PRELIMINARY SITE ASSESSMENT

PARCEL 003 NAEGELE OUTDOOR ADVERTISING, INC. PROPERTY 1730 BURNETT BLVD. WILMINGTON, NEW HANOVER COUNTY, NORTH CAROLINA

INTERSECTION OF SR 1436 / US 421 TRUCK (FRONT STREET) AND SR 1140 (BURNETT BLVD.) SOUTH OF WILLARD STREET WBS ELEMENT: 17BP.3.R.28

CATLIN PROJECT NO. 214037

PREPARED FOR:



NCDOT GEOTECHNICAL ENGINEERING UNIT-GEOENVIRONMENTAL SECTION 1589 MSC RALEIGH, NORTH CAROLINA 27699-1589

JUNE 25, 2014

PREPARED BY:

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1.0 INTRODUCTION

The North Carolina Department of Transportation (NCDOT) is planning construction activities and acquisition of the right-of-way (ROW) is necessary for intersection improvements at the Greenfield Lake Spillway Culvert (above referenced WBS Element 17BP.3.R.28). NCDOT has indicated site investigations are necessary to determine the presence of contaminated groundwater and/or soil at three (3) sites along the proposed construction area.

2.0 PURPOSE OF INVESTIGATION AND DESCRIPTION

Catlin Engineers & Scientists (CATLIN) was retained by the NCDOT Geotechnical Engineering Unit to provide field investigations concluding with Preliminary Site Assessment (PSA) reports for three (3) sites associated with WBS Element 17BP.3.R.28. In response to a Request for Technical and Cost Proposal (RFP) dated March 17, 2014, CATLIN submitted a proposal for conducting PSAs at the three (3) sites. This report documents the investigation at Parcel 003, Naegele Outdoor Advertising, Inc. Property, 1730 Burnett Blvd. in Wilmington, North Carolina 28401. The property is a vacant lot except for an advertisement billboard. The general location is illustrated on Sheet 1. CATLIN personnel conducted a field investigation at the property on May 16, 2014. This PSA report documents activities and findings.

According to the RFP, this heavily vegetated property is undeveloped. The proposed ROW will be expanded up to 35 feet onto this parcel. This parcel does not appear on the North Carolina Department of Environment and Natural Resources (NCDENR) Underground Storage Tank (UST) Section registry or Ground Water Incident database.

The requested area of investigation is within the proposed right of way (ROW) and easement along Front Street near the intersection of Burnett Blvd. Borings were proposed within the proposed ROW and along planned

drainage features including catch basins and drainage piping. The NCDOT conventional plan sheet symbols are provided on Sheet 2 and the site layout including proposed features are illustrated on Sheet 3.

The NCDOT has requested an investigation to determine if contamination is present at the site. The purpose of this investigation was to:

- Locate all USTs and determine approximate size and contents (if any).
- Determine if contaminated soils are present.
- If contamination is evident, estimate the quantity of impacted soils and indicate the approximate area of soil contamination on a site map.
- Provide a MicroStation file with the location of USTs, soil contamination and monitoring wells.
- Prepare a report including field activities, findings, and recommendations for this site and submit to this office in triplicate.

3.0 METHODS

Proposed boring/sample locations were illustrated on a Plan Sheet provided by NCDOT and agreed upon before beginning investigations. Borings/samples were approved by NCDOT at proposed drainage catch basin locations, along the proposed drainage features, and near the edge of the proposed ROW.

CATLIN coordinated geophysical activities with Pyramid Environmental and Engineering (Pyramid). The geophysical investigation methods are detailed in the Pyramid geophysical report provided in Appendix A.

CATLIN proposed utilizing QROS On-Site Rapid measurement Techniques and Tools (QED[™] Analyzer) to evaluate potential for petroleum impacts to soil in a cost effective manner. Soil samples collected from above the approximate water table depth with concentrations greater than 10 milligrams per kilogram (mg/kg) diesel range organics (DRO) or gasoline range organics (GRO) will be considered contaminated for estimated contaminated vadose soil volume calculations. Contaminated soil volume is estimated from the surface to the water table and/or the midpoint distance between a clean sample location and dirty sample location or the property line and ROW/easement. Saturated soils were encountered two (2) to three (3) feet below land surface (BLS).

Borings advanced during this investigation are identified with the parcel number prefix ("3") and numbered sequentially "##". Proposed boring locations and nomenclature were established before beginning field work. During field work, it was determined that proposed borings 3-01 and 3-02 could not be accessed due to heavy vegetation (wooded area) so therefore, there are no borings/samples numbered "-01" or "-02". Soil samples for

analysis per QROS QED[™] Analyzer were identified by parcel number, boring number, and depth [example: 3-03 (2')].

3.1 FIELD METHODS

All field work was conducted in general accordance with state and federal guidelines and industry standards.

Underground utility locating was coordinated by CATLIN personnel. The North Carolina One Call Center (NC-1-Call) was contacted for underground utility location. The areas around the proposed boring locations were checked and underground utilities were indicated by NC-1-Call personnel.

CATLIN personnel gathered subsurface soil data by Direct Push Technology boring advancement using an AMS PowerProbe[™] 9600D (PowerProbe) and a hand auger (at three locations, 3-03, 3-04, and 3-07). When using the PowerProbe, the borings are advanced to depth by static force and a 90-pound hydraulic percussion hammer. Two and one-quarter inch diameter by four-foot length steel is used as casing. Soil samples are continuously collected in one and one-half inch clear liners. Liners are removed from the casing and then cut in half longitudinally to allow for visual/manual classification utilizing the Unified Soil Classification System (USCS). Boring information was recorded on field logs and transferred to Boring Logs (see Appendix B). Soil samples were collected and packed in the appropriate glassware for analysis.

Following removal of the PowerProbe tooling at boring 3-08, groundwater was pumped from the open bore hole directly into the appropriate laboratory provided glassware utilizing new polypropylene tubing and a peristaltic pump.

New disposable nitrile gloves were worn during sampling activities. Soils selected for QROS QED[™] analysis were placed into new glassware provided by QROS. All samples were placed on ice in an insulated cooler for transportation to the laboratory. Sample integrity was maintained by following proper Chain of Custody procedures. A copy of the Chain of Custody is provided following the analytical report in Appendix C.

The groundwater sample was placed in a cooler and transported following proper Chain of Custody procedures to Pace Analytical Services, Inc. (Pace) laboratory in Huntersville, North Carolina by a Pace courier. A copy of the Chain of Custody is provided following the analytical report in Appendix D. Boreholes were abandoned to the surface in grassy areas and just below existing asphalt in asphalt areas using three-eighth inch bentonite chips. Bentonite and water were poured into the borehole simultaneously to facilitate hydration. Boreholes in asphalt were finished with asphalt patch to the surface.

3.2 ANALYTICAL TESTING

<u>Soil</u>

The QROS QED[™] Analyzer methods have been approved by the NCDENR for petroleum contamination determination. Complete QROS QED[™] procedures are on file with the NCDENR and are available upon request. The QROS QED[™] analysis was conducted by QROS personnel at their laboratory in Wilmington, North Carolina.

QROS QED[™] analysis provides total Benzene, Toluene, Ethylbenzene, and Toluene (BTEX), DRO, GRO, total petroleum hydrocarbon (TPH), total aromatics (C-10-C35) and (total) 16 Environmental Protection Agency (EPA) Poly Aromatic Hydrocarbons (PAHs) concentrations. Soil sample DRO and GRO results greater than 10 mg/kg are considered contaminated for this investigation.

Groundwater

One groundwater sample was collected near the proposed replacement culvert under Front Street. The groundwater sample was submitted to Pace (NC Certification #12) for volatile and semi volatile organics analysis per Standard Method (SM) 6200B and EPA 625, respectively.

