42</td><td>75-00-3</td><td></td></td<>	Chloroethane	ND ug/l	-	1.0	1		06/05/14 16:42	75-00-3	
Onlocitient         ND         ug/L         0.55         1         06/05/14         16:42         74-87-3           2-Chlorotoluene         ND         ug/L         0.50         1         06/05/14         16:42         74-87-3           2-Chlorotoluene         ND         ug/L         0.50         1         06/05/14         16:42         106-43-4           1,2-Dibromo-3-chloropropane         ND         ug/L         1.0         1         06/05/14         16:42         106-43-4           1,2-Dibromo-3-chloropropane         ND         ug/L         0.50         1         06/05/14         16:42         124-48-1           1,2-Dibromoethane (EDB)         ND         ug/L         0.50         1         06/05/14         16:42         74-95-3           1,2-Dichlorobenzene         ND         ug/L         0.50         1         06/05/14         16:42         74-95-3           1,2-Dichlorobenzene         ND         ug/L         0.50         1         06/05/14         16:42         74-95-3           1,2-Dichlorobenzene         ND         ug/L         0.50         1         06/05/14         16:42         74-87-3           1,3-Dichlorobenzene         ND         ug/L         0.50 <td< td=""><td>Chloroform</td><td>ND ug/L</td><td>-</td><td>0.50</td><td>1</td><td></td><td>06/05/14 16:42</td><td>67-66-3</td><td></td></td<>	Chloroform	ND ug/L	-	0.50	1		06/05/14 16:42	67-66-3	
Ontoincidiality         ND         ug/L         1.5         1         Obtoine 1000 14 16:42         14 0 1           2-Chlorotoluene         ND         ug/L         0.50         1         06/05/14 16:42         106-43-4           1,2-Dibromo-3-chloropropane         ND         ug/L         1.0         1         06/05/14 16:42         96-12-8           Dibromochloromethane         ND         ug/L         0.50         1         06/05/14 16:42         124-48-1           1,2-Dibromoethane (EDB)         ND         ug/L         0.50         1         06/05/14 16:42         74-95-3           Dibromomethane         ND         ug/L         0.50         1         06/05/14 16:42         74-95-3           1,2-Dichlorobenzene         ND         ug/L         0.50         1         06/05/14 16:42         74-95-3           1,3-Dichlorobenzene         ND         ug/L         0.50         1         06/05/14 16:42         74-95-3           1,4-Dichlorobenzene         ND         ug/L         0.50         1         06/05/14 16:42         74-95-3           1,4-Dichlorobenzene         ND         ug/L         0.50         1         06/05/14 16:42         75-71-8           1,4-Dichlorobenzene         ND	Chloromethane	ND ug/L	-	1.0	1		06/05/14 16:42	74-87-3	
Productive         ND ug/L         0.50         1         00/05/14 16:42         50-95-0           4-Chlorotoluene         ND ug/L         0.50         1         06/05/14 16:42         106-43-4           1,2-Dibromo-3-chloropropane         ND ug/L         0.50         1         06/05/14 16:42         96-12-8           Dibromochloromethane         ND ug/L         0.50         1         06/05/14 16:42         124-48-1           1,2-Dibromoethane (EDB)         ND ug/L         0.50         1         06/05/14 16:42         74-95-3           1,2-Dichlorobenzene         ND ug/L         0.50         1         06/05/14 16:42         95-50-1           1,3-Dichlorobenzene         ND ug/L         0.50         1         06/05/14 16:42         95-50-1           1,4-Dichlorobenzene         ND ug/L         0.50         1         06/05/14 16:42         54-67           1,4-Dichlorobenzene         ND ug/L         0.50         1         06/05/14 16:42         57-71-8           1,1-Dichloroethane         ND ug/L         0.50         1         06/05/14 16:42         75-34-3           1,2-Dichloroethane         ND ug/L         0.50         1         06/05/14 16:42         75-35-4           1,1-Dichloroethene         ND ug/L	2-Chlorotoluene	ND ug/L	-	0.50	1		06/05/14 16:42	05-40-8	
1,2-Dibromo-3-chloropropane       ND ug/L       1.0       1       06/05/14 16:42       96-12-8         Dibromochloromethane       ND ug/L       0.50       1       06/05/14 16:42       124-48-1         1,2-Dibromochloromethane       ND ug/L       0.50       1       06/05/14 16:42       124-48-1         1,2-Dibromoethane (EDB)       ND ug/L       0.50       1       06/05/14 16:42       74-95-3         1,2-Dichlorobenzene       ND ug/L       0.50       1       06/05/14 16:42       95-50-1         1,3-Dichlorobenzene       ND ug/L       0.50       1       06/05/14 16:42       95-50-1         1,3-Dichlorobenzene       ND ug/L       0.50       1       06/05/14 16:42       75-71-8         1,4-Dichlorobenzene       ND ug/L       0.50       1       06/05/14 16:42       75-71-8         1,1-Dichloroethane       ND ug/L       0.50       1       06/05/14 16:42       75-34-3         1,2-Dichloroethane       ND ug/L       0.50       1       06/05/14 16:42       75-35-4         1,1-Dichloroethene       ND ug/L       0.50       1       06/05/14 16:42       75-35-4         1,2-Dichloroethene       ND ug/L       0.50       1       06/05/14 16:42       75-35-4		ND ug/L	-	0.50	1		06/05/14 16:42	106-43-4	
ND ug/L       0.50       1       06/05/14 16:42       124-48-1         Dibromochhane (EDB)       ND ug/L       0.50       1       06/05/14 16:42       124-48-1         Dibromomethane       ND ug/L       0.50       1       06/05/14 16:42       124-48-1         Dibromomethane       ND ug/L       0.50       1       06/05/14 16:42       124-48-1         J.2-Dichlorobenzene       ND ug/L       0.50       1       06/05/14 16:42       95-50-1         1,3-Dichlorobenzene       ND ug/L       0.50       1       06/05/14 16:42       95-50-1         1,4-Dichlorobenzene       ND ug/L       0.50       1       06/05/14 16:42       541-73-1         1,4-Dichlorobenzene       ND ug/L       0.50       1       06/05/14 16:42       75-71-8         1,4-Dichloroethane       ND ug/L       0.50       1       06/05/14 16:42       75-71-8         1,1-Dichloroethane       ND ug/L       0.50       1       06/05/14 16:42       75-34-3         1,2-Dichloroethane       ND ug/L       0.50       1       06/05/14 16:42       107-06-2         1,1-Dichloroethene       ND ug/L       0.50       1       06/05/14 16:42       156-59-2         trans-1,2-Dichloroethene       ND ug/L	1 2-Dibromo-3-chloropropape	ND ug/L	-	1.0	1		06/05/14 16:42	96-12-8	
Initial constraints       ND ug/L       0.50       1       06/05/14 16.42       124-40 1         1,2-Dibromoethane (EDB)       ND ug/L       0.50       1       06/05/14 16:42       74-95-3         1,2-Dichlorobenzene       ND ug/L       0.50       1       06/05/14 16:42       74-95-3         1,3-Dichlorobenzene       ND ug/L       0.50       1       06/05/14 16:42       95-50-1         1,3-Dichlorobenzene       ND ug/L       0.50       1       06/05/14 16:42       541-73-1         1,4-Dichlorobenzene       ND ug/L       0.50       1       06/05/14 16:42       75-71-8         1,1-Dichloroethane       ND ug/L       0.50       1       06/05/14 16:42       75-71-8         1,1-Dichloroethane       ND ug/L       0.50       1       06/05/14 16:42       75-71-8         1,2-Dichloroethane       ND ug/L       0.50       1       06/05/14 16:42       75-34-3         1,2-Dichloroethene       ND ug/L       0.50       1       06/05/14 16:42       75-35-4         1,1-Dichloroethene       ND ug/L       0.50       1       06/05/14 16:42       75-35-4         1,2-Dichloroethene       ND ug/L       0.50       1       06/05/14 16:42       75-35-4         1,2-Dichlor	Dibromochloromethane	ND ug/L	-	0.50	1		06/05/14 16:42	124-48-1	
ND ug/L       0.50       1       06/05/14 16:42       74-95-3         J.2-Dichlorobenzene       ND ug/L       0.50       1       06/05/14 16:42       95-50-1         J.3-Dichlorobenzene       ND ug/L       0.50       1       06/05/14 16:42       95-50-1         J.3-Dichlorobenzene       ND ug/L       0.50       1       06/05/14 16:42       95-50-1         J.4-Dichlorobenzene       ND ug/L       0.50       1       06/05/14 16:42       95-71-8         J.4-Dichlorobenzene       ND ug/L       0.50       1       06/05/14 16:42       75-71-8         J.1-Dichloroethane       ND ug/L       0.50       1       06/05/14 16:42       75-34-3         J.2-Dichloroethane       ND ug/L       0.50       1       06/05/14 16:42       75-35-4         J.1-Dichloroethene       ND ug/L       0.50       1       06/05/14 16:42       75-35-4         Lis-J.2-Dichloroethene       ND ug/L       0.50       1       06/05/14 16:42       75-92         Lis-J.2-Dichloroethene       ND ug/L       0.50       1       06/05/14 16:42       78-87-5         J.2-Dichloropropane       ND ug/L       0.50       1       06/05/14 16:42       78-87-5         J.3-Dichloropropane       ND ug/L <td>1 2-Dibromoethane (EDB)</td> <td>ND ug/L</td> <td>-</td> <td>0.50</td> <td>1</td> <td></td> <td>06/05/14 16:42</td> <td>106-93-4</td> <td></td>	1 2-Dibromoethane (EDB)	ND ug/L	-	0.50	1		06/05/14 16:42	106-93-4	
1,2-Dichlorobenzene       ND ug/L       0.50       1       06/05/14 16:42       95-50-1         1,3-Dichlorobenzene       ND ug/L       0.50       1       06/05/14 16:42       95-50-1         1,3-Dichlorobenzene       ND ug/L       0.50       1       06/05/14 16:42       541-73-1         1,4-Dichlorobenzene       ND ug/L       0.50       1       06/05/14 16:42       541-73-1         1,4-Dichlorobenzene       ND ug/L       0.50       1       06/05/14 16:42       75-71-8         1,1-Dichloroethane       ND ug/L       0.50       1       06/05/14 16:42       75-34-3         1,2-Dichloroethane       ND ug/L       0.50       1       06/05/14 16:42       107-06-2         1,1-Dichloroethane       ND ug/L       0.50       1       06/05/14 16:42       107-06-2         1,1-Dichloroethene       ND ug/L       0.50       1       06/05/14 16:42       156-59-2         1,1-Dichloroethene       ND ug/L       0.50       1       06/05/14 16:42       156-59-2         trans-1,2-Dichloroethene       ND ug/L       0.50       1       06/05/14 16:42       156-60-5         1,2-Dichloropropane       ND ug/L       0.50       1       06/05/14 16:42       142-28-9         2,2	Dibromomethane	ND ug/L	-	0.50	1		06/05/14 16:42	74-95-3	
1,3-Dichlorobenzene       ND ug/L       0.50       1       06/05/14 16:42       541-73-1         1,3-Dichlorobenzene       ND ug/L       0.50       1       06/05/14 16:42       541-73-1         1,4-Dichlorobenzene       ND ug/L       0.50       1       06/05/14 16:42       106-46-7         Dichlorodifluoromethane       ND ug/L       0.50       1       06/05/14 16:42       75-71-8         1,1-Dichloroethane       ND ug/L       0.50       1       06/05/14 16:42       75-34-3         1,2-Dichloroethane       ND ug/L       0.50       1       06/05/14 16:42       75-35-4         1,1-Dichloroethane       ND ug/L       0.50       1       06/05/14 16:42       75-35-4         1,1-Dichloroethene       ND ug/L       0.50       1       06/05/14 16:42       156-59-2         1,1-Dichloroethene       ND ug/L       0.50       1       06/05/14 16:42       156-60-5         1,2-Dichloroethene       ND ug/L       0.50       1       06/05/14 16:42       156-60-5         1,2-Dichloropropane       ND ug/L       0.50       1       06/05/14 16:42       142-28-9         2,2-Dichloropropane       ND ug/L       0.50       1       06/05/14 16:42       594-20-7         1,1-D	1 2-Dichlorobenzene	ND ug/L	-	0.50	1		06/05/14 16:42	95-50-1	
1,4-Dichlorobenzene       ND ug/L       0.50       1       06/05/14 16:42       341 75 1         1,4-Dichlorobenzene       ND ug/L       0.50       1       06/05/14 16:42       106-46-7         Dichlorodifluoromethane       ND ug/L       0.50       1       06/05/14 16:42       75-71-8         1,1-Dichloroethane       ND ug/L       0.50       1       06/05/14 16:42       75-34-3         1,2-Dichloroethane       ND ug/L       0.50       1       06/05/14 16:42       107-06-2         1,1-Dichloroethene       ND ug/L       0.50       1       06/05/14 16:42       75-35-4         cis-1,2-Dichloroethene       ND ug/L       0.50       1       06/05/14 16:42       156-59-2         trans-1,2-Dichloroethene       ND ug/L       0.50       1       06/05/14 16:42       156-60-5         1,2-Dichloropropane       ND ug/L       0.50       1       06/05/14 16:42       156-60-5         1,2-Dichloropropane       ND ug/L       0.50       1       06/05/14 16:42       142-28-9         2,2-Dichloropropane       ND ug/L       0.50       1       06/05/14 16:42       142-28-9         2,2-Dichloropropane       ND ug/L       0.50       1       06/05/14 16:42       594-20-7	1.3-Dichlorobenzene	ND ug/L	-	0.50	1		06/05/14 16:42	541-73-1	
1,4-Dichlorodifluoromethane       ND ug/L       0.50       1       06/05/14 16:42       75-71-8         1,1-Dichloroethane       ND ug/L       0.50       1       06/05/14 16:42       75-34-3         1,2-Dichloroethane       ND ug/L       0.50       1       06/05/14 16:42       75-34-3         1,2-Dichloroethane       ND ug/L       0.50       1       06/05/14 16:42       107-06-2         1,1-Dichloroethene       ND ug/L       0.50       1       06/05/14 16:42       75-35-4         cis-1,2-Dichloroethene       ND ug/L       0.50       1       06/05/14 16:42       156-59-2         trans-1,2-Dichloroethene       ND ug/L       0.50       1       06/05/14 16:42       156-60-5         1,2-Dichloropropane       ND ug/L       0.50       1       06/05/14 16:42       156-60-5         1,2-Dichloropropane       ND ug/L       0.50       1       06/05/14 16:42       142-28-9         1,3-Dichloropropane       ND ug/L       0.50       1       06/05/14 16:42       142-28-9         2,2-Dichloropropane       ND ug/L       0.50       1       06/05/14 16:42       594-20-7         1,1-Dichloropropane       ND ug/L       0.50       1       06/05/14 16:42       594-20-7	1.4-Dichlorobenzene	ND ug/L	-	0.50	1		06/05/14 16:42	106-46-7	
1,1-Dichloroethane       ND ug/L       0.50       1       06/05/14 16:42       75-34-3         1,2-Dichloroethane       ND ug/L       0.50       1       06/05/14 16:42       75-34-3         1,2-Dichloroethane       ND ug/L       0.50       1       06/05/14 16:42       107-06-2         1,1-Dichloroethene       ND ug/L       0.50       1       06/05/14 16:42       75-35-4         1,1-Dichloroethene       ND ug/L       0.50       1       06/05/14 16:42       156-59-2         trans-1,2-Dichloroethene       ND ug/L       0.50       1       06/05/14 16:42       156-60-5         1,2-Dichloropropane       ND ug/L       0.50       1       06/05/14 16:42       78-87-5         1,3-Dichloropropane       ND ug/L       0.50       1       06/05/14 16:42       142-28-9         2,2-Dichloropropane       ND ug/L       0.50       1       06/05/14 16:42       594-20-7         1,1-Dichloropropane       ND ug/L       0.50       1       06/05/14 16:42       594-20-7         1,1-Dichloropropane       ND ug/L       0.50       1       06/05/14 16:42       563-58-6	Dichlorodifluoromethane	ND ug/L	-	0.50	1		06/05/14 16:42	75-71-8	
1,2-Dichloroethane       ND ug/L       0.50       1       06/05/14 16:42       107-06-2         1,2-Dichloroethane       ND ug/L       0.50       1       06/05/14 16:42       107-06-2         1,1-Dichloroethane       ND ug/L       0.50       1       06/05/14 16:42       75-35-4         cis-1,2-Dichloroethane       ND ug/L       0.50       1       06/05/14 16:42       156-59-2         trans-1,2-Dichloroethane       ND ug/L       0.50       1       06/05/14 16:42       156-60-5         1,2-Dichloropropane       ND ug/L       0.50       1       06/05/14 16:42       78-87-5         1,3-Dichloropropane       ND ug/L       0.50       1       06/05/14 16:42       142-28-9         2,2-Dichloropropane       ND ug/L       0.50       1       06/05/14 16:42       594-20-7         1,1-Dichloropropane       ND ug/L       0.50       1       06/05/14 16:42       594-20-7         1,1-Dichloropropane       ND ug/L       0.50       1       06/05/14 16:42       563-58-6		ND ug/L	-	0.50	1		06/05/14 16:42	75-34-3	
1,1-Dichloroethene       ND ug/L       0.50       1       06/05/14 16:42       75-35-4         1,1-Dichloroethene       ND ug/L       0.50       1       06/05/14 16:42       75-35-4         cis-1,2-Dichloroethene       ND ug/L       0.50       1       06/05/14 16:42       156-59-2         trans-1,2-Dichloroethene       ND ug/L       0.50       1       06/05/14 16:42       156-60-5         1,2-Dichloropropane       ND ug/L       0.50       1       06/05/14 16:42       78-87-5         1,3-Dichloropropane       ND ug/L       0.50       1       06/05/14 16:42       142-28-9         2,2-Dichloropropane       ND ug/L       0.50       1       06/05/14 16:42       594-20-7         1,1-Dichloropropane       ND ug/L       0.50       1       06/05/14 16:42       593-58-6	1.2-Dichloroethane	ND ug/L	-	0.50	1		06/05/14 16:42	107-06-2	
ND ug/L       0.50       1       06/05/14       16.42       75-53-4         cis-1,2-Dichloroethene       ND ug/L       0.50       1       06/05/14       16:42       156-59-2         trans-1,2-Dichloroethene       ND ug/L       0.50       1       06/05/14       16:42       156-60-5         1,2-Dichloropropane       ND ug/L       0.50       1       06/05/14       16:42       78-87-5         1,3-Dichloropropane       ND ug/L       0.50       1       06/05/14       16:42       142-28-9         2,2-Dichloropropane       ND ug/L       0.50       1       06/05/14       16:42       594-20-7         1,-Dichloropropane       ND ug/L       0.50       1       06/05/14       16:42       594-20-7         1,-Dichloropropane       ND ug/L       0.50       1       06/05/14       16:42       594-20-7	1 1-Dichloroethene		-	0.50	1		06/05/14 16:42	75-35-4	
ND ug/L       0.50       1       06/05/14       10.42       156-03-2         trans-1,2-Dichloroethene       ND ug/L       0.50       1       06/05/14       16:42       156-60-5         1,2-Dichloropropane       ND ug/L       0.50       1       06/05/14       16:42       78-87-5         1,3-Dichloropropane       ND ug/L       0.50       1       06/05/14       16:42       142-28-9         2,2-Dichloropropane       ND ug/L       0.50       1       06/05/14       16:42       594-20-7         1,1-Dichloropropane       ND ug/L       0.50       1       06/05/14       16:42       563-58-6	cis-1 2-Dichloroethene		-	0.50	1		06/05/14 16:42	156-50-2	
ND ug/L     0.50     1     06/05/14 16:42     136-00-5       1,2-Dichloropropane     ND ug/L     0.50     1     06/05/14 16:42     78-87-5       1,3-Dichloropropane     ND ug/L     0.50     1     06/05/14 16:42     142-28-9       2,2-Dichloropropane     ND ug/L     0.50     1     06/05/14 16:42     594-20-7       1,1-Dichloropropane     ND ug/L     0.50     1     06/05/14 16:42     593-58-6	trans-1 2-Dichloroethene		-	0.50	1		06/05/14 16:42	156-60-5	
ND ug/L     0.50     1     06/05/14 10.42     78-07-5       1,3-Dichloropropane     ND ug/L     0.50     1     06/05/14 16:42     142-28-9       2,2-Dichloropropane     ND ug/L     0.50     1     06/05/14 16:42     594-20-7       1,1-Dichloropropane     ND ug/L     0.50     1     06/05/14 16:42     594-20-7	1 2-Dichloropropape		-	0.50	1		06/05/14 10.42	78-87-5	
ND ug/L         0.50         1         06/05/14         10.42         142-26-9           2,2-Dichloropropane         ND ug/L         0.50         1         06/05/14         16:42         594-20-7           1.1-Dichloropropane         ND ug/L         0.50         1         06/05/14         16:42         594-20-7	1 3-Dichloropropane		-	0.50	1		06/05/14 16:42	142-28-0	
1.1-Dichloropropene ND ug/L 0.50 1 06/05/14 10.42 594-20-7	2 2-Dichloropropane		-	0.50	1		06/05/14 16:42	594-20-7	
	1.1-Dichloropropene	ND ug/l	-	0.50	1		06/05/14 16:42	563-58-6	



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Pace Project No.: 92204085

Sample: 8-4 (TW)	Lab ID: 9220408500	2 Collected: 06/03/14 08	:05 Received:	06/04/14 17:30 N	latrix: Water	
Parameters	ResultsUnits	Report Limit DF	Prepared	Analyzed	CAS No.	Qual
6200B MSV	Analytical Method: SM	6200B				
cis-1,3-Dichloropropene	ND ug/L	0.50 1		06/05/14 16:42	10061-01-5	
trans-1,3-Dichloropropene	ND ug/L	0.50 1		06/05/14 16:42	10061-02-6	
Diisopropyl ether	ND ug/L	0.50 1		06/05/14 16:42	108-20-3	
Ethylbenzene	ND ug/L	0.50 1		06/05/14 16:42	100-41-4	
Hexachloro-1,3-butadiene	ND ug/L	2.0 1		06/05/14 16:42	87-68-3	
Isopropylbenzene (Cumene)	ND ug/L	0.50 1		06/05/14 16:42	98-82-8	
Methylene Chloride	ND ug/L	2.0 1		06/05/14 16:42	75-09-2	
Methyl-tert-butyl ether	ND ug/L	0.50 1		06/05/14 16:42	1634-04-4	
Naphthalene	ND ug/L	2.0 1		06/05/14 16:42	91-20-3	
n-Propylbenzene	ND ug/L	0.50 1		06/05/14 16:42	103-65-1	
Styrene	ND ug/L	0.50 1		06/05/14 16:42	100-42-5	
1,1,1,2-Tetrachloroethane	ND ug/L	0.50 1		06/05/14 16:42	630-20-6	
1,1,2,2-Tetrachloroethane	ND ug/L	0.50 1		06/05/14 16:42	79-34-5	
Tetrachloroethene	ND ug/L	0.50 1		06/05/14 16:42	127-18-4	
Toluene	ND ug/L	0.50 1		06/05/14 16:42	108-88-3	
1,2,3-Trichlorobenzene	ND ug/L	2.0 1		06/05/14 16:42	87-61-6	
1,2,4-Trichlorobenzene	ND ug/L	2.0 1		06/05/14 16:42	120-82-1	
1,1,1-Trichloroethane	ND ug/L	0.50 1		06/05/14 16:42	71-55-6	
1,1,2-Trichloroethane	ND ug/L	0.50 1		06/05/14 16:42	79-00-5	
Trichloroethene	ND ug/L	0.50 1		06/05/14 16:42	79-01-6	
Trichlorofluoromethane	ND ug/L	1.0 1		06/05/14 16:42	75-69-4	
1,2,3-Trichloropropane	ND ug/L	0.50 1		06/05/14 16:42	96-18-4	
1,2,4-Trimethylbenzene	ND ug/L	0.50 1		06/05/14 16:42	95-63-6	
1,3,5-Trimethylbenzene	ND ug/L	0.50 1		06/05/14 16:42	108-67-8	
Vinyl chloride	ND ug/L	1.0 1		06/05/14 16:42	75-01-4	
m&p-Xylene	ND ug/L	1.0 1		06/05/14 16:42	179601-23-1	
o-Xylene	ND ug/L	0.50 1		06/05/14 16:42	95-47-6	
Surrogates	5					
1,2-Dichloroethane-d4 (S)	97 %	70-130 1		06/05/14 16:42	17060-07-0	
4-Bromofluorobenzene (S)	97 %	70-130 1		06/05/14 16:42	460-00-4	
Toluene-d8 (S)	101 %	70-130 1		06/05/14 16:42	2037-26-5	



## **QUALITY CONTROL DATA**

Project: 2014-093 Johnston 34182.1.2

Pace Project No.: 92204085

Pace Project No.: 92204085					
QC Batch: MSV/27102		Analysis Meth	od: SN	/I 6200B	
QC Batch Method: SM 6200B		Analysis Desc	ription: 62	00B MSV	
Associated Lab Samples: 922040	85002				
METHOD BLANK: 1214892		Matrix:	Water		
Associated Lab Samples: 922040	85002				
		Blank	Reporting		
Parameter	Units	Result	Limit	Analyzed	Qualifiers
1,1,1,2-Tetrachloroethane	ug/L	ND	0.50	06/05/14 14:46	
1,1,1-Trichloroethane	ug/L	ND	0.50	06/05/14 14:46	
1,1,2,2-Tetrachloroethane	ug/L	ND	0.50	06/05/14 14:46	
1,1,2-Trichloroethane	ug/L	ND	0.50	06/05/14 14:46	
1,1-Dichloroethane	ua/L	ND	0.50	06/05/14 14:46	