4.0 FIELD ACTIVITIES

4.1 CURRENT SITE CONDITIONS AND FIELD OBSERVATIONS

As previously mentioned, the site is a vacant lot with an advertising billboard and portions of the site are covered with heavy vegetation (wooded). No signs of USTs were observed. Photographs taken during the geophysical investigation are included in the geophysical report provided in Appendix A.

The site vicinity is illustrated on Sheet 1 and Sheet 3 illustrates the current site map with soil boring and sample locations.

4.2 SOIL SAMPLING

A total of six (6) borings were installed as part of the investigation. At least one (1) soil sample interval was collected from each boring and submitted for analysis. Boring/sample locations are illustrated on Sheet 3. Boring logs are included in Appendix B.

PowerProbe borings were advanced to approximately four (4) feet deep and terminated in saturated soils except boring 3-08 was terminated at eight (8) feet BLS. One hand auger boring (3-07) was terminated at one foot deep. Soils were collected continuously to boring termination. After retrieving the drive, soil was visually/manually classified for USCS classification. Soil samples collected from each boring for analysis were packed in the appropriate glassware, labeled, and placed in a cooler on ice. Borings 3-03 and 3-06 were located at proposed catch basins and two (2) soil samples were collected for analysis at these locations (one soil sample at 2 feet BLS and one at 4 feet BLS). A total of eight (8) soil samples were submitted to QROS for QED[™] analysis. Chain of Custody documentation is included in Appendix C.

4.3 GROUNDWATER SAMPLING

Boring 3-08 was advanced for groundwater sample collection. Groundwater was encountered at approximately three (3) feet BLS. A groundwater sample was pumped directly from the open bore hole into laboratory provided glassware. The groundwater samples was submitted to Pace for analysis per SM 6200B and EPA Method 625.

4.4 SURVEYING

Boring/sample locations were recorded utilizing a Trimble[®] global positioning survey instrument and data collector. Boring coordinates are shown on the Boring Logs provided in Appendix B. Borings locations are indicated on plan sheets provided by NCDOT and are included as Sheet 3.

5.0 RESULTS

Geophysical Investigation

The complete geophysical investigation report is included in Appendix A. As indicated in the Pyramid Report, the investigation did not reveal any evidence of metallic USTs in the survey area.

<u>Soil</u>

Soil sample results from the recent assessment activities utilizing QROS QED[™] analysis are provided on Table 1. Soil sample locations, summarized results and estimated extent of TPH impacted soils are illustrated on Sheet 3. The complete QROS QED[™] report is provided in Appendix C.

Soils encountered across the site were predominately sands with gravel. Saturated soils were encountered approximately two (2) to three (3) feet BLS. Soils from the surface to two (2) feet BLS are considered vadose zone.

No GRO soil concentrations were reported above the detection limits. Soils collected from proposed catch basin (CB) 0425 (boring 3-03) revealed DRO concentrations greater than 10 mg/kg in the samples collected from two (2) and four (4) feet BLS (results = 127.1 mg/kg and 689.3 mg/kg DRO, respectively). Soil samples 3-06 (2') and 3-06 (4') were collected at CB 0423 and revealed DRO concentrations of 11.62 mg/kg and 10.38 mg/kg, respectively. The boring 3-08 advanced at proposed catch basin 0435 did not reveal DRO or GRO concentrations above 10 mg/kg in the soil sample collected from two (2) feet BLS.

The soil samples collected from two (2) feet BLS along the proposed drainage pipe at borings 3-04 and 3-05 revealed DRO soil contamination of 54.7 mg/kg and 20.22 mg/kg, respectively. The sample collected from approximately one foot deep at boring 3-07 and near the end of the proposed drainage line and ditch did not reveal soil contamination above 10 mg/kg.

The estimated volume of petroleum impacted soils as illustrated on Sheet 3 includes the area around all borings except borings 3-07 and 3-08. The approximate water table at both these locations is two (2) to three (3) feet BLS. The approximate area is 4,380 feet² and a total volume of impacted soils is approximately 320 yds³.

Groundwater

The one groundwater sample collected at boring 3-08 located near the end of the culvert under Front Street and CB 0435 did not reveal any SM 6200B or EPA Method 625 parameters above the method detection limits. The laboratory analytical report including a list of all parameters and method detection limits is provided in Appendix D.

6.0 SUMMARY AND CONCLUSIONS

The site is currently a vacant lot except for an advertising billboard. No USTs are suspected at the area of investigation. Six (6) borings were advanced for soil sample collection at proposed drainage features and within the proposed

ROW. There were no borings advanced that numbered "01" or "02" due to proposed boring locations being inaccessible in a heavily wooded area.

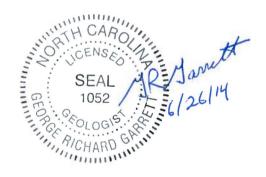
Contaminated soils were revealed by QROS QED[™] in borings 3-03 and 3-06 at proposed catch basins 0425 and 0423, respectively, in samples collected from two (2) and four (4) feet BLS. The borings 3-04 and 3-05 were advanced along the proposed drainage line and samples collected at two (2) feet BLS from these locations also revealed contamination. No soil or groundwater contamination was detected in samples collected from boring 3-08 near the culvert under Front Street and CB 0435. No soil contaminant concentrations were detected above 10 mg/kg in the soil sample collected from boring 3-07 (where the proposed drainage line ends at a ditch).

A total estimated contaminated soil volume of 320 yds³ may be encountered in vadose zone soils across the site and around borings 3-03 through 3-06. Any detectable concentrations in excavated soils may require handling and disposal as an impacted waste.