Parameter	Units	Result	Limit	Analyzed	Qualifiers
1,1,1,2-Tetrachloroethane	ug/L		0.50	06/05/14 14:46	
1,1,1-Trichloroethane	ug/L	ND	0.50	06/05/14 14:46	
1,1,2,2-Tetrachloroethane	ug/L	ND	0.50	06/05/14 14:46	
1,1,2-Trichloroethane	ug/L	ND	0.50	06/05/14 14:46	
1,1-Dichloroethane	ug/L	ND	0.50	06/05/14 14:46	
1,1-Dichloroethene	ug/L	ND	0.50	06/05/14 14:46	
1,1-Dichloropropene	ug/L	ND	0.50	06/05/14 14:46	
1,2,3-Trichlorobenzene	ug/L	ND	2.0	06/05/14 14:46	
1,2,3-Trichloropropane	ug/L	ND	0.50	06/05/14 14:46	
1,2,4-Trichlorobenzene	ug/L	ND	2.0	06/05/14 14:46	
1,2,4-Trimethylbenzene	ug/L	ND	0.50	06/05/14 14:46	
1,2-Dibromo-3-chloropropane	ug/L	ND	1.0	06/05/14 14:46	
1,2-Dibromoethane (EDB)	ug/L	ND	0.50	06/05/14 14:46	
1,2-Dichlorobenzene	ug/L	ND	0.50	06/05/14 14:46	
1,2-Dichloroethane	ug/L	ND	0.50	06/05/14 14:46	
1,2-Dichloropropane	ug/L	ND	0.50	06/05/14 14:46	
1,3,5-Trimethylbenzene	ug/L	ND	0.50	06/05/14 14:46	
1,3-Dichlorobenzene	ug/L	ND	0.50	06/05/14 14:46	
1,3-Dichloropropane	ug/L	ND	0.50	06/05/14 14:46	
1,4-Dichlorobenzene	ug/L	ND	0.50	06/05/14 14:46	
2,2-Dichloropropane	ug/L	ND	0.50	06/05/14 14:46	
2-Chlorotoluene	ug/L	ND	0.50	06/05/14 14:46	
4-Chlorotoluene	ug/L	ND	0.50	06/05/14 14:46	
Benzene	ug/L	ND	0.50	06/05/14 14:46	
Bromobenzene	ug/L	ND	0.50	06/05/14 14:46	
Bromochloromethane	ug/L	ND	0.50	06/05/14 14:46	
Bromodichloromethane	ug/L	ND	0.50	06/05/14 14:46	
Bromoform	ug/L	ND	0.50	06/05/14 14:46	
Bromomethane	ug/L	ND	5.0	06/05/14 14:46	
Carbon tetrachloride	ug/L	ND	0.50	06/05/14 14:46	
Chlorobenzene	ug/L	ND	0.50	06/05/14 14:46	
Chloroethane	ug/L	ND	1.0	06/05/14 14:46	
Chloroform	ug/L	ND	0.50	06/05/14 14:46	
Chloromethane	ug/L	ND	1.0	06/05/14 14:46	
cis-1,2-Dichloroethene	ug/L	ND	0.50	06/05/14 14:46	
cis-1,3-Dichloropropene	ug/L	ND	0.50	06/05/14 14:46	
Dibromochloromethane	ug/L	ND	0.50	06/05/14 14:46	
Dibromomethane	ug/L	ND	0.50	06/05/14 14:46	
Dichlorodifluoromethane	ug/L	ND	0.50	06/05/14 14:46	
Diisopropyl ether	ug/L	ND	0.50	06/05/14 14:46	
Ethylbenzene	ug/L	ND	0.50	06/05/14 14:46	
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## **REPORT OF LABORATORY ANALYSIS**

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## **QUALITY CONTROL DATA**

Matrix: Water

Project: 2014-093 Johnston 34182.1.2

# Pace Project No.: 92204085

METHOD BLANK:	1214892	

Associated Lab Samples: 92204085002

Blank Reporting Parameter Units Result Limit Analyzed Qualifiers Hexachloro-1,3-butadiene ug/L ND 2.0 06/05/14 14:46 Isopropylbenzene (Cumene) ND 0.50 06/05/14 14:46 ug/L ND 06/05/14 14:46 m&p-Xylene ug/L 1.0 Methyl-tert-butyl ether ug/L ND 0.50 06/05/14 14:46 ug/L Methylene Chloride ND 2.0 06/05/14 14:46 n-Butylbenzene ug/L ND 0.50 06/05/14 14:46 n-Propylbenzene ug/L ND 0.50 06/05/14 14:46 Naphthalene ug/L ND 2.0 06/05/14 14:46 o-Xylene ug/L ND 0.50 06/05/14 14:46 sec-Butylbenzene ND 0.50 06/05/14 14:46 ug/L Styrene ug/L ND 0.50 06/05/14 14:46 tert-Butylbenzene ug/L ND 0.50 06/05/14 14:46 Tetrachloroethene ug/L ND 0.50 06/05/14 14:46 Toluene ND 0.50 06/05/14 14:46 ug/L trans-1,2-Dichloroethene ND ug/L 0.50 06/05/14 14:46 trans-1,3-Dichloropropene ug/L ND 0.50 06/05/14 14:46 Trichloroethene ug/L ND 0.50 06/05/14 14:46 Trichlorofluoromethane ug/L ND 1.0 06/05/14 14:46 Vinyl chloride ug/L ND 1.0 06/05/14 14:46 1,2-Dichloroethane-d4 (S) % 94 70-130 06/05/14 14:46 % 97 4-Bromofluorobenzene (S) 70-130 06/05/14 14:46 Toluene-d8 (S) % 100 70-130 06/05/14 14:46

## LABORATORY CONTROL SAMPLE: 1214893

		Spike	LCS	LCS	% Rec	
Parameter	Units	Conc.	Result	% Rec	Limits	Qualifiers
1,1,1,2-Tetrachloroethane	ug/L	50	51.8	104	60-140	
1,1,1-Trichloroethane	ug/L	50	48.8	98	60-140	
1,1,2,2-Tetrachloroethane	ug/L	50	51.5	103	60-140	
1,1,2-Trichloroethane	ug/L	50	53.9	108	60-140	
1,1-Dichloroethane	ug/L	50	49.1	98	60-140	
1,1-Dichloroethene	ug/L	50	51.5	103	60-140	
1,1-Dichloropropene	ug/L	50	53.4	107	60-140	
1,2,3-Trichlorobenzene	ug/L	50	53.3	107	60-140	
1,2,3-Trichloropropane	ug/L	50	49.1	98	60-140	
1,2,4-Trichlorobenzene	ug/L	50	52.7	105	60-140	
1,2,4-Trimethylbenzene	ug/L	50	54.8	110	60-140	
1,2-Dibromo-3-chloropropane	ug/L	50	47.0	94	60-140	
1,2-Dibromoethane (EDB)	ug/L	50	53.0	106	60-140	
1,2-Dichlorobenzene	ug/L	50	52.1	104	60-140	
1,2-Dichloroethane	ug/L	50	46.0	92	60-140	
1,2-Dichloropropane	ug/L	50	51.0	102	60-140	
1,3,5-Trimethylbenzene	ug/L	50	53.9	108	60-140	
1.3-Dichlorobenzene	ua/l	50	52.4	105	60-140	

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## **REPORT OF LABORATORY ANALYSIS**

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Project: 2014-093 Johnston 34182.1.2

Pace Project No.: 92204085

#### LABORATORY CONTROL SAMPLE: 1214893

		Spike	LCS	LCS	% Rec	
Parameter	Units	Conc.	Result	% Rec	Limits	Qualifiers
1,3-Dichloropropane	ug/L		52.7	105	60-140	
1,4-Dichlorobenzene	ug/L	50	51.2	102	60-140	
2,2-Dichloropropane	ug/L	50	51.4	103	60-140	
2-Chlorotoluene	ug/L	50	51.9	104	60-140	
4-Chlorotoluene	ug/L	50	50.0	100	60-140	
Benzene	ug/L	50	56.4	113	60-140	
Bromobenzene	ug/L	50	53.2	106	60-140	
Bromochloromethane	ug/L	50	53.0	106	60-140	
Bromodichloromethane	ug/L	50	51.0	102	60-140	
Bromoform	ug/L	50	45.5	91	60-140	
Bromomethane	ug/L	50	63.7	127	60-140	
Carbon tetrachloride	ug/L	50	52.8	106	60-140	
Chlorobenzene	ug/L	50	52.1	104	60-140	
Chloroethane	ug/L	50	54.8	110	60-140	
Chloroform	ug/L	50	51.8	104	60-140	
Chloromethane	ug/L	50	56.9	114	60-140	
cis-1.2-Dichloroethene	ug/L	50	51.2	102	60-140	
cis-1.3-Dichloropropene	ug/L	50	55.0	110	60-140	
Dibromochloromethane	ug/L	50	52.4	105	60-140	
Dibromomethane	ug/L	50	52.1	104	60-140	
Dichlorodifluoromethane	ug/L	50	52.9	106	60-140	
Diisopropyl ether	ug/L	50	52.7	105	60-140	
Ethylbenzene	ug/L	50	52.2	104	60-140	
Hexachloro-1.3-butadiene	ug/L	50	50.3	101	60-140	
Isopropylbenzene (Cumene)	ug/L	50	54.4	109	60-140	
m&p-Xvlene	ug/L	100	108	108	60-140	
Methyl-tert-butyl ether	ug/L	50	51.6	103	60-140	
Methylene Chloride	ug/L	50	51.3	103	60-140	
n-Butvlbenzene	ug/L	50	56.1	112	60-140	
n-Propylbenzene	ug/L	50	53.7	107	60-140	
Naphthalene	ug/L	50	53.6	107	60-140	
o-Xvlene	ua/L	50	52.5	105	60-140	
sec-Butylbenzene	ua/L	50	53.6	107	60-140	
Styrene	ua/L	50	57.4	115	60-140	
tert-Butylbenzene	ug/L	50	52.7	105	60-140	
Tetrachloroethene	ug/L	50	52.6	105	60-140	
Toluene	ua/L	50	52.2	104	60-140	
trans-1.2-Dichloroethene	ua/L	50	51.7	103	60-140	
trans-1.3-Dichloropropene	ug/L	50	54.3	109	60-140	
Trichloroethene	ua/L	50	52.5	105	60-140	
Trichlorofluoromethane	ug/L	50	50.2	100	60-140	
Vinvl chloride	ug/L	50	64.0	128	60-140	
1.2-Dichloroethane-d4 (S)	~ <u>9</u> , <u>–</u> %	50	00	0	70-130	
4-Bromofluorobenzene (S)	%			100	70-130	
Toluene-d8 (S)	%			102	70-130	

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# **REPORT OF LABORATORY ANALYSIS**



Project: 2014-093 Johnston 34182.1.2

Pace Project No.: 92204085

MATRIX SPIKE & MATRIX SPIK	IATRIX SPIKE & MATRIX SPIKE DUPLICATE: 1215546 1215547										
			MS	MSD							
	922	04081001	Spike	Spike	MS	MSD	MS	MSD	% Rec		
Parameter	Units	Result	Conc.	Conc.	Result	Result	% Rec	% Rec	Limits	RPD	Qual
1,1,1,2-Tetrachloroethane	ug/L	ND	20	20	21.0	21.4	105	107	60-140	2	
1,1,1-Trichloroethane	ug/L	ND	20	20	21.3	21.3	107	106	60-140	0	
1,1,2,2-Tetrachloroethane	ug/L	ND	20	20	20.6	20.8	103	104	60-140	1	
1,1,2-Trichloroethane	ug/L	ND	20	20	21.5	21.6	107	108	60-140	1	
1,1-Dichloroethane	ug/L	ND	20	20	20.4	20.9	102	104	60-140	2	
1,1-Dichloroethene	ug/L	ND	20	20	22.4	23.3	112	116	60-140	4	
1,1-Dichloropropene	ug/L	ND	20	20	23.2	23.6	116	118	60-140	1	
1,2,3-Trichlorobenzene	ug/L	ND	20	20	19.4	20.4	97	102	60-140	5	
1,2,3-Trichloropropane	ug/L	ND	20	20	20.2	20.2	101	101	60-140	0	
1,2,4-Trichlorobenzene	ug/L	ND	20	20	20.1	20.5	101	103	60-140	2	
1,2,4-Trimethylbenzene	ug/L	ND	20	20	22.5	22.8	113	114	60-140	1	
1,2-Dibromo-3-chloropropane	ug/L	ND	20	20	17.7	18.5	89	92	60-140	4	
1,2-Dibromoethane (EDB)	ug/L	ND	20	20	20.9	21.7	104	108	60-140	4	
1,2-Dichlorobenzene	ug/L	ND	20	20	21.0	21.5	105	107	60-140	2	
1,2-Dichloroethane	ug/L	ND	20	20	19.0	19.5	95	98	60-140	3	
1,2-Dichloropropane	ug/L	ND	20	20	20.8	21.1	104	105	60-140	1	
1,3,5-Trimethylbenzene	ug/L	ND	20	20	22.2	22.8	111	114	60-140	3	
1,3-Dichlorobenzene	ug/L	ND	20	20	21.4	21.3	107	107	60-140	0	
1,3-Dichloropropane	ug/L	ND	20	20	21.1	21.3	105	106	60-140	1	
1,4-Dichlorobenzene	ug/L	ND	20	20	20.6	21.3	103	106	60-140	3	
2,2-Dichloropropane	ug/L	ND	20	20	21.9	22.3	109	111	60-140	2	
2-Chlorotoluene	ug/L	ND	20	20	21.2	21.9	106	109	60-140	3	
4-Chlorotoluene	ug/L	ND	20	20	20.4	21.1	102	106	60-140	3	
Benzene	ug/L	ND	20	20	23.1	23.3	115	117	60-140	1	
Bromobenzene	ug/L	ND	20	20	21.3	21.5	107	107	60-140	1	
Bromochloromethane	ug/L	ND	20	20	21.9	22.5	110	112	60-140	3	
Bromodichloromethane	ug/L	ND	20	20	20.4	20.1	102	101	60-140	1	
Bromoform	ug/L	ND	20	20	17.6	17.8	88	89	60-140	1	
Bromomethane	ug/L	ND	20	20	23.0	24.1	115	121	60-140	5	
Carbon tetrachloride	ug/L	ND	20	20	22.1	22.4	110	112	60-140	2	
Chlorobenzene	ug/L	ND	20	20	21.1	21.7	105	108	60-140	3	
Chloroethane	ug/L	ND	20	20	25.1	24.5	125	122	60-140	2	
Chloroform	ug/L	ND	20	20	21.4	22.1	107	110	60-140	3	
Chloromethane	ug/L	ND	20	20	21.6	25.9	108	130	60-140	18	
cis-1,2-Dichloroethene	ug/L	ND	20	20	21.9	21.7	109	108	60-140	1	
cis-1,3-Dichloropropene	ug/L	ND	20	20	20.8	21.4	104	107	60-140	3	
Dibromochloromethane	ug/L	ND	20	20	19.1	20.2	96	101	60-140	6	
Dibromomethane	ug/L	ND	20	20	20.3	20.9	101	105	60-140	3	
Dichlorodifluoromethane	ug/L	ND	20	20	24.0	24.5	120	122	60-140	2	
Diisopropyl ether	ug/L	ND	20	20	21.4	22.0	107	110	60-140	3	
Ethylbenzene	ug/L	ND	20	20	21.9	22.2	110	111	60-140	1	
Hexachloro-1,3-butadiene	ug/L	ND	20	20	20.2	21.0	101	105	60-140	4	
Isopropylbenzene (Cumene)	ug/L	ND	20	20	22.6	23.3	113	116	60-140	3	
m&p-Xylene	ug/L	ND	40	40	45.0	45.4	112	114	60-140	1	
Methyl-tert-butyl ether	ug/L	ND	20	20	20.5	21.7	102	108	60-140	6	
Methylene Chloride	ug/L	ND	20	20	18.9	19.4	94	97	60-140	3	

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# **REPORT OF LABORATORY ANALYSIS**



Project: 2014-093 Johnston 34182.1.2

Pace Project No.: 92204085

MATRIX SPIKE & MATRIX SPIKE DUPLICATE: 1215546 121554					1215547						
			MS	MSD							
	922	204081001	Spike	Spike	MS	MSD	MS	MSD	% Rec		
Parameter	Units	Result	Conc.	Conc.	Result	Result	% Rec	% Rec	Limits	RPD	Qual
n-Butylbenzene	ug/L	ND	20	20	22.9	23.3	114	117	60-140	2	
n-Propylbenzene	ug/L	ND	20	20	22.4	22.9	112	114	60-140	2	
Naphthalene	ug/L	ND	20	20	20.0	20.7	100	103	60-140	3	
o-Xylene	ug/L	ND	20	20	21.5	21.9	108	109	60-140	2	
sec-Butylbenzene	ug/L	ND	20	20	22.5	22.8	113	114	60-140	1	
Styrene	ug/L	ND	20	20	22.9	23.7	115	118	60-140	3	
tert-Butylbenzene	ug/L	ND	20	20	22.5	22.8	112	114	60-140	1	
Tetrachloroethene	ug/L	ND	20	20	22.4	22.4	112	112	60-140	0	
Toluene	ug/L	ND	20	20	21.7	22.3	109	111	60-140	3	
trans-1,2-Dichloroethene	ug/L	ND	20	20	22.0	23.0	110	115	60-140	5	
trans-1,3-Dichloropropene	ug/L	ND	20	20	20.9	21.0	105	105	60-140	1	
Trichloroethene	ug/L	ND	20	20	22.4	22.3	112	111	60-140	1	
Trichlorofluoromethane	ug/L	ND	20	20	23.4	23.7	117	119	60-140	1	
Vinyl chloride	ug/L	ND	20	20	27.8	29.5	139	147	60-140	6 M	J
1,2-Dichloroethane-d4 (S)	%						91	92	70-130		
4-Bromofluorobenzene (S)	%						100	99	70-130		
Toluene-d8 (S)	%						100	100	70-130		

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# **REPORT OF LABORATORY ANALYSIS**



Project: 2014-093 Johnston 34182.1.2

EPA 8260

Pace Project No.: 92204085

QC Batch Method:

QC Batch: MSV/27125

Analysis Method:

Analysis Description:

EPA 8260 8260 MSV 5035A Volatile Organics

Associated Lab Samples: 92204085001

METHOD BLANK: 1216668 Matrix: Solid Associated Lab Samples: 92204085001 Blank Reporting Result Limit Qualifiers Parameter Units Analyzed 1,1,1,2-Tetrachloroethane ND 06/09/14 12:28 ug/kg 5.0 1,1,1-Trichloroethane ND 50 06/09/14 12:28 ug/kg 1,1,2,2-Tetrachloroethane ug/kg ND 5.0 06/09/14 12:28 ND 1,1,2-Trichloroethane ug/kg 5.0 06/09/14 12:28 1,1-Dichloroethane ug/kg ND 5.0 06/09/14 12:28 1,1-Dichloroethene ug/kg ND 5.0 06/09/14 12:28 1,1-Dichloropropene ug/kg ND 5.0 06/09/14 12:28 ND 5.0 06/09/14 12:28 1,2,3-Trichlorobenzene ug/kg 1,2,3-Trichloropropane ug/kg ND 5.0 06/09/14 12:28 1.2.4-Trichlorobenzene ug/kg ND 5.0 06/09/14 12:28 1,2,4-Trimethylbenzene ND 5.0 06/09/14 12:28 ug/kg ND 5.0 06/09/14 12:28 1,2-Dibromo-3-chloropropane ug/kg 1.2-Dibromoethane (EDB) ND 5.0 06/09/14 12:28 ug/kg 1,2-Dichlorobenzene ND 5.0 06/09/14 12:28 ug/kg ND 1,2-Dichloroethane ug/kg 5.0 06/09/14 12:28 1,2-Dichloropropane ug/kg ND 5.0 06/09/14 12:28 1,3,5-Trimethylbenzene ug/kg ND 5.0 06/09/14 12:28 1,3-Dichlorobenzene ug/kg ND 5.0 06/09/14 12:28 ND 5.0 06/09/14 12:28 1,3-Dichloropropane ug/kg 06/09/14 12:28 1,4-Dichlorobenzene ug/kg ND 5.0 ug/kg 2,2-Dichloropropane ND 5.0 06/09/14 12:28 2-Butanone (MEK) ND 100 06/09/14 12:28 ug/kg 2-Chlorotoluene ND 5.0 06/09/14 12:28 ug/kg 2-Hexanone ND 50.0 06/09/14 12:28 ug/kg 4-Chlorotoluene ND 5.0 06/09/14 12:28 ug/kg 4-Methyl-2-pentanone (MIBK) ug/kg ND 50.0 06/09/14 12:28 Acetone ug/kg ND 100 06/09/14 12:28 Benzene ug/kg ND 5.0 06/09/14 12:28 Bromobenzene ug/kg ND 5.0 06/09/14 12:28 ND 06/09/14 12:28 Bromochloromethane ug/kg 5.0 Bromodichloromethane ug/kg ND 5.0 06/09/14 12:28 Bromoform ug/kg ND 5.0 06/09/14 12:28 Bromomethane ug/kg ND 10.0 06/09/14 12:28 Carbon tetrachloride ug/kg ND 50 06/09/14 12:28 5.0 Chlorobenzene ug/kg ND 06/09/14 12:28 Chloroethane ug/kg ND 10.0 06/09/14 12:28 Chloroform ND ug/kg 5.0 06/09/14 12:28 ND Chloromethane ug/kg 10.0 06/09/14 12:28 cis-1,2-Dichloroethene ug/kg ND 5.0 06/09/14 12:28 cis-1,3-Dichloropropene ND 5.0 06/09/14 12:28 ug/kg Dibromochloromethane ug/kg ND 5.0 06/09/14 12:28

Results presented on this page are in the units indicated by the "Units" column except where an alternate unit is presented to the right of the result.

# **REPORT OF LABORATORY ANALYSIS**



Project: 2014-093 Johnston 34182.1.2

# Pace Project No.: 92204085

# METHOD BLANK: 1216668

Associated Lab Samples: 92204085001

Matrix: Solid

		Blank	Reporting		
Parameter	Units	Result	Limit	Analyzed	Qualifiers
Dibromomethane	ug/kg	ND	5.0	06/09/14 12:28	
Dichlorodifluoromethane	ug/kg	ND	10.0	06/09/14 12:28	
Diisopropyl ether	ug/kg	ND	5.0	06/09/14 12:28	
Ethylbenzene	ug/kg	ND	5.0	06/09/14 12:28	
Hexachloro-1,3-butadiene	ug/kg	ND	5.0	06/09/14 12:28	
Isopropylbenzene (Cumene)	ug/kg	ND	5.0	06/09/14 12:28	
m&p-Xylene	ug/kg	ND	10.0	06/09/14 12:28	
Methyl-tert-butyl ether	ug/kg	ND	5.0	06/09/14 12:28	
Methylene Chloride	ug/kg	ND	20.0	06/09/14 12:28	
n-Butylbenzene	ug/kg	ND	5.0	06/09/14 12:28	
n-Propylbenzene	ug/kg	ND	5.0	06/09/14 12:28	
Naphthalene	ug/kg	ND	5.0	06/09/14 12:28	
o-Xylene	ug/kg	ND	5.0	06/09/14 12:28	
p-Isopropyltoluene	ug/kg	ND	5.0	06/09/14 12:28	
sec-Butylbenzene	ug/kg	ND	5.0	06/09/14 12:28	
Styrene	ug/kg	ND	5.0	06/09/14 12:28	
tert-Butylbenzene	ug/kg	ND	5.0	06/09/14 12:28	
Tetrachloroethene	ug/kg	ND	5.0	06/09/14 12:28	
Toluene	ug/kg	ND	5.0	06/09/14 12:28	
trans-1,2-Dichloroethene	ug/kg	ND	5.0	06/09/14 12:28	
trans-1,3-Dichloropropene	ug/kg	ND	5.0	06/09/14 12:28	
Trichloroethene	ug/kg	ND	5.0	06/09/14 12:28	
Trichlorofluoromethane	ug/kg	ND	5.0	06/09/14 12:28	
Vinyl acetate	ug/kg	ND	50.0	06/09/14 12:28	
Vinyl chloride	ug/kg	ND	10.0	06/09/14 12:28	
Xylene (Total)	ug/kg	ND	10.0	06/09/14 12:28	
1,2-Dichloroethane-d4 (S)	%	98	70-132	06/09/14 12:28	
4-Bromofluorobenzene (S)	%	105	70-130	06/09/14 12:28	
Toluene-d8 (S)	%	106	70-130	06/09/14 12:28	

#### LABORATORY CONTROL SAMPLE: 1216669

Parameter	Units	Spike Conc	LCS Result	LCS % Rec	% Rec	Qualifiers
						Qualment
1,1,1,2-Tetrachloroethane	ug/kg	49.6	53.3	107	74-137	
1,1,1-Trichloroethane	ug/kg	49.6	49.6	100	67-140	
1,1,2,2-Tetrachloroethane	ug/kg	49.6	52.9	107	72-141	
1,1,2-Trichloroethane	ug/kg	49.6	53.9	109	78-138	
1,1-Dichloroethane	ug/kg	49.6	48.4	97	69-134	
1,1-Dichloroethene	ug/kg	49.6	51.1	103	67-138	
1,1-Dichloropropene	ug/kg	49.6	56.3	113	69-139	
1,2,3-Trichlorobenzene	ug/kg	49.6	55.2	111	70-146	
1,2,3-Trichloropropane	ug/kg	49.6	51.6	104	69-144	
1,2,4-Trichlorobenzene	ug/kg	49.6	56.4	114	68-148	
1,2,4-Trimethylbenzene	ug/kg	49.6	57.9	117	74-137	

Results presented on this page are in the units indicated by the "Units" column except where an alternate unit is presented to the right of the result.