7.0 SIGNATURES



Benjamin J. Ashba, P.G. Project Manager

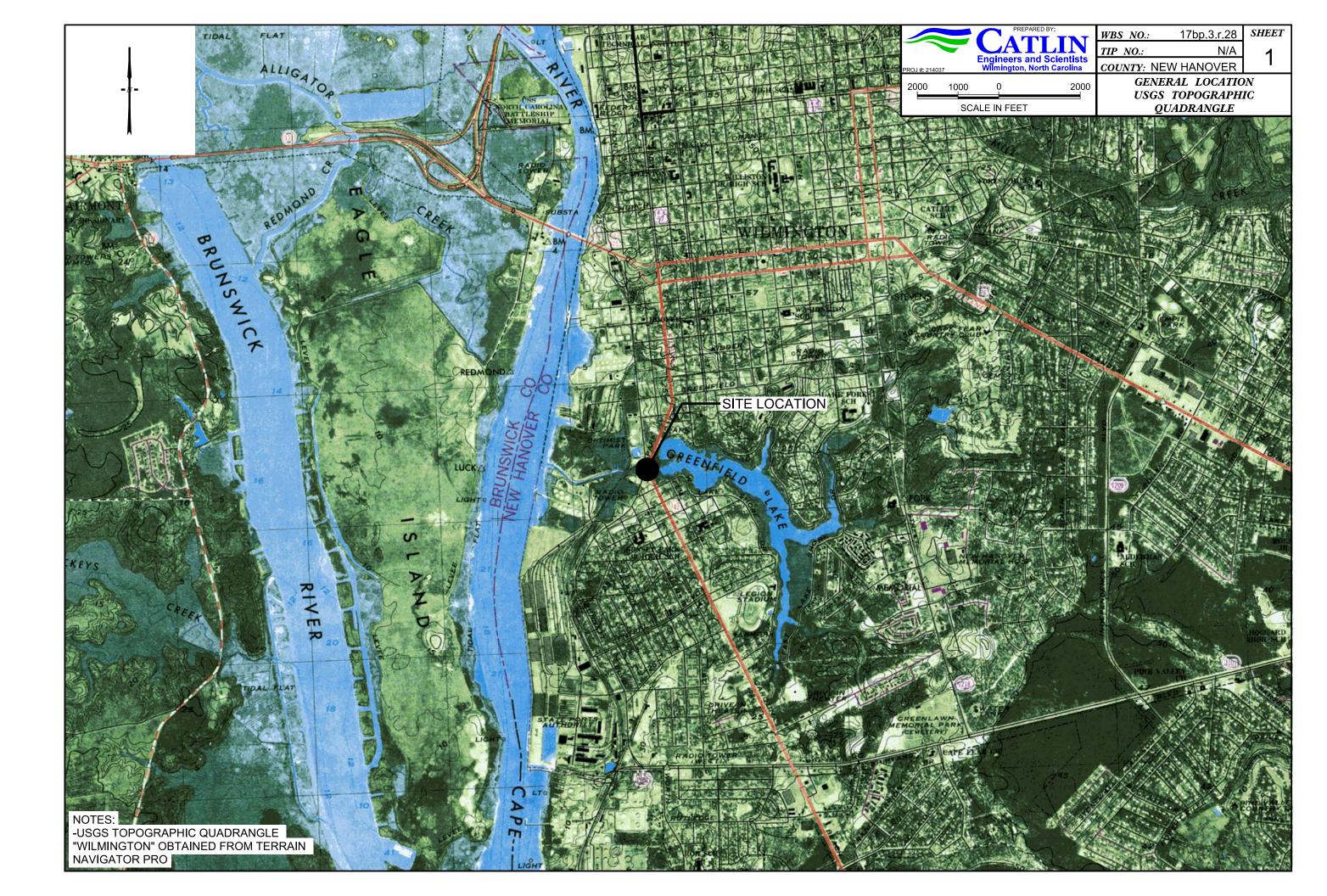


G. Richard Garrett, P.G. Contract Manager

TABLES

\square														DROS
					Hvdroca	arbon Ar	nalvsis R	esults						
											Samples taken Samples extracted Samples analysed Operator			Friday, May 16, 2014 Friday, May 16, 2014 Monday, May 19, 2014 Rachel Menoher
Project:		NCDOT Front St. and Burnett Blv oject No. 214037	vd - WBS	: 17BP.3	.R.28									
							Total		Ratios			HC Fingerprint Match		
Matrix	Sample ID	Location	Dilution used	BTEX (C6 - C9)	GRO (C5 - C10)	DRO (C10 - C35)	TPH (C5 - C35)	Aromatics (C10-C35)	16 EPA PAHs	BaP	% light	% mid	% heavy	
S	3-03 (2')	Proposed Catch	21.0	<1	<1	127.1	127.1	35.19	2.18	<0.021	57.9	37.4	4.7	Waste Oil (FCM) 97.7%
S	3-03 (4')	Basin (CB) 0425	256.0	<12.8	<12.8	689.3	689.3	161.2	5.67	<0.256	85.5	12.1	2.3	Waste Oil (FCM) 70.8%
S	3-04 (2')	North of CB 0425 along proposed drainage line	20.0	<1	<1	54.7	54.7	40.72	2.09	0.037	27.3	58.5	14.3	V.Deg.PHC 77.6%
S	3-05 (2')	Along proposed drainage line north of boring 3-03	19.0	<1	<1	20.22	20.22	19.12	1.02	<0.019	51.6	39.8	8.6	V.Deg.PHC 90.4%
S	3-06 (2')	CB 0423	22.0	<1.1	<1.1	11.62	11.62	10.65	0.71	<0.022	36.9	47.5	15.6	V.Deg.PHC 74.2%
S	3-06 (4')		19.0	<0.9	<0.9	10.38	10.38	10.35	0.64	<0.019	50.8	38.1	11.1	V.Deg.PHC 83.7%
	3-07 (1')	End of proposed drainage line at ditch	22.0	<1.1	<1.1	9.57	9.57	8.77	0.58	<0.022	40.9	44.1	15	V.Deg.PHC 76.8%
S	3-08 (2')	Near end of culvert under Front St. and CB 0435	21.0	<1.1	<1.1	5.52	5.52	5.06	0.34	<0.021	46.4	37.3	16.3	V.Deg.PHC 76.1%
s	, ,			QC check	OK					Final FC		Chock	OK	98.3%

SHEETS



Note: Not to Scale

*S.U.E. = Subsurface Utility Engineering

STATE OF NORTH CAROLINA DIVISION OF HIGHWAYS

CONVENTIONAL PLAN SHEET SYMBOLS

BOUNDARIES AND PROPERTY:

State Line	
County Line ———	
Township Line	
City Line	
Reservation Line	
Property Line	
Existing Iron Pin	
Property Corner	
Property Monument ————	
Parcel/Sequence Number	• • • • • • • • • • • • • • • • • • • •
Existing Fence Line	-×
Proposed Woven Wire Fence	
Proposed Chain Link Fence	
Proposed Barbed Wire Fence	
Existing Wetland Boundary	
Proposed Wetland Boundary	k3
Existing Endangered Animal Boundary	
Existing Endangered Plant Boundary	0
Known Soil Contamination: Area or Site	
Potential Soil Contamination: Area or Site	-xx - xx
Potential Soil Contamination: Area or Site — BUILDINGS AND OTHER CULTU	000
	000
BUILDINGS AND OTHER CULTU	VRE:
BUILDINGS AND OTHER CULTU Gas Pump Vent or U/G Tank Cap Sign Well	<i>VRE:</i>
BUILDINGS AND OTHER CULTU Gas Pump Vent or U/G Tank Cap Sign	<i>VRE:</i>
BUILDINGS AND OTHER CULTU Gas Pump Vent or U/G Tank Cap Sign Well	<i>VRE:</i>
BUILDINGS AND OTHER CULTU Gas Pump Vent or U/G Tank Cap Sign Well Small Mine	<i>VRE:</i>
BUILDINGS AND OTHER CULTUR Gas Pump Vent or U/G Tank Cap Sign Well Small Mine Foundation Area Outline Cemetery	<i>VRE:</i>
BUILDINGS AND OTHER CULTUR Gas Pump Vent or U/G Tank Cap Sign Well Small Mine Foundation Area Outline	<i>VRE:</i>
BUILDINGS AND OTHER CULTUR Gas Pump Vent or U/G Tank Cap Sign Well Small Mine Foundation Area Outline Cemetery	<i>VRE:</i>
BUILDINGS AND OTHER CULTUR Gas Pump Vent or U/G Tank Cap Sign Well Small Mine Foundation Area Outline Cemetery Building	<i>VRE:</i>
BUILDINGS AND OTHER CULTUR Gas Pump Vent or U/G Tank Cap Sign Well Small Mine Foundation Area Outline Cemetery Building School	<i>VRE:</i>
BUILDINGS AND OTHER CULTUR Gas Pump Vent or U/G Tank Cap Sign Well Small Mine Foundation Area Outline Cemetery Building School Church	<i>VRE:</i>
BUILDINGS AND OTHER CULTUR Gas Pump Vent or U/G Tank Cap Sign Well Small Mine Foundation Area Outline Cemetery Building School Church Dam	
BUILDINGS AND OTHER CULTUR Gas Pump Vent or U/G Tank Cap Sign Well Small Mine Foundation Area Outline Cemetery Building School Church Dam	
BUILDINGS AND OTHER CULTUR Gas Pump Vent or U/G Tank Cap Sign Well Small Mine Foundation Area Outline Cemetery Building School Church Dam HYDROLOGY: Stream or Body of Water Hydro, Pool or Reservoir	
BUILDINGS AND OTHER CULTUR Gas Pump Vent or U/G Tank Cap Sign Well Small Mine Foundation Area Outline Cemetery Building School Church Dam HYDROLOGY: Stream or Body of Water	

Existing from Fin	
Property Corner ————	* RIGHT OF WAY:
Property Monument ———	🔜 🛛 Baseline Control Point ————————————————————————————————————
Parcel/Sequence Number	🤨 Existing Right of Way Marker — 🗛 🛆
Existing Fence Line ————————————————————————————————————	
Proposed Woven Wire Fence	
Proposed Chain Link Fence	Proposed Right of Way Line with
Proposed Barbed Wire Fence	Iron Pin and Cap Marker
Existing Wetland Boundary	Proposed Right of Way Line with Concrete or Granite Marker
Proposed Wetland Boundary ————————————————————————————————————	Existing Control of Access
Existing Endangered Animal Boundary ———	
Existing Endangered Plant Boundary	
Known Soil Contamination: Area or Site ———— 🎅	
Potential Soil Contamination: Area or Site	
BUILDINGS AND OTHER CULTURE:	Proposed Permanent Drainage Easement PDE
Gas Pump Vent or U/G Tank Cap	Proposed Permanent Drainage / Utility Easement
Sign	Proposed Permanent Utility Easement PUE
Well	Proposed Temporary Utility Easement TUE
Small Mine	Proposed Aerial Utility Easement
Foundation ———	
Area Outline ———	Proposed Permanent Easement with Iron Pin and Cap Marker
Cemetery	ROADS AND RELATED FEATURES :
Building — C	Existing Edge of Pavement — — — — — — — — — — — — — — — — — — —
School	Existing Curb — — — — — — — — — — — — — — — — — — —
Church ———	Proposed Slope Stakes Cut
Dam	Proposed Slope Stakes Fill
	Proposed Curb Ramp CR
HYDROLOGY:	Curk Cut Euture Rame
Stream or Body of Water	Existing Metal Guardrail ————
Hydro, Pool or Reservoir ————————————————————————————————————	Proposed Guardrail
Jurisdictional Streams Buffer Zone 1	Evicting Cable Guiderail
Buffer Zone 2	
Flow Arrow	
Disappearing Stream	
Spring — O	VEGETATION:
Wetland	single Tree 😌 😚
Proposed Lateral, Tail, Head Ditch ———— 🏼 🚬	Single Shrub
False Sump	Hedge
- also comp	Woods Line

RAILROADS:

Standard Gauge -

RR Abandoned -

RR Dismantled

Switch -

RR Signal Milepost —

CS# TRANSPORTATION O MULEPOST 35 SUITCH	Orchard ————————————————————————————————————	ලි ලි ලි ලි Vineyard
	EXISTING STRUCTURES:	
	MAJOR:	
	Bridge, Tunnel or Box Culvert	CONC
•	Bridge Wing Wall, Head Wall and End Wall-) CONC ## (
Δ	MINOR:	•
	Head and End Wall	CONC HW
	Pipe Culvert	
	Footbridge	
	Drainage Box: Catch Basin, DI or JB	CB
	Paved Ditch Gutter	
— 	Storm Sewer Manhole	9
— ——	Storm Sewer	s
——E——		

UTILITIES:

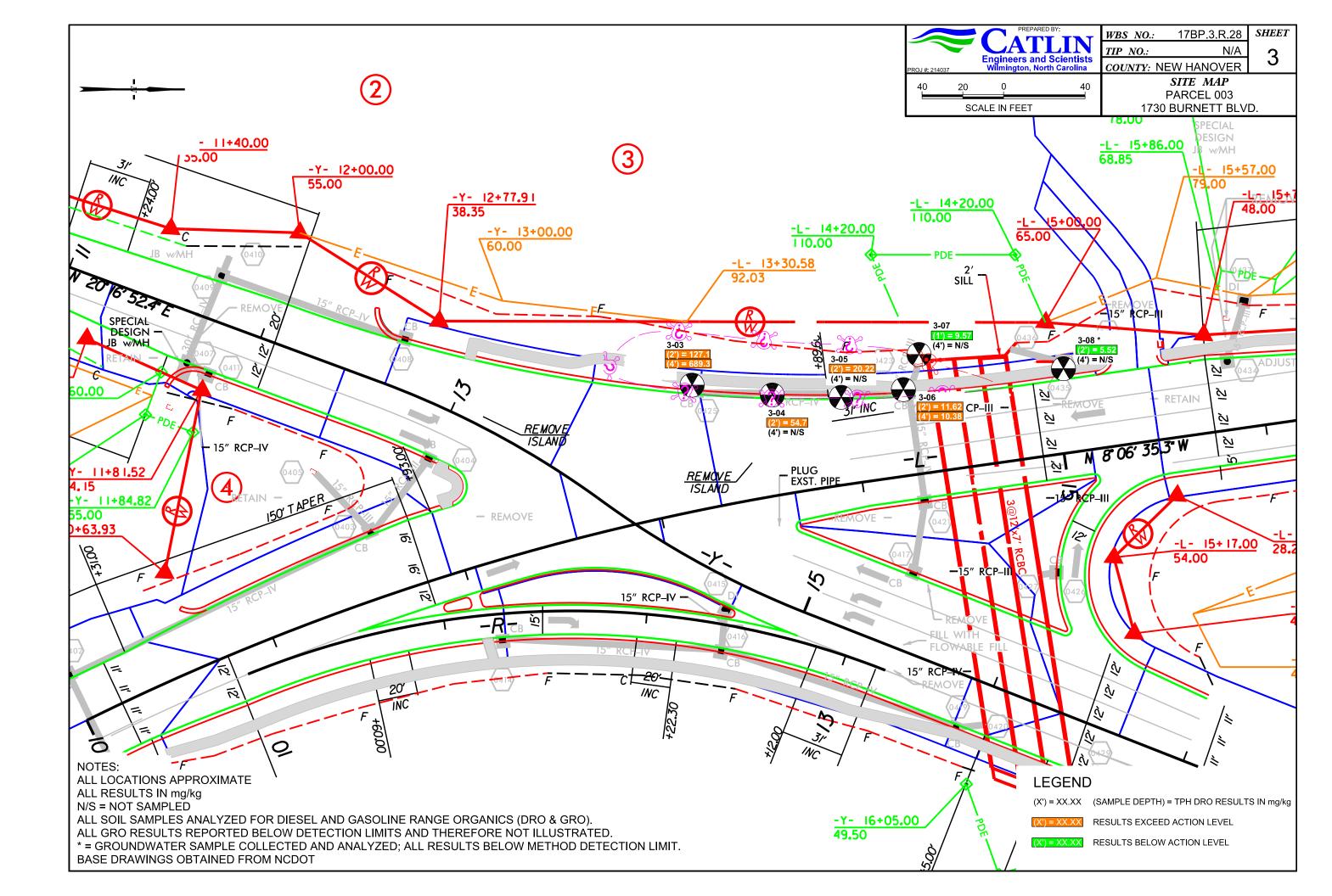
P

POWER:	
Existing Power Pole	•
Proposed Power Pole	6
Existing Joint Use Pole	-
Proposed Joint Use Pole	•
Power Manhole	ø
Power Line Tower	\boxtimes
Power Transformer	
U/G Power Cable Hand Hole	
H-Frame Pole	••
Recorded U/G Power Line	P
Designated U/G Power Line (S.U.E.*)	

TELEPHONE:

Existing Telephone Pole
Proposed Telephone Pole 🔶
Telephone Manhole 🛛 🕜
Telephone Booth 1
Telephone Pedestal 🔟
Telephone Cell Tower ———— 👗
U/G Telephone Cable Hand Hole 🔣
Recorded U/G Telephone Cable
Designated U/G Telephone Cable (S.U.E.*)
Recorded U/G Telephone Conduit
Designated U/G Telephone Conduit (S.U.E.*)
Recorded U/G Fiber Optics Cable
Designated U/G Fiber Optics Cable (S.U.E.*)