# **REPORT OF LABORATORY ANALYSIS**



Project: 2014-093 Johnston 34182.1.2

Pace Project No.: 92204085

#### LABORATORY CONTROL SAMPLE: 1216669

		Spike	LCS	LCS	% Rec	
Parameter	Units	Conc.	Result	% Rec	Limits	Qualifiers
1,2-Dibromo-3-chloropropane	ug/kg	49.6	54.8		65-140	
1,2-Dibromoethane (EDB)	ug/kg	49.6	53.9	109	77-135	
1,2-Dichlorobenzene	ug/kg	49.6	55.0	111	77-141	
1,2-Dichloroethane	ug/kg	49.6	51.9	105	65-137	
1,2-Dichloropropane	ug/kg	49.6	53.5	108	72-136	
1,3,5-Trimethylbenzene	ug/kg	49.6	56.2	113	76-133	
1,3-Dichlorobenzene	ug/kg	49.6	55.1	111	74-138	
1,3-Dichloropropane	ug/kg	49.6	55.1	111	71-139	
1.4-Dichlorobenzene	ua/ka	49.6	55.7	112	76-138	
2,2-Dichloropropane	ug/kg	49.6	48.5	98	68-137	
2-Butanone (MEK)	ug/kg	99.2	92.7J	93	58-147	
2-Chlorotoluene	ug/kg	49.6	55.5	112	73-139	
2-Hexanone	ua/ka	99.2	106	107	62-145	
4-Chlorotoluene	ug/kg	49.6	55.4	112	76-141	
4-Methyl-2-pentanone (MIBK)	ug/kg	99.2	112	113	64-149	
Acetone	ua/ka	99.2	97.2J	98	53-153	
Benzene	ua/ka	49.6	53.5	108	73-135	
Bromobenzene	ua/ka	49.6	56.7	114	75-133	
Bromochloromethane	ua/ka	49.6	51.5	104	73-134	
Bromodichloromethane	ua/ka	49.6	58.7	118	71-135	
Bromoform	ua/ka	49.6	56.3	114	66-141	
Bromomethane	ua/ka	49.6	54.8	110	53-160	
Carbon tetrachloride	ua/ka	49.6	50.6	102	60-145	
Chlorobenzene	ua/ka	49.6	52.0	105	78-130	
Chloroethane	ua/ka	49.6	48.7	98	64-149	
Chloroform	ua/ka	49.6	58.2	117	70-134	
Chloromethane	ua/ka	49.6	40.7	82	52-150	
cis-1.2-Dichloroethene	ua/ka	49.6	54.0	109	70-133	
cis-1,3-Dichloropropene	ua/ka	49.6	54.1	109	68-134	
Dibromochloromethane	ug/kg	49.6	60.7	122	71-138	
Dibromomethane	ua/ka	49.6	51.8	105	74-130	
Dichlorodifluoromethane	ua/ka	49.6	39.2	79	40-160	
Diisopropyl ether	ua/ka	49.6	56.7	114	69-141	
Ethylbenzene	ua/ka	49.6	53.8	108	75-133	
Hexachloro-1.3-butadiene	ua/ka	49.6	52.7	106	68-143	
Isopropylbenzene (Cumene)	ua/ka	49.6	54.1	109	76-143	
m&p-Xvlene	ua/ka	99.2	106	107	75-136	
Methyl-tert-butyl ether	ua/ka	49.6	52.3	105	68-144	
Methylene Chloride	ua/ka	49.6	50.2	101	45-154	
n-Butylbenzene	ua/ka	49.6	57.5	116	72-137	
n-Propylbenzene	ua/ka	49.6	56.5	114	76-136	
Naphthalene	ua/ka	49.6	56.3	114	68-151	
o-Xvlene	ug/ka	49.6	52.7	106	76-141	
p-Isopropyltoluene	ug/kg	49.6	55.5	112	76-140	
sec-Butylbenzene	ug/kg	49.6	55.5	112	79-139	
Styrene	ug/kg	49.6	55 3	112	79-137	
tert-Butylbenzene	ug/kg	40.0 40 A	55 F	112	74-143	
tort Buty bonzono	49/119	-0.0	55.0	114	1-1-5	

Results presented on this page are in the units indicated by the "Units" column except where an alternate unit is presented to the right of the result.

# **REPORT OF LABORATORY ANALYSIS**



Project: 2014-093 Johnston 34182.1.2

Pace Project No.: 92204085

#### LABORATORY CONTROL SAMPLE: 1216669

		Spike	LCS	LCS	% Rec	
Parameter	Units	Conc.	Result	% Rec	Limits	Qualifiers
Tetrachloroethene	ug/kg	49.6	52.0	105	71-138	
Toluene	ug/kg	49.6	52.4	106	74-131	
trans-1,2-Dichloroethene	ug/kg	49.6	52.1	105	67-135	
trans-1,3-Dichloropropene	ug/kg	49.6	54.2	109	65-146	
Trichloroethene	ug/kg	49.6	52.6	106	67-135	
Trichlorofluoromethane	ug/kg	49.6	45.6	92	59-144	
Vinyl acetate	ug/kg	99.2	103	103	40-160 F	=3
Vinyl chloride	ug/kg	49.6	43.1	87	56-141	
Xylene (Total)	ug/kg	149	159	107	76-137	
1,2-Dichloroethane-d4 (S)	%			94	70-132	
4-Bromofluorobenzene (S)	%			94	70-130	
Toluene-d8 (S)	%			97	70-130	

MATRIX SPIKE SAMPLE:	1217247						
		92204084001	Spike	MS	MS	% Rec	
Parameter	Units	Result	Conc.	Result	% Rec	Limits	Qualifiers
1,1-Dichloroethene	ug/kg	ND	45.1	46.2	103	49-180	
Benzene	ug/kg	ND	45.1	45.9	102	50-166	
Chlorobenzene	ug/kg	ND	45.1	42.6	95	43-169	
Toluene	ug/kg	ND	45.1	46.1	102	52-163	
Trichloroethene	ug/kg	ND	45.1	45.7	101	49-167	
1,2-Dichloroethane-d4 (S)	%				118	70-132	
4-Bromofluorobenzene (S)	%				98	70-130	
Toluene-d8 (S)	%				101	70-130	

## SAMPLE DUPLICATE: 1217468

		92204085001	Dup		
Parameter	Units	Result	Result	RPD	Qualifiers
1,1,1,2-Tetrachloroethane	ug/kg	ND	ND		
1,1,1-Trichloroethane	ug/kg	ND	ND		
1,1,2,2-Tetrachloroethane	ug/kg	ND	ND		
1,1,2-Trichloroethane	ug/kg	ND	ND		
1,1-Dichloroethane	ug/kg	ND	ND		
1,1-Dichloroethene	ug/kg	ND	ND		
1,1-Dichloropropene	ug/kg	ND	ND		
1,2,3-Trichlorobenzene	ug/kg	ND	ND		
1,2,3-Trichloropropane	ug/kg	ND	ND		
1,2,4-Trichlorobenzene	ug/kg	ND	ND		
1,2,4-Trimethylbenzene	ug/kg	ND	ND		
1,2-Dibromo-3-chloropropane	ug/kg	ND	ND		
1,2-Dibromoethane (EDB)	ug/kg	ND	ND		
1,2-Dichlorobenzene	ug/kg	ND	ND		
1,2-Dichloroethane	ug/kg	ND	ND		
1,2-Dichloropropane	ug/kg	ND	ND		

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# **REPORT OF LABORATORY ANALYSIS**



Project: 2014-093 Johnston 34182.1.2

Pace Project No.: 92204085

# SAMPLE DUPLICATE: 1217468

		92204085001	Dup		
Parameter	Units	Result	Result	RPD	Qualifiers
1,3,5-Trimethylbenzene	ug/kg	ND	ND		_
1,3-Dichlorobenzene	ug/kg	ND	ND		
1,3-Dichloropropane	ug/kg	ND	ND		
1,4-Dichlorobenzene	ug/kg	ND	ND		
2,2-Dichloropropane	ug/kg	ND	ND		
2-Butanone (MEK)	ug/kg	ND	ND		
2-Chlorotoluene	ug/kg	ND	ND		
2-Hexanone	ug/kg	ND	ND		
4-Chlorotoluene	ug/kg	ND	ND		
4-Methyl-2-pentanone (MIBK)	ug/kg	ND	ND		
Acetone	ug/kg	ND	19.5J		
Benzene	ug/kg	ND	ND		
Bromobenzene	ug/kg	ND	ND		
Bromochloromethane	ua/ka	ND	ND		
Bromodichloromethane	ug/kg	ND	ND		
Bromoform	ug/kg	ND	ND		
Bromomethane	ua/ka	ND	ND		
Carbon tetrachloride	ua/ka	ND	ND		
Chlorobenzene	ua/ka	ND	ND		
Chloroethane	ug/kg	ND	ND		
Chloroform	ug/kg	ND	ND		
Chloromethane	ug/kg	ND	ND		
cis-1,2-Dichloroethene	ug/kg	ND	ND		
cis-1.3-Dichloropropene	ua/ka	ND	ND		
Dibromochloromethane	ug/kg	ND	ND		
Dibromomethane	ug/kg	ND	ND		
Dichlorodifluoromethane	ug/kg	ND	ND		
Diisopropyl ether	ug/kg	ND	ND		
Ethylbenzene	ug/kg	ND	ND		
Hexachloro-1,3-butadiene	ug/kg	ND	ND		
Isopropylbenzene (Cumene)	ug/kg	ND	ND		
m&p-Xylene	ug/kg	ND	ND		
Methyl-tert-butyl ether	ug/kg	ND	ND		
Methylene Chloride	ug/kg	ND	ND		
n-Butylbenzene	ug/kg	ND	ND		
n-Propylbenzene	ug/kg	ND	ND		
Naphthalene	ug/kg	ND	ND		
o-Xylene	ug/kg	ND	ND		
p-Isopropyltoluene	ug/kg	ND	ND		
sec-Butylbenzene	ug/kg	ND	ND		
Styrene	ug/kg	ND	ND		
tert-Butylbenzene	ug/kg	ND	ND		
Tetrachloroethene	ug/kg	ND	ND		
Toluene	ug/kg	ND	ND		
trans-1,2-Dichloroethene	ug/kg	ND	ND		
trans-1,3-Dichloropropene	ug/kg	ND	ND		
Trichloroethene	ug/kg	ND	ND		

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# **REPORT OF LABORATORY ANALYSIS**



Dup

Project: 2014-093 Johnston 34182.1.2

Pace Project No.: 92204085

SAMPLE DUPLICATE:	1217468	
Paramotor	Linite	

Parameter	Units	Result	Result	RPD	Qualifiers
Trichlorofluoromethane	ug/kg	ND	ND		
Vinyl acetate	ug/kg	ND	ND		
Vinyl chloride	ug/kg	ND	ND		
Xylene (Total)	ug/kg	ND	ND		
1,2-Dichloroethane-d4 (S)	%	113	112	12	
4-Bromofluorobenzene (S)	%	102	104	9	
Toluene-d8 (S)	%	101	102	10	

92204085001

Results presented on this page are in the units indicated by the "Units" column except where an alternate unit is presented to the right of the result.



Project: 2014-093 Johnston 34182.1.2

Pace Project No.: 92204085

Face Floject No 92204065					
QC Batch: OEXT/2818	9	Analysis Met	hod: EF	PA 625	
QC Batch Method: EPA 625		Analysis Des	cription: 62	25 MSS	
Associated Lab Samples: 9220	4085002				
METHOD BLANK: 1218104		Matrix:	Water		
Associated Lab Samples: 9220	4085002				
		Blank	Reporting		
Parameter	Units	Result	Limit	Analyzed	Qualifiers
1,2,4-Trichlorobenzene	ug/L	ND	5.0	06/12/14 08:35	
2,4,6-Trichlorophenol	ug/L	ND	10.0	06/12/14 08:35	
2,4-Dichlorophenol	ug/L	ND	5.0	06/12/14 08:35	
2,4-Dimethylphenol	ug/L	ND	10.0	06/12/14 08:35	
2,4-Dinitrophenol	ug/L	ND	50.0	06/12/14 08:35	
2,4-Dinitrotoluene	ug/L	ND	5.0	06/12/14 08:35	
2,6-Dinitrotoluene	ug/L	ND	5.0	06/12/14 08:35	
2-Chloronaphthalene	ua/L	ND	5.0	06/12/14 08:35	

		Blank Reporting			
Parameter	Units	Result	Limit	Analyzed	Qualifiers
1,2,4-Trichlorobenzene	ug/L	ND	5.0	06/12/14 08:35	
2,4,6-Trichlorophenol	ug/L	ND	10.0	06/12/14 08:35	
2,4-Dichlorophenol	ug/L	ND	5.0	06/12/14 08:35	
2,4-Dimethylphenol	ug/L	ND	10.0	06/12/14 08:35	
2,4-Dinitrophenol	ug/L	ND	50.0	06/12/14 08:35	
2,4-Dinitrotoluene	ug/L	ND	5.0	06/12/14 08:35	
2,6-Dinitrotoluene	ug/L	ND	5.0	06/12/14 08:35	
2-Chloronaphthalene	ug/L	ND	5.0	06/12/14 08:35	
2-Chlorophenol	ug/L	ND	5.0	06/12/14 08:35	
2-Nitrophenol	ug/L	ND	5.0	06/12/14 08:35	
3,3'-Dichlorobenzidine	ug/L	ND	25.0	06/12/14 08:35	
4,6-Dinitro-2-methylphenol	ug/L	ND	20.0	06/12/14 08:35	
4-Bromophenylphenyl ether	ug/L	ND	5.0	06/12/14 08:35	
4-Chloro-3-methylphenol	ug/L	ND	5.0	06/12/14 08:35	
4-Chlorophenylphenyl ether	ug/L	ND	5.0	06/12/14 08:35	
4-Nitrophenol	ug/L	ND	50.0	06/12/14 08:35	
Acenaphthene	ug/L	ND	5.0	06/12/14 08:35	
Acenaphthylene	ug/L	ND	5.0	06/12/14 08:35	
Anthracene	ug/L	ND	5.0	06/12/14 08:35	
Benzo(a)anthracene	ug/L	ND	5.0	06/12/14 08:35	
Benzo(a)pyrene	ug/L	ND	5.0	06/12/14 08:35	
Benzo(b)fluoranthene	ug/L	ND	5.0	06/12/14 08:35	
Benzo(g,h,i)perylene	ug/L	ND	5.0	06/12/14 08:35	
Benzo(k)fluoranthene	ug/L	ND	5.0	06/12/14 08:35	
bis(2-Chloroethoxy)methane	ug/L	ND	10.0	06/12/14 08:35	
bis(2-Chloroethyl) ether	ug/L	ND	5.0	06/12/14 08:35	
bis(2-Chloroisopropyl) ether	ug/L	ND	5.0	06/12/14 08:35	
bis(2-Ethylhexyl)phthalate	ug/L	ND	5.0	06/12/14 08:35	
Butylbenzylphthalate	ug/L	ND	5.0	06/12/14 08:35	
Chrysene	ug/L	ND	5.0	06/12/14 08:35	
Di-n-butylphthalate	ug/L	ND	5.0	06/12/14 08:35	
Di-n-octylphthalate	ug/L	ND	5.0	06/12/14 08:35	
Dibenz(a,h)anthracene	ug/L	ND	5.0	06/12/14 08:35	
Diethylphthalate	ug/L	ND	5.0	06/12/14 08:35	
Dimethylphthalate	ug/L	ND	5.0	06/12/14 08:35	
Fluoranthene	ug/L	ND	5.0	06/12/14 08:35	
Fluorene	ug/L	ND	5.0	06/12/14 08:35	
Hexachloro-1,3-butadiene	ug/L	ND	5.0	06/12/14 08:35	

Results presented on this page are in the units indicated by the "Units" column except where an alternate unit is presented to the right of the result.