	PROJECT REFERENCE	E NO.	SHEET NO.
	17BP.3.R.28 (N/A)	2
		NABY	DI ANG
	PRELIMI DO NOT US	E FOR CONST	
	WATER:		
	Water Manhole		W
~	Water Meter		0
0	Water Valve		8
	Water Hydrant		
			•
	Recorded U/G Water Line		-•
	Designated U/G Water Line (S.U.E.*)		
	Above Ground Water Line	A/	3 Noter
	_ <i>.</i>		
	TV:		
	TV Satellite Dish —————		R
	TV Pedestal		C
<	TV Tower		\otimes
	U/G TV Cable Hand Hole		1
	Recorded U/G TV Cable		- 1y
	Designated U/G TV Cable (S.U.E.*)		-1y
	Recorded U/G Fiber Optic Cable		
_	Designated U/G Fiber Optic Cable (S.U.E.*)-		
	Designated 0/G Fiber Optic Cable (3.0.2.)		
	GAS:		
	Gas Valve		٥
	Gas Meter		0
	Recorded U/G Gas Line		-6
	Designated U/G Gas Line (S.U.E.*)		-6
			G Gos
	Above Ground Gas Line		
	SANITARY SEWER:		
	Sanitary Sewer Manhole		A
	Sanitary Sewer Cleanout		•
	U/G Sanitary Sewer Line		
			- 55
	Above Ground Sanitary Sewer		
	Recorded SS Forced Main Line		- 155
	Designated SS Forced Main Line (S.U.E.*) —		
	MISCELLANEOUS:		
	Utility Pole		•
	Utility Pole with Base		
	Utility Located Object		0
	Utility Traffic Signal Box		5
	Utility Unknown U/G Line		
	U/G Tank; Water, Gas, Oil		
_	Underground Storage Tank, Approx. Loc. —		
-	A/G Tank; Water, Gas, Oil		
	Geoenvironmental Boring		*
	U/G Test Hole (S.U.E.*) ————		•
_	Abandoned According to Utility Records —	A	ATUR
	End of Information	E	.O.I.



APPENDIX A

PYRAMID GEOPHYSICAL REPORT



GEOPHYSICAL SURVEY

PARCEL 003 – VICINITY OF FRONT STREET & BURNETT BOULEVARD NCDOT PROJECT WBS: 17BP.3.R.28

WILMINGTON, NEW HANOVER COUNTY, NC

MAY 12, 2014

Report prepared for:

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EXECUTIVE SUMMARY

Project Description: Pyramid Environmental conducted a geophysical investigation for Catlin Engineers & Scientists at NCDOT Parcel 3 located along Burnett Blvd. near the Front Street intersection in Wilmington, New Hanover County, NC. The survey was part of a North Carolina Department of Transportation (NCDOT) overhead rail line project (NCDOT Project WBS 17BP.3.R.28). Catlin Engineers & Scientists directed Pyramid as to the geophysical survey boundaries at the project site, which were designed to include the area between the existing edge of pavement and the NCDOT proposed ROW and/or easement. The geophysical investigation consisted of an electromagnetic (EM) induction-metal detection survey.

Geophysical Results: All of the EM61 anomalies detected could be attributed to known cultural features such as underground utilities and metal guy wires. The geophysical investigation <u>did not</u> record any evidence of metallic UST at the property.

INTRODUCTION

Pyramid Environmental conducted a geophysical investigation for Catlin Engineers & Scientists at NCDOT Parcel 3 located along Burnett Blvd. near the Front Street intersection in Wilmington, New Hanover County, NC. The survey was part of a North Carolina Department of Transportation (NCDOT) overhead rail line project (NCDOT Project WBS 17BP.3.R.28). Catlin Engineers & Scientists directed Pyramid as to the geophysical survey boundaries at the project site, which were designed to include the area between the existing edge of pavement and the NCDOT proposed ROW and/or easement. The survey grid spanned approximately 110 feet from north to south and a maximum of 40 feet from west to east, and included the accessible areas of the parcel. Conducted on May 9, 2014, the geophysical investigation was performed to determine if unknown, metallic underground storage tanks (USTs) were present beneath the survey area.

The consisted open grassy areas, heavily vegetated forest area, and a large drainage ditch. Dense tree cover was present on the south side of the parcel that was not accessible by the geophysical equipment. Similarly, the large drainage ditch and an area containing debris prevented the survey from extending to the north boundary of the parcel. Aerial photographs showing the survey area boundaries and ground-level photographs are shown in **Figure 1**.

FIELD METHODOLOGY

The geophysical investigation consisted of an electromagnetic (EM) induction-metal detection survey. Pyramid collected the EM data using a Geonics EM61 metal detector integrated with a Trimble AG-114 GPS antenna. 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 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 along north-south trending or east-west trending, generally 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 not required due to all anomalies being directly attributed to known cultural features (see discussions below).

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 that was observed along the entire east boundary of the survey area adjacent to the road was associated with a large natural gas pipeline located in this area. The pipeline had been marked by utility locators, and was also visible above ground. The EM response at northwest corner of the survey area was associated with a metal guy wire. The remaining survey area did not exhibit any significant EM responses that would be indicative of USTs or other structures.

The geophysical investigation did not record any evidence of metallic USTs at the property.

SUMMARY & CONCLUSIONS

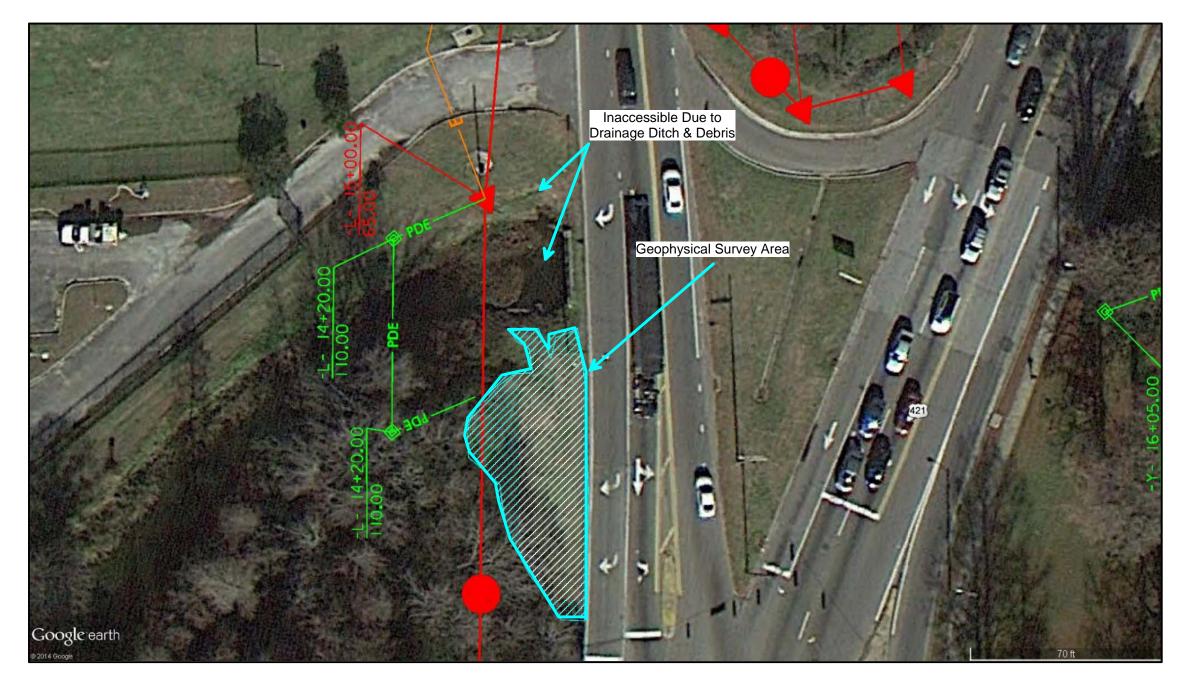
Our evaluation of the EM61 data collected at Parcel 003 along Burnett Blvd. in Wilmington, New Hanover County, North Carolina, provides the following summary and conclusions:

- The EM61 survey provided reliable results for the detection of metallic USTs within the accessible portions of the geophysical survey area.
- All of the EM61 anomalies detected could be attributed to known cultural features such as underground utilities and metal guy wires.
- The geophysical investigation <u>did not record any evidence of metallic UST at the</u> <u>property</u>.