5.0 06/12/14 08:35

5.0 06/12/14 08:35

10.0 06/12/14 08:35

ND

ND

ND

# **REPORT OF LABORATORY ANALYSIS**

Hexachlorobenzene

Hexachloroethane

Hexachlorocyclopentadiene

ug/L

ug/L

ug/L



 Project:
 2014-093 Johnston 34182.1.2

 Pace Project No.:
 92204085

# METHOD BLANK: 1218104

Associated Lab Samples: 92204085002

Matrix: Water

Parameter	Units	Blank Result	Reporting Limit	Analyzed	Qualifiers
Indeno(1,2,3-cd)pyrene	ug/L		5.0	06/12/14 08:35	
Isophorone	ug/L	ND	10.0	06/12/14 08:35	
N-Nitroso-di-n-propylamine	ug/L	ND	5.0	06/12/14 08:35	
N-Nitrosodimethylamine	ug/L	ND	5.0	06/12/14 08:35	
N-Nitrosodiphenylamine	ug/L	ND	10.0	06/12/14 08:35	
Naphthalene	ug/L	ND	5.0	06/12/14 08:35	
Nitrobenzene	ug/L	ND	5.0	06/12/14 08:35	
Pentachlorophenol	ug/L	ND	10.0	06/12/14 08:35	
Phenanthrene	ug/L	ND	5.0	06/12/14 08:35	
Phenol	ug/L	ND	5.0	06/12/14 08:35	
Pyrene	ug/L	ND	5.0	06/12/14 08:35	
2,4,6-Tribromophenol (S)	%	110	10-137	06/12/14 08:35	
2-Fluorobiphenyl (S)	%	76	15-120	06/12/14 08:35	
2-Fluorophenol (S)	%	40	10-120	06/12/14 08:35	
Nitrobenzene-d5 (S)	%	97	10-120	06/12/14 08:35	
Phenol-d6 (S)	%	29	10-120	06/12/14 08:35	
Terphenyl-d14 (S)	%	80	11-131	06/12/14 08:35	

#### LABORATORY CONTROL SAMPLE: 1218105

		Spike	LCS	LCS	% Rec	
Parameter	Units	Conc.	Result	% Rec	Limits	Qualifiers
1,2,4-Trichlorobenzene	ug/L		40.5	81	44-142	
2,4,6-Trichlorophenol	ug/L	50	44.8	90	37-144	
2,4-Dichlorophenol	ug/L	50	41.1	82	1-191	
2,4-Dimethylphenol	ug/L	50	36.9	74	32-119	
2,4-Dinitrophenol	ug/L	250	114	46	1-181	
2,4-Dinitrotoluene	ug/L	50	66.7	133	39-139	
2,6-Dinitrotoluene	ug/L	50	59.4	119	50-158	
2-Chloronaphthalene	ug/L	50	39.4	79	60-118	
2-Chlorophenol	ug/L	50	34.2	68	23-134	
2-Nitrophenol	ug/L	50	46.2	92	29-182	
3,3'-Dichlorobenzidine	ug/L	100	79.3	79	1-262	
4,6-Dinitro-2-methylphenol	ug/L	100	46.5	46	1-181	
4-Bromophenylphenyl ether	ug/L	50	35.7	71	53-127	
4-Chloro-3-methylphenol	ug/L	100	95.7	96	22-147	
4-Chlorophenylphenyl ether	ug/L	50	45.1	90	25-158	
4-Nitrophenol	ug/L	250	207	83	1-132	
Acenaphthene	ug/L	50	40.1	80	47-145	
Acenaphthylene	ug/L	50	39.6	79	33-145	
Anthracene	ug/L	50	44.2	88	1-166	
Benzo(a)anthracene	ug/L	50	42.9	86	33-143	
Benzo(a)pyrene	ug/L	50	44.2	88	17-163	
Benzo(b)fluoranthene	ug/L	50	39.6	79	24-159	
Benzo(g,h,i)perylene	ug/L	50	32.1	64	1-219	

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# **REPORT OF LABORATORY ANALYSIS**



Project: 2014-093 Johnston 34182.1.2

Pace Project No.: 92204085

#### LABORATORY CONTROL SAMPLE: 1218105

		Spike	LCS	LCS	% Rec	
Parameter	Units	Conc.	Result	% Rec	Limits	Qualifiers
Benzo(k)fluoranthene	ug/L	50	45.3	91	11-162	
bis(2-Chloroethoxy)methane	ug/L	50	43.0	86	33-184	
bis(2-Chloroethyl) ether	ug/L	50	40.8	82	12-158	
bis(2-Chloroisopropyl) ether	ug/L	50	42.2	84	36-166	
bis(2-Ethylhexyl)phthalate	ug/L	50	41.8	84	8-158	
Butylbenzylphthalate	ug/L	50	46.2	92	1-152	
Chrysene	ug/L	50	44.3	89	17-168	
Di-n-butylphthalate	ug/L	50	39.5	79	1-118	
Di-n-octylphthalate	ug/L	50	49.4	99	4-146	
Dibenz(a,h)anthracene	ug/L	50	34.4	69	1-227	
Diethylphthalate	ug/L	50	45.1	90	1-114	
Dimethylphthalate	ug/L	50	44.4	89	1-112	
Fluoranthene	ug/L	50	53.2	106	26-137	
Fluorene	ug/L	50	47.7	95	59-121	
Hexachloro-1,3-butadiene	ug/L	50	44.3	89	24-116	
Hexachlorobenzene	ug/L	50	35.1	70	1-152	
Hexachlorocyclopentadiene	ug/L	50	22.4	45	25-150	
Hexachloroethane	ug/L	50	41.8	84	40-113	
Indeno(1,2,3-cd)pyrene	ug/L	50	27.7	55	1-171	
Isophorone	ug/L	50	52.1	104	21-196	
N-Nitroso-di-n-propylamine	ug/L	50	38.6	77	1-230	
N-Nitrosodimethylamine	ug/L	50	25.3	51	25-150	
N-Nitrosodiphenylamine	ug/L	50	31.3	63	25-150	
Naphthalene	ug/L	50	37.3	75	21-133	
Nitrobenzene	ug/L	50	56.0	112	35-180	
Pentachlorophenol	ug/L	100	85.3	85	14-176	
Phenanthrene	ug/L	50	42.3	85	54-120	
Phenol	ug/L	50	19.2	38	5-112	
Pyrene	ug/L	50	38.5	77	52-115	
2,4,6-Tribromophenol (S)	%			124	10-137	
2-Fluorobiphenyl (S)	%			79	15-120	
2-Fluorophenol (S)	%			45	10-120	
Nitrobenzene-d5 (S)	%			103	10-120	
Phenol-d6 (S)	%			31	10-120	
Terphenyl-d14 (S)	%			80	11-131	

MATRIX SPIKE & MATRIX SPIKE	PIKE DUPLICAT	E: 12181	06		1218107						
			MS	MSD							
	922	04282003	Spike	Spike	MS	MSD	MS	MSD	% Rec		
Parameter	Units	Result	Conc.	Conc.	Result	Result	% Rec	% Rec	Limits	RPD	Qual
1,2,4-Trichlorobenzene	ug/L	ND	100	100	74.8	76.4	75	76	44-142	2	
2,4,6-Trichlorophenol	ug/L	ND	100	100	84.1	85.0	84	85	37-144	1	
2,4-Dichlorophenol	ug/L	ND	100	100	81.6	83.1	82	83	1-191	2	
2,4-Dimethylphenol	ug/L	ND	100	100	72.1	74.6	72	75	32-119	3	
2,4-Dinitrophenol	ug/L	ND	500	500	266	182	53	36	1-181	38 R	1

Results presented on this page are in the units indicated by the "Units" column except where an alternate unit is presented to the right of the result.

# **REPORT OF LABORATORY ANALYSIS**



Project: 2014-093 Johnston 34182.1.2

Pace Project No.: 92204085

MATRIX SPIKE & MATRIX SPIKE DUPLICATE: 1218106 1218107											
M		MS	MSD								
	9220	04282003	Spike	Spike	MS	MSD	MS	MSD	% Rec		
Parameter	Units	Result	Conc.	Conc.	Result	Result	% Rec	% Rec	Limits	RPD	Qual
2,4-Dinitrotoluene	ug/L	ND	100	100	122	116	122	116	39-139	5	
2,6-Dinitrotoluene	ug/L	ND	100	100	113	110	113	110	50-158	2	
2-Chloronaphthalene	ug/L	ND	100	100	74.2	72.3	74	72	60-118	3	
2-Chlorophenol	ug/L	ND	100	100	71.6	71.2	72	71	23-134	1	
2-Nitrophenol	ug/L	ND	100	100	86.2	89.9	86	90	29-182	4	
3,3'-Dichlorobenzidine	ug/L	ND	200	200	156	160	78	80	1-262	2	
4,6-Dinitro-2-methylphenol	ug/L	ND	200	200	108	105	54	53	1-181	3	
4-Bromophenylphenyl ether	ug/L	ND	100	100	71.3	69.0	71	69	53-127	3	
4-Chloro-3-methylphenol	ug/L	ND	200	200	199	216	99	108	22-147	8	
4-Chlorophenylphenyl ether	ug/L	ND	100	100	87.8	85.9	88	86	25-158	2	
4-Nitrophenol	ug/L	ND	500	500	564	539	113	108	1-132	4	
Acenaphthene	ug/L	ND	100	100	77.9	76.5	78	76	47-145	2	
Acenaphthylene	ug/L	ND	100	100	76.7	75.5	77	76	33-145	2	
Anthracene	ug/L	ND	100	100	86.2	84.6	86	85	1-166	2	
Benzo(a)anthracene	ug/L	ND	100	100	82.4	81.8	82	82	33-143	1	
Benzo(a)pyrene	ug/L	ND	100	100	86.2	87.0	86	87	17-163	1	
Benzo(b)fluoranthene	ug/L	ND	100	100	80.5	81.9	81	82	24-159	2	
Benzo(g.h.i)pervlene	ua/L	ND	100	100	61.0	62.4	61	62	1-219	2	
Benzo(k)fluoranthene	ug/L	ND	100	100	85.5	85.7	86	86	11-162	0	
bis(2-Chloroethoxy)methane	ua/L	ND	100	100	81.5	80.5	82	81	33-184	1	
bis(2-Chloroethyl) ether	ua/L	ND	100	100	80.2	79.1	80	79	12-158	1	
bis(2-Chloroisopropyl) ether	ua/L	ND	100	100	82.4	82.0	82	82	36-166	1	
bis(2-Ethylhexyl)phthalate	ua/L	ND	100	100	80.4	81.3	80	81	8-158	1	
Butylbenzylphthalate	ua/L	ND	100	100	89.2	87.8	89	88	1-152	2	
Chrysene	ua/L	ND	100	100	85.0	85.2	85	85	17-168	0	
Di-n-butylphthalate	ua/L	ND	100	100	76.4	75.4	76	75	1-118	1	
Di-n-octylphthalate	ug/l	ND	100	100	93.6	94.8	94	95	4-146	1	
Dibenz(a,h)anthracene	ug/l	ND	100	100	65.2	67.3	65	67	1-227	3	
Diethylphthalate	ug/L	ND	100	100	85.9	83.1	86	83	1-114	3	
Dimethylphthalate	ug/L	ND	100	100	86.8	84.8	87	85	1-112	2	
Eluoranthene	ug/L	ND	100	100	104	106	104	106	26-137	- 3	
Fluorene	ug/L	ND	100	100	93.8	90.7	94	.00	59-121	3	
Hexachloro-1.3-butadiene	ug/L	ND	100	100	79.2	77.3	79	77	24-116	2	
Hexachlorobenzene	ug/L	ND	100	100	69.8	67.3	70	67	1-152	4	
Hexachlorocyclopentadiene	ug/L	ND	100	100	42.4	39.3	42	39	25-150	8	
Hexachloroethane	ug/L	ND	100	100	73.7	77.1	74	77	40-113	5	
Indeno(1 2 3-cd)pyrene	ug/L	ND	100	100	52.4	54.2	52	54	1-171	3	
Isophorope	ug/L	ND	100	100	95.5	101	96	101	21-196	5	
N-Nitroso-di-n-propylamine	ug/L	ND	100	100	76.0	81.4	76	81	1-230	7	
N-Nitrosodimethylamine	ug/L	ND	100	100	58.8	56.6	59	57	25-150	4	
N-Nitrosodinbenylamine	ug/L	ND	100	100	60.0	60.0	60	60	25-150	0	
Nanhthalene	ug/L	ND	100	100	60.0	70.7	70	71	21-132	2	
Nitrobenzene	ug/L	ND	100	100	106	10.7	106	104	25-180	2	
Pentachlorophenol	ug/L		200	200	100	162	79	104 Q1	14-176	2	
Phononthrono	ug/L		100	100	82.0	80 B	20 20	Q1	54-120	1	
Phonol	ug/L		100	100	52.U	51 7	52	50	5 110	1	
	uy/L		100	100	JJ.1	51.7	54	52	0-112	4	

Results presented on this page are in the units indicated by the "Units" column except where an alternate unit is presented to the right of the result.