LIMITATIONS

Geophysical surveys have been performed and this report prepared for Catlin Engineers & Scientists 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 Location of the Geophysical Survey Area With NCDOT Proposed ROW/Easement Overlay



View of Geophysical Survey Area (Facing Approximately North)

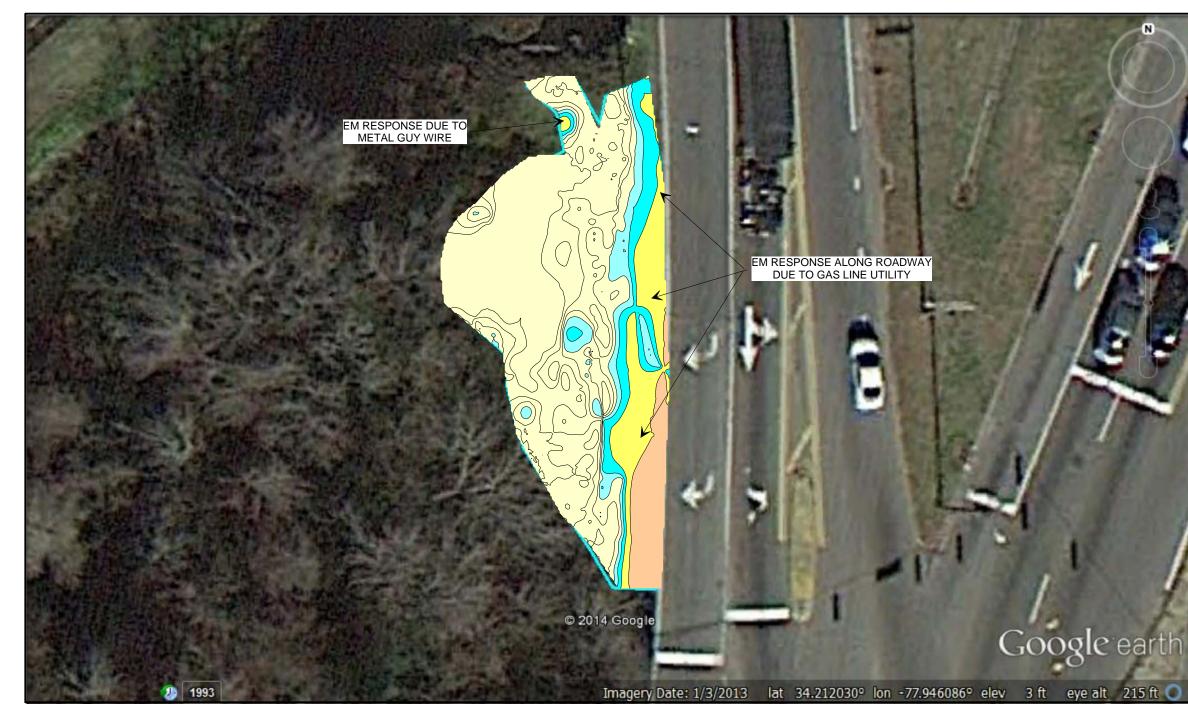


View of Inaccessible Drainage Area and Debris

	EL 003: RVEY BOUNDARIES IOTOGRAPHS
	URNETT BLVD. S 17BP.3.R.28) ANOVER COUNTY, NC
	503 INDUSTRIAL AVENUE GREENSBORO, NC 27460 36) 335-3174 (p) (336) 691-0648 (f) we # C1251 Eng. / License # C257 Geology
DATE 5/9/2014	CLIENT CATLIN ENGINEERS
PYRAMID 2014- PROJECT #:	FIGURE 1