## **REPORT OF LABORATORY ANALYSIS**



Project: 2014-093 Johnston 34182.1.2

Pace Project No.: 92204085

MATRIX SPIKE & MATRIX SPI	KE DUPLICATE	: 12181	06		1218107						
			MS	MSD							
	922	04282003	Spike	Spike	MS	MSD	MS	MSD	% Rec		
Parameter	Units	Result	Conc.	Conc.	Result	Result	% Rec	% Rec	Limits	RPD	Qual
Pyrene	ug/L	ND	100	100	77.3	72.1	77	72	52-115	7	
2,4,6-Tribromophenol (S)	%						113	111	10-137		
2-Fluorobiphenyl (S)	%						71	70	15-120		
2-Fluorophenol (S)	%						55	53	10-120		
Nitrobenzene-d5 (S)	%						96	97	10-120		
Phenol-d6 (S)	%						46	46	10-120		
Terphenyl-d14 (S)	%						78	75	11-131		

Results presented on this page are in the units indicated by the "Units" column except where an alternate unit is presented to the right of the result.



Analysis Method:

Matrix: Solid

Project: 2014-093 Johnston 34182.1.2

Pace Project No.: 92204085

QC Batch: OEXT/28108

QC Batch Method: EPA 3546 92204085001

Analysis Description:

EPA 8270 8270 Solid MSSV Microwave

Associated Lab Samples:

METHOD BLANK: 1214764

Associated Lab Samples: 92204	085001				
		Blank	Reporting		
Parameter	Units	Result	Limit	Analyzed	Qualifiers
1,2,4-Trichlorobenzene	ug/kg	ND	330	06/09/14 11:28	
1,2-Dichlorobenzene	ug/kg	ND	330	06/09/14 11:28	
1,3-Dichlorobenzene	ug/kg	ND	330	06/09/14 11:28	
1,4-Dichlorobenzene	ug/kg	ND	330	06/09/14 11:28	
1-Methylnaphthalene	ug/kg	ND	330	06/09/14 11:28	
2,4,5-Trichlorophenol	ug/kg	ND	330	06/09/14 11:28	
2,4,6-Trichlorophenol	ug/kg	ND	330	06/09/14 11:28	
2,4-Dichlorophenol	ug/kg	ND	330	06/09/14 11:28	
2,4-Dimethylphenol	ug/kg	ND	330	06/09/14 11:28	
2,4-Dinitrophenol	ug/kg	ND	1650	06/09/14 11:28	
2,4-Dinitrotoluene	ug/kg	ND	330	06/09/14 11:28	
2,6-Dinitrotoluene	ug/kg	ND	330	06/09/14 11:28	
2-Chloronaphthalene	ug/kg	ND	330	06/09/14 11:28	
2-Chlorophenol	ug/kg	ND	330	06/09/14 11:28	
2-Methylnaphthalene	ug/kg	ND	330	06/09/14 11:28	
2-Methylphenol(o-Cresol)	ug/kg	ND	330	06/09/14 11:28	
2-Nitroaniline	ug/kg	ND	1650	06/09/14 11:28	
2-Nitrophenol	ug/kg	ND	330	06/09/14 11:28	
3&4-Methylphenol(m&p Cresol)	ug/kg	ND	330	06/09/14 11:28	
3,3'-Dichlorobenzidine	ug/kg	ND	1650	06/09/14 11:28	
3-Nitroaniline	ug/kg	ND	1650	06/09/14 11:28	
4,6-Dinitro-2-methylphenol	ug/kg	ND	660	06/09/14 11:28	
4-Bromophenylphenyl ether	ug/kg	ND	330	06/09/14 11:28	
4-Chloro-3-methylphenol	ug/kg	ND	660	06/09/14 11:28	
4-Chloroaniline	ug/kg	ND	1650	06/09/14 11:28	
4-Chlorophenylphenyl ether	ug/kg	ND	330	06/09/14 11:28	
4-Nitroaniline	ug/kg	ND	660	06/09/14 11:28	
4-Nitrophenol	ug/kg	ND	1650	06/09/14 11:28	
Acenaphthene	ug/kg	ND	330	06/09/14 11:28	
Acenaphthylene	ug/kg	ND	330	06/09/14 11:28	
Aniline	ug/kg	ND	330	06/09/14 11:28	
Anthracene	ug/kg	ND	330	06/09/14 11:28	
Benzo(a)anthracene	ug/kg	ND	330	06/09/14 11:28	
Benzo(a)pyrene	ug/kg	ND	330	06/09/14 11:28	
Benzo(b)fluoranthene	ug/kg	ND	330	06/09/14 11:28	
Benzo(g,h,i)perylene	ug/kg	ND	330	06/09/14 11:28	
Benzo(k)fluoranthene	ug/kg	ND	330	06/09/14 11:28	
Benzoic Acid	ug/kg	ND	1650	06/09/14 11:28	
Benzyl alcohol	ug/kg	ND	660	06/09/14 11:28	
bis(2-Chloroethoxy)methane	ug/kg	ND	330	06/09/14 11:28	
bis(2-Chloroethyl) ether	ug/kg	ND	330	06/09/14 11:28	

Results presented on this page are in the units indicated by the "Units" column except where an alternate unit is presented to the right of the result.

# **REPORT OF LABORATORY ANALYSIS**



Project: 2014-093 Johnston 34182.1.2

# Pace Project No.: 92204085

METHOD BLANK:	1214764	

Associated Lab Samples: 92204085001

Matrix: Solid

	Blank Reporting		Reporting		
Parameter	Units	Result	Limit	Analyzed	Qualifiers
bis(2-Chloroisopropyl) ether	ug/kg	ND	330	06/09/14 11:28	
bis(2-Ethylhexyl)phthalate	ug/kg	ND	330	06/09/14 11:28	
Butylbenzylphthalate	ug/kg	ND	330	06/09/14 11:28	
Chrysene	ug/kg	ND	330	06/09/14 11:28	
Di-n-butylphthalate	ug/kg	ND	330	06/09/14 11:28	
Di-n-octylphthalate	ug/kg	ND	330	06/09/14 11:28	
Dibenz(a,h)anthracene	ug/kg	ND	330	06/09/14 11:28	
Dibenzofuran	ug/kg	ND	330	06/09/14 11:28	
Diethylphthalate	ug/kg	ND	330	06/09/14 11:28	
Dimethylphthalate	ug/kg	ND	330	06/09/14 11:28	
Fluoranthene	ug/kg	ND	330	06/09/14 11:28	
Fluorene	ug/kg	ND	330	06/09/14 11:28	
Hexachloro-1,3-butadiene	ug/kg	ND	330	06/09/14 11:28	
Hexachlorobenzene	ug/kg	ND	330	06/09/14 11:28	
Hexachlorocyclopentadiene	ug/kg	ND	330	06/09/14 11:28	
Hexachloroethane	ug/kg	ND	330	06/09/14 11:28	
Indeno(1,2,3-cd)pyrene	ug/kg	ND	330	06/09/14 11:28	
Isophorone	ug/kg	ND	330	06/09/14 11:28	
N-Nitroso-di-n-propylamine	ug/kg	ND	330	06/09/14 11:28	
N-Nitrosodimethylamine	ug/kg	ND	330	06/09/14 11:28	
N-Nitrosodiphenylamine	ug/kg	ND	330	06/09/14 11:28	
Naphthalene	ug/kg	ND	330	06/09/14 11:28	
Nitrobenzene	ug/kg	ND	330	06/09/14 11:28	
Pentachlorophenol	ug/kg	ND	1650	06/09/14 11:28	
Phenanthrene	ug/kg	ND	330	06/09/14 11:28	
Phenol	ug/kg	ND	330	06/09/14 11:28	
Pyrene	ug/kg	ND	330	06/09/14 11:28	
2,4,6-Tribromophenol (S)	%	38	27-110	06/09/14 11:28	
2-Fluorobiphenyl (S)	%	40	30-110	06/09/14 11:28	
2-Fluorophenol (S)	%	40	13-110	06/09/14 11:28	
Nitrobenzene-d5 (S)	%	37	23-110	06/09/14 11:28	
Phenol-d6 (S)	%	41	22-110	06/09/14 11:28	
Terphenyl-d14 (S)	%	48	28-110	06/09/14 11:28	

#### LABORATORY CONTROL SAMPLE: 1214765

		Spike	LCS	LCS	% Rec	
Parameter	Units	Conc.	Result	% Rec	Limits	Qualifiers
1,2,4-Trichlorobenzene	ug/kg	1670	1230	74	39-101	
1,2-Dichlorobenzene	ug/kg	1670	1180	71	36-110	
1,3-Dichlorobenzene	ug/kg	1670	1160	70	35-110	
1,4-Dichlorobenzene	ug/kg	1670	1210	72	35-110	
1-Methylnaphthalene	ug/kg	1670	1110	66	45-105	
2,4,5-Trichlorophenol	ug/kg	1670	1240	75	48-109	
2,4,6-Trichlorophenol	ug/kg	1670	1240	75	45-111	

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# **REPORT OF LABORATORY ANALYSIS**



Project: 2014-093 Johnston 34182.1.2

Pace Project No.: 92204085

#### LABORATORY CONTROL SAMPLE: 1214765

		Spike	LCS	LCS	% Rec	
Parameter	Units	Conc.	Result	% Rec	Limits	Qualifiers
2,4-Dichlorophenol	ug/kg		1270	76	51-116	
2,4-Dimethylphenol	ug/kg	1670	1220	73	42-103	
2,4-Dinitrophenol	ug/kg	8330	4440	53	28-103	
2,4-Dinitrotoluene	ug/kg	1670	1430	86	46-114	
2,6-Dinitrotoluene	ug/kg	1670	1450	87	48-112	
2-Chloronaphthalene	ug/kg	1670	1250	75	44-105	
2-Chlorophenol	ug/kg	1670	1270	76	36-110	
2-Methylnaphthalene	ug/kg	1670	1210	72	39-112	
2-Methylphenol(o-Cresol)	ug/kg	1670	1210	72	39-101	
2-Nitroaniline	ug/kg	3330	2660	80	44-111	
2-Nitrophenol	ug/kg	1670	1240	75	41-100	
3&4-Methylphenol(m&p Cresol)	ug/kg	1670	1180	71	43-103	
3,3'-Dichlorobenzidine	ug/kg	3330	2450	73	10-150	
3-Nitroaniline	ug/kg	3330	2530	76	35-110	
4,6-Dinitro-2-methylphenol	ug/kg	3330	2430	73	38-118	
4-Bromophenylphenyl ether	ug/kg	1670	1330	80	47-115	
4-Chloro-3-methylphenol	ug/kg	3330	2400	72	43-127	
4-Chloroaniline	ug/kg	3330	2540	76	34-109	
4-Chlorophenylphenyl ether	ug/kg	1670	1260	75	44-115	
4-Nitroaniline	ug/kg	3330	2660	80	37-111	
4-Nitrophenol	ug/kg	8330	5630	68	21-152	
Acenaphthene	ug/kg	1670	1200	72	38-117	
Acenaphthylene	ug/kg	1670	1220	73	46-107	
Aniline	ug/kg	1670	1110	66	29-110	
Anthracene	ug/kg	1670	1250	75	50-110	
Benzo(a)anthracene	ug/kg	1670	1280	77	47-116	
Benzo(a)pyrene	ug/kg	1670	1310	79	47-106	
Benzo(b)fluoranthene	ug/kg	1670	1260	76	47-109	
Benzo(g,h,i)perylene	ug/kg	1670	1200	72	39-115	
Benzo(k)fluoranthene	ug/kg	1670	1280	77	45-117	
Benzoic Acid	ug/kg	8330	3690	44	16-110	
Benzyl alcohol	ug/kg	3330	2390	72	38-105	
bis(2-Chloroethoxy)methane	ug/kg	1670	1210	73	39-110	
bis(2-Chloroethyl) ether	ug/kg	1670	1220	73	19-119	
bis(2-Chloroisopropyl) ether	ug/kg	1670	1020	61	21-110	
bis(2-Ethylhexyl)phthalate	ug/kg	1670	1410	85	35-116	
Butylbenzylphthalate	ug/kg	1670	1440	87	38-110	
Chrysene	ug/kg	1670	1280	77	49-110	
Di-n-butylphthalate	ug/kg	1670	1240	74	43-109	
Di-n-octylphthalate	ug/kg	1670	1400	84	37-109	
Dibenz(a,h)anthracene	ug/kg	1670	1280	77	43-116	
Dibenzofuran	ug/kg	1670	1300	78	45-106	
Diethylphthalate	ug/kg	1670	1160	70	41-114	
Dimethylphthalate	ug/kg	1670	1190	71	43-110	
Fluoranthene	ug/kg	1670	1180	71	50-114	
Fluorene	ug/kg	1670	1210	72	46-114	
Hexachloro-1,3-butadiene	ug/kg	1670	1190	71	28-111	

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## **REPORT OF LABORATORY ANALYSIS**



Project: 2014-093 Johnston 34182.1.2

Pace Project No.: 92204085

#### LABORATORY CONTROL SAMPLE: 1214765

		Spike	LCS	LCS	% Rec	
Parameter	Units	Conc.	Result	% Rec	Limits	Qualifiers
Hexachlorobenzene	ug/kg	1670	1230	74	46-120	
Hexachlorocyclopentadiene	ug/kg	1670	1350	81	18-119	
Hexachloroethane	ug/kg	1670	1150	69	33-110	
Indeno(1,2,3-cd)pyrene	ug/kg	1670	972	58	42-115	
Isophorone	ug/kg	1670	1240	74	44-109	
N-Nitroso-di-n-propylamine	ug/kg	1670	1050	63	43-104	
N-Nitrosodimethylamine	ug/kg	1670	1170	70	29-110	
N-Nitrosodiphenylamine	ug/kg	1670	1290	78	48-113	
Naphthalene	ug/kg	1670	1180	71	41-110	
Nitrobenzene	ug/kg	1670	1380	83	38-110	
Pentachlorophenol	ug/kg	3330	2100	63	32-128	
Phenanthrene	ug/kg	1670	1200	72	50-110	
Phenol	ug/kg	1670	1200	72	28-106	
Pyrene	ug/kg	1670	1380	83	45-114	
2,4,6-Tribromophenol (S)	%			79	27-110	
2-Fluorobiphenyl (S)	%			77	30-110	
2-Fluorophenol (S)	%			75	13-110	
Nitrobenzene-d5 (S)	%			71	23-110	
Phenol-d6 (S)	%			73	22-110	
Terphenyl-d14 (S)	%			85	28-110	

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# **REPORT OF LABORATORY ANALYSIS**



Project:	2014-093 Johnsto	n 34182.1.2					
Pace Project No.:	92204085						
QC Batch:	PMST/6693		Analysis Meth	nod: A	ASTM D2974-87	7	
QC Batch Method:	ASTM D2974-87	,	Analysis Desc	cription: [	Dry Weight/Perc	ent Moisture	
Associated Lab San	nples: 92204085	001					
SAMPLE DUPLICA	TE: 1219653						
			92204084001	Dup			
Paran	neter	Units	Result	Result	RPD	Qualifiers	
Percent Moisture		%	14.1	13.8	8	2	
SAMPLE DUPLICA	TE: 1219654						
			92204947001	Dup			
Paran	neter	Units	Result	Result	RPD	Qualifiers	

Results presented on this page are in the units indicated by the "Units" column except where an alternate unit is presented to the right of the result.



## QUALIFIERS

Project: 2014-093 Johnston 34182.1.2

Pace Project No.: 92204085

#### DEFINITIONS

DF - Dilution Factor, if reported, represents the factor applied to the reported data due to changes in sample preparation, dilution of the sample aliquot, or moisture content.

ND - Not Detected at or above adjusted reporting limit.

J - Estimated concentration above the adjusted method detection limit and below the adjusted reporting limit.

MDL - Adjusted Method Detection Limit.

PRL - Pace Reporting Limit.

RL - Reporting Limit.

S - Surrogate

1,2-Diphenylhydrazine (8270 listed analyte) decomposes to Azobenzene.

Consistent with EPA guidelines, unrounded data are displayed and have been used to calculate % recovery and RPD values.

LCS(D) - Laboratory Control Sample (Duplicate)

MS(D) - Matrix Spike (Duplicate)

DUP - Sample Duplicate

RPD - Relative Percent Difference

NC - Not Calculable.

SG - Silica Gel - Clean-Up

U - Indicates the compound was analyzed for, but not detected.

N-Nitrosodiphenylamine decomposes and cannot be separated from Diphenylamine using Method 8270. The result reported for each analyte is a combined concentration.

Acid preservation may not be appropriate for 2-Chloroethylvinyl ether, Styrene, and Vinyl chloride.

Pace Analytical is TNI accredited. Contact your Pace PM for the current list of accredited analytes.

TNI - The NELAC Institute.

#### LABORATORIES

PASI-C Pace Analytical Services - Charlotte

#### ANALYTE QUALIFIERS

- F3 The recovery of the second source standard used to verify the initial calibration curve for this analyte is outside the laboratory's control limits. The result is estimated.
- M0 Matrix spike recovery and/or matrix spike duplicate recovery was outside laboratory control limits.
- R1 RPD value was outside control limits.



# QUALITY CONTROL DATA CROSS REFERENCE TABLE

 Project:
 2014-093 Johnston 34182.1.2

 Pace Project No.:
 92204085

Lab ID	Sample ID	QC Batch Method	QC Batch	Analytical Method	Analytical Batch
92204085002	8-4 (TW)	EPA 625	OEXT/28189	EPA 625	MSSV/9235
92204085001	8-2 (4-6)	EPA 3546	OEXT/28108	EPA 8270	MSSV/9220
92204085002	8-4 (TW)	SM 6200B	MSV/27102		
92204085001	8-2 (4-6)	EPA 8260	MSV/27125		
92204085001	8-2 (4-6)	ASTM D2974-87	PMST/6693		

		Document Name:			Document Revised: A	oril 07. 2014	
Pace Analytical*	Sample C	Condition U	pon R	eceipt (SCL	JR)	Issuing Autho	z rity:
(		F-CHR-CS-	003-re	er. ev.14	10	Pace Huntersville Qu	ality Office
Client Name: lypmid E							
	PS Client		rcial	Pace O	ther	Optional	1
Custody Seal on Cooler/Box Presen	t: 🖵 yes	no 🖓	Seals i	intact:	yes 🗌	no Proj. Due Da Proj. Name:	te:
Packing Material: 🗌 Bubble Wrap	Bubble B	Bags 🗌 No	ne 🗌	Other	/		
Thermometer Used: IR Gun T1102	T1401	Type of Ice:	Wet	Blue Non	ie 🛛	Samples on ice, cooling pro	ocess has begun
Temp Correction Factor T1102	No Correct	ion <b>T1</b> :	301: N	lo Correction	n Č		to to a
Corrected Cooler Temp.:	°C	<b>Biological T</b>	issue i	is Frozen: Y	es No (N/A)	contents: C (2/4/	lu
Temp should be above freezing to 6°C				Comments:	Ŭ		
Chain of Custody Present:		TYes DNo	□n/A	1.			
Chain of Custody Filled Out:		ØYes □No	□n/A	2.			
Chain of Custody Relinquished:		PYes □No	□n/A	3.	_		
Sampler Name & Signature on COC:		Pres DNo	□n/A	4.	-		
Samples Arrived within Hold Time:		ØYes □No	□n/a	5.	4	÷	
Short Hold Time Analysis (<72hr):		□Yes □No	□n/a	6.			
Rush Turn Around Time Requested:			□n/a	7.			
Sufficient Volume:		Yes No	□n/A	8.			
Correct Containers Used:		ØYes □No		9.			
-Pace Containers Used:		ZYes DNo					
Containers Intact:			□n/A	10.			
Filtered volume received for Dissolved	tests	□Yes □No		11.			
Sample Labels match COC:	10010	ØYes □No		12.			
-Includes date/time/ID/Analysis	Matrix:		_				
All containers needing preservation have been	h checked.	□Yes □No		13.			
All containers needing preservation are four compliance with EPA recommendation.	nd to be in	□Yes □No					
exceptions: VOA, coliform, TOC, O&G, WI-DRO	(water)	ZYes DNo					
Samples checked for dechlorination:		Yes DNo	□n/A	14.			
Headspace in VOA Vials ( >6mm):		□Yes ØNo	□n/A	15.			
Trip Blank Present:		□Yes □No		16.			
Trip Blank Custody Seals Present		□Yes □No					
Pace Trip Blank Lot # (if purchased):_		-	V				
Client Notification/ Resolution:						Field Data Required?	Y / N
Person Contacted:			Date/	Time:		unu nu no della ministra della de	
Comments/ Resolution:							
					10.11	0000	6000
SCURF Review:	Date:	614/14			WU# :	9220408	5
SRF Review: JP3	Date:	6414					
samples, a copy of this form will be se	necting North ( nt to the North	Carolina compli Carolina DEHN ative, out of ten	IR NR		92204085		

samples, a copy of this form will be sent to the North Carolina DEHNR Certification Office ( i.e out of hold, incorrect preservative, out of temp, incorrect containers)



# APPENDIX F

	FIELD PERSONNEL LOG				
<b>PROJECT NAME</b> : NC PARCELS 1, 2, 5, 8	DOT Johnston County PS.	As <b>PROJECT NO.:</b> I-3318BB			
Name: Eric Cross	<b>Date:</b> 5/13/14 N	Mon Tue Wed Th Fri Sat Sun			
TASKS PERFORMED	:				
E. Cross: Mobilize to site. Perform Leave site: ~3:30PM	ed site visit reconnaissance				

# FIELD PERSONNEL LOG

**PROJECT NAME**: NCDOT Johnston County PSAs PARCELS 1, 2, 5, 8

PROJECT NO.: I-3318BB

Name: Eric Cross & Alan McFaddenDate: 5/21/14Mon Tue WedTh Fri Sat Sun

# **TASKS PERFORMED:**

*E. Cross & A. McFadden:* Mobilize to site. Performed geophysical surveys and data processing in evening. Leave site: ~5:30PM

FIELD	) PERSONNEL I	LOG
<b>PROJECT NAME</b> : NCDOT Johnston ( PARCELS 1, 2, 5, 8	County PSAs	PROJECT NO.: I-3318BB
Name: Eric Cross & Alan McFadden	<b>Date:</b> 5/22/14	Mon Tue Wed Th Fri Sat Sun
TASKS PERFORMED:		
<i>E. Cross &amp; A. McFadden:</i> Mobilize to site. Performed geophysical Leave site: ~12:30PM	surveys.	
·		

	FIELD PERS	SONNEL 1	LOG
<b>PROJECT NAME</b> : NCI PARCELS 1, 2, 5, 8	DOT Johnston County	PSAs	PROJECT NO.: I-3318BB
Name: Eric Cross	<b>Date:</b> 5/23/14	Mon Tu	e Wed Th Fri Sat Sun
TASKS PERFORMED	:		
<i>E. Cross:</i> Mobilize to site. Performe Leave site: ~5:00PM	ed geophysical surveys	(GPR) and	d data processing.

FIELD PERSONNEL LOG				
<b>PROJECT NAME</b> : NCDOT Johnston Co PARCELS 1, 2, 5, 8	ounty PSAs	PROJECT NO.: I-3318BB		
Name: Eric Cross & Tim Leatherman	<b>Date:</b> 5/29/14	Mon Tue Wed Th Fri Sat Sun		
TASKS PERFORMED:				
<i>E. Cross &amp; T. Leatherman:</i> Mobilize to site. Placed proposed boring le Leave site: ~5:00PM	ocations and supe	ervised private utility locating.		

FIELD PERSONNEL LOG				
<b>PROJECT NAME:</b> NCDOT Johnston County PSAs <b>PROJECT NO.:</b> I-3318BBPARCELS 1, 2, 5, 8				
Name: Tim Leatherman & Ryan Kramer Date: 6/2/14 Mon Tue Wed Th Fri Sat Sun				
TASKS PERFORMED:				
<i>T. Leatherman &amp; R. Kramer:</i> Mobilize to site. Supervised Geoprobe sampling, performed QED analysis (some in evening). Leave site: ~5:30PM				

# FIELD PERSONNEL LOG

**PROJECT NAME**: NCDOT Johnston County PSAs PARCELS 1, 2, 5, 8

PROJECT NO.: I-3318BB

Name: Tim Leatherman & Ryan Kramer Date: 6/3/14 Mon Tue Wed Th Fri Sat Sun

# TASKS PERFORMED:

*T. Leatherman & R. Kramer:* Mobilize to site. Supervised Geoprobe sampling, performed QED analysis (some in evening). Leave site: ~4:30PM

# FIELD PERSONNEL LOG

**PROJECT NAME**: NCDOT Johnston County PSAs PARCELS 1, 2, 5, 8

PROJECT NO.: I-3318BB

Name: Tim Leatherman & Ryan Kramer Date: 6/4/14 Mon Tue Wed Th Fri Sat Sun

# **TASKS PERFORMED:**

*T. Leatherman & R. Kramer:* Mobilize to site. Supervised Geoprobe sampling, performed QED analysis. Leave site: ~3:00PM