EM61 Differential Results



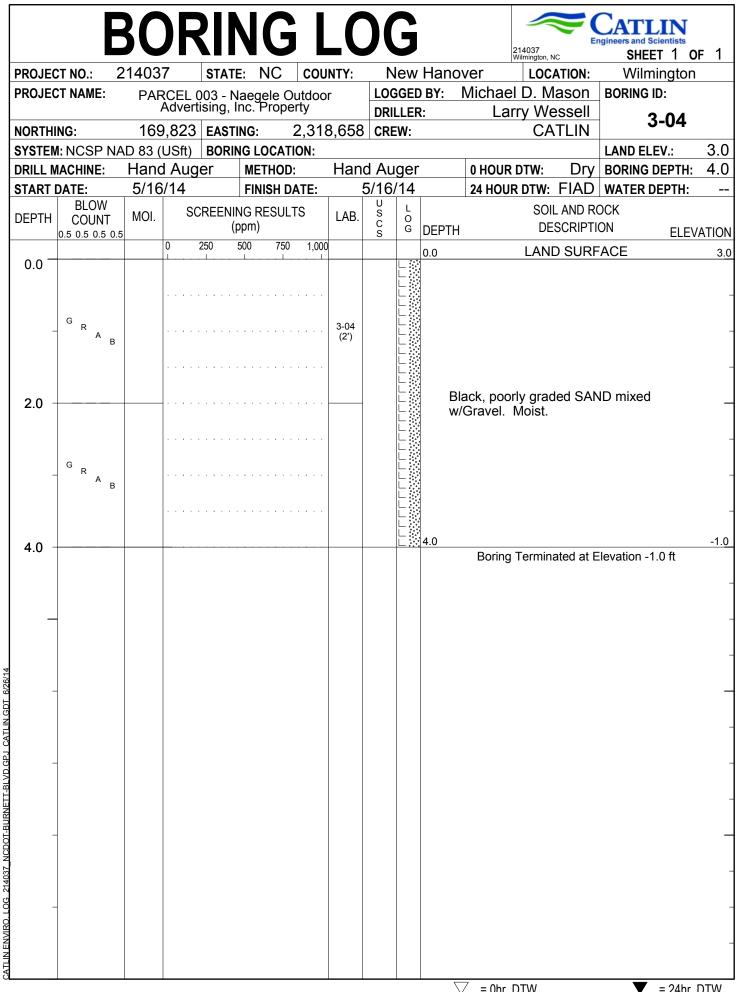
NO EVIDENCE OF METALLIC USTs OBSERVED

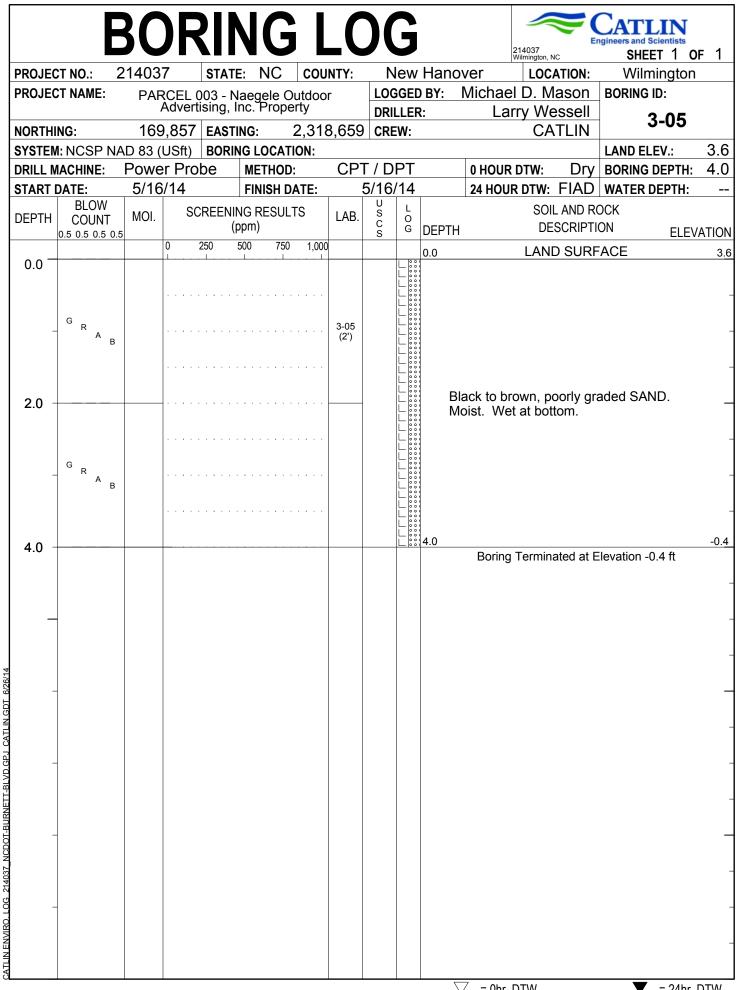
The contour plots show the differential results of the EM61instrument in millivolts (mV). The differential response focuses on larger, buried metallic objects such as drums and USTs and ignores smaller miscellaneous buried, metal debris. The EM61 data were collected on May 9, 2014, using a Geonics EM61 instrument. Ground penetrating radar (GPR) data were not required due to all anomalies being directly attributed to visible objects at ground surface or known utility lines. EM61 Metal Detection Response (millivolts) TITLE PARCEL 003: EM61 DIFFERENTIAL **RESULTS CONTOUR MAP** PROJECT FRONT ST. & BURNETT BLVD. (NCDOT WBS 17BP.3.R.28) WILMINGTON, NEW HANOVER COUNTY, NC 503 INDUSTRIAL AVENUE GREENSBORO, NC 27460 (336) 335-3174 (p) (336) 691-0648 (f) License # C1251 Eng. / License # C257 Geology PYRAMID ENVIRONMENTAL & ENGINEERING, P.C. CLIENT CATLIN ENGINEERS DATE 5/9/2014 PYRAMID 2014-FIGURE 2 PROJECT #:

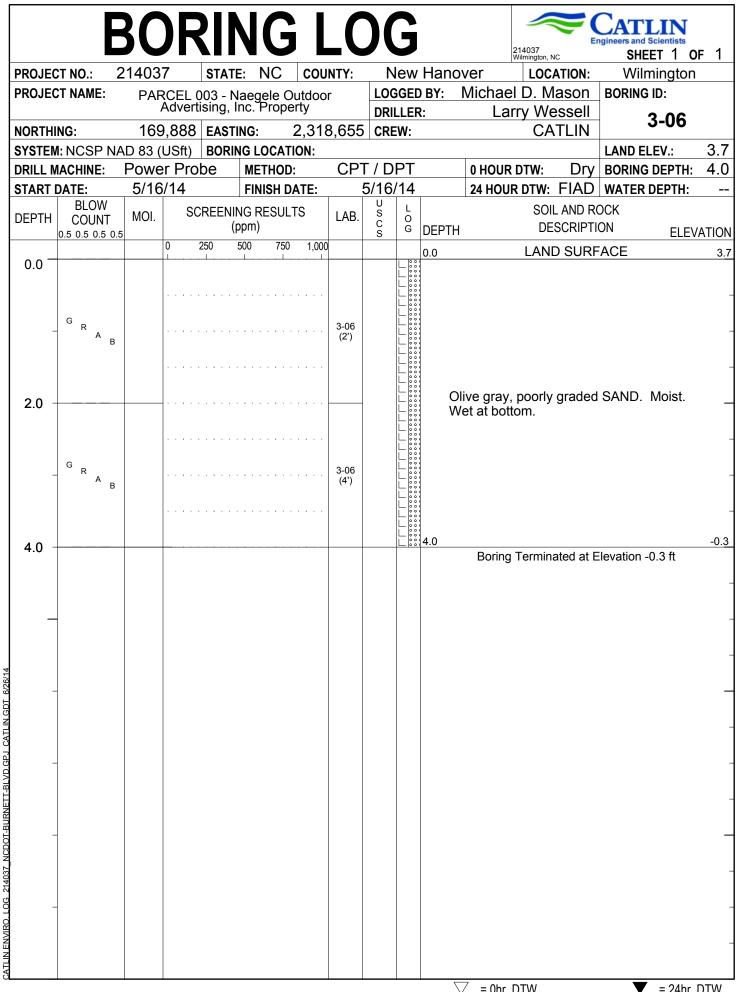
APPENDIX B

BORING LOGS

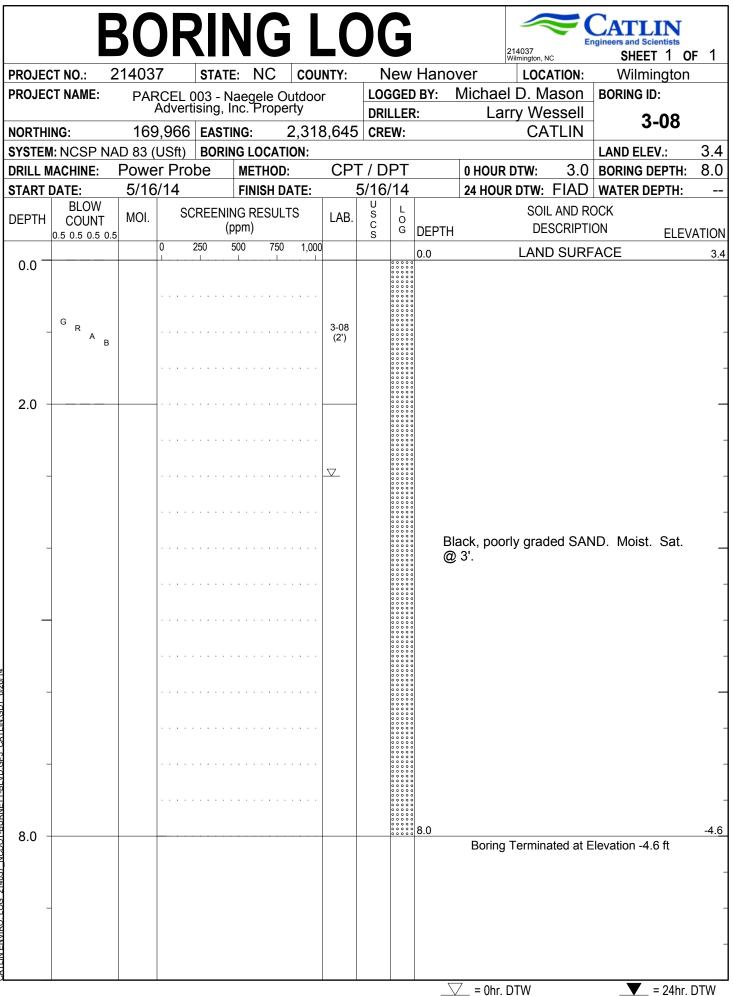
PROJE	CT NAME:	PAF	RCEL 00 Advertisi	3 - Naeg ng, Inc.	jele Oi Prope	utdoo rty	r		GEE LLEF) BY: R:	Michael D. M Larry We		BORING ID:	
NORTH	ING:	169),784 e	ASTING:	2	2,318	8,653	-				TLIN	3-03	
	I: NCSP NA												LAND ELEV.:	NN
	ACHINE:		d Auger		THOD:		Hand		_	•	0 HOUR DTW:			3.5
START	BLOW	5/16			NISH DA			5/16/ U	/14 L				WATER DEPTH:	-
DEPTH	COUNT 0.5 0.5 0.5 0.5	MOI.		EENING F (ppm)			LAB.	S C S	O G	DEPT		AND RO		VATIO
0.0 -			0 250) 500	750	1,000				0.0	LANE) SURF	ACE	
0.0										4 0 0				
				•••••						•				
-	G R						3-03				Black, poorly grad	ed SAN	ND w/Gravel.	
	A B						(2')			•	Moist.			
				• • • • • •						*				
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2.0														
	G			• • • • • •						4 9 9				
	R A B						3-03 (4')				S.A.A. w/HCO. W	/et.		
-										0 0 0				
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											Boring Terminated 3.5 ft du	e to obs	truction.	n
-	-													
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ROJEC	T NAME:	PAF	RCEL 00 Advertisi	3 - Ņa	egele	Outdoo	r	LOC	GEE	BY: I	Michael D. M		BORING ID:	
IORTHI	NG:		9,895 E				3,638		LLEF EW:	R:	Larry We	essell TLIN	3-07	
	I: NCSP NA			BORING			-,						LAND ELEV.:	3.0
	IACHINE:	,	d Auger		NETHO		Han	d Au	ger		0 HOUR DTW:	Dry	BORING DEPTH:	1.(
TART [DATE:	5/16	6/14	F	INISH	DATE:		5/16	_		24 HOUR DTW:	FIAD	WATER DEPTH:	-
EPTH	BLOW COUNT	MOI.	SCRI	EENING (ppi		ILTS	LAB.	U S C	L O G	DEPTH		L AND RO	ON	
	0.5 0.5 0.5 0.5		0 250			50 1,000)	S	0					
0.0										0.0	LANI	D SURF	ACE	3
	G R B						3-07 (1')			Bla	ack, TOPSOIL a	and orga	anic (grass).	
1.0										1.0	Poring Tormin	atad at [Elevation 2.0 ft	2
											Boring Termin			
_														
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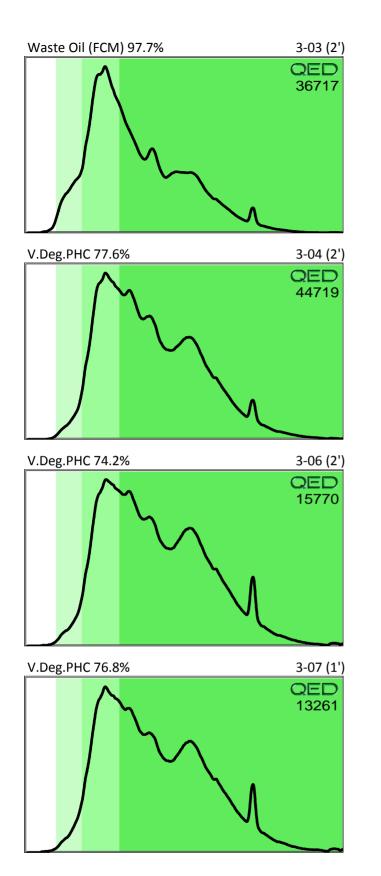
CATLIN ENVIRO. LOG 214037 NCDOT-BURNETT-BLVD.GPJ CATLIN.GDT

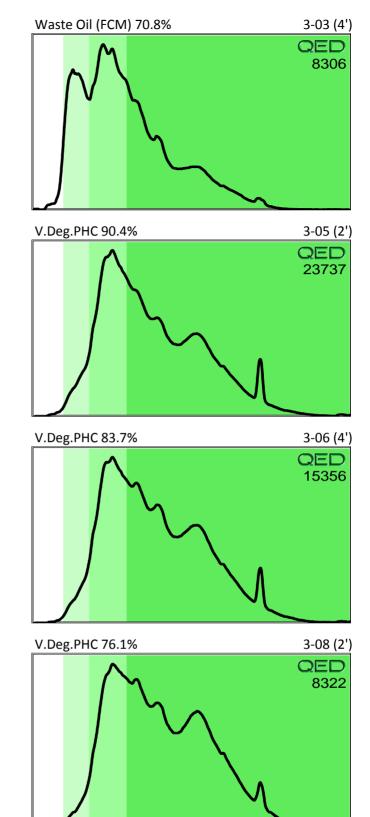
__ = 24hr. DTW

APPENDIX C

QROS QED™ REPORT

Q	ED												QROS
				Hydroca	arbon An	alysis R	esults						
	Catlin Wilmington, NC								Sar Sample Sampl		racted		Friday, May 16, 2014 Friday, May 16, 2014 Monday, May 19, 2014
Contact: Ben Ashba Derator Rachel Menoher													
•	CATLIN Project No. 214037 Sample ID	Sample ID Dilution BTEX GRO DRO TPH Total 16 EPA BaP Rational contraction				Ratios HC Fingerprint Ma		HC Fingerprint Match					
			(,	(,	(,	(,	(C10-C35)			% light	% mid	% heavy	
S	3-03 (2')	21.0	<1	<1	127.1	127.1	35.19	2.18	<0.021	57.9	37.4	4.7	Waste Oil (FCM) 97.7%
S	3-03 (4')	256.0		<12.8			161.2	5.67	<0.256		12.1		Waste Oil (FCM) 70.8%
	3-04 (2')	20.0		<1	54.7	54.7	40.72	2.09	0.037	27.3	58.5		V.Deg.PHC 77.6%
	3-05 (2')	19.0		<1	20.22	20.22		1.02	< 0.019	51.6	39.8	-	V.Deg.PHC 90.4%
	3-06 (2')	22.0		<1.1	11.62		10.65	0.71	< 0.022	36.9	47.5		V.Deg.PHC 74.2%
	3-06 (4') 3-07 (1')	19.0 22.0		<0.9 <1.1		10.38 9.57	10.35 8.77	0.64 0.58	<0.019 <0.022	50.8 40.9	38.1 44.1		V.Deg.PHC 83.7% V.Deg.PHC 76.8%
	3-08 (2')	22.0		<1.1	9.57 5.52	5.52	5.06	0.34	<0.022	40.9	37.3		V.Deg.PHC 76.1%
-	(-)				0.02	0.02	0.00	5.01					
	Initial (Calibrator	QC check	OK					Final FC	CM QC	Check	OK	98.3
ingerprints	erated by a QED HC-1 analyser. Concer provide a tentative hydrocarbon identificat 3S) = Site Specific or Library Background S	ion. The abbi	reviations ar	e:- FCM = R	esults calcula	ted using Fur	ndamental Calil	oration Mod	e : % = conf				







Sample Collection Sample ID TAT Requested QED UVF Date Time Initials 24 Hour 48 Hour MM 1350 3-03(2) 5.16.14 MM 1355 3-03(41) mm 3-04(2) 1325 im 3.05(2) 1320 1300 MM 3-06 (21) mm 3-06 (4) 130 imm 1250 071 Mmi VČ 1205 Q 2.

Chain of Custody Record and Analytical Request Form

Client: CATLW
Contact: Ben Ashba
Phone: 910-452-5861
Email: benashbaccattin Usaco
Project Reference:
214037
PO#140519-1
Each Sample will be analyzed for total

BTEX, GRO, DRO, TPH, and PAH

Each Sample will generate a fingerprint representative of the petroleum product within the sample. Electronic Data will be submitted to the email above.

		,				
Relinquished by	Date/time	Accepted by	Date/time			
Muhallun	5/19/14 840	Brack	5-19.14 840			
Relinquished by	Date/time	Accepted by	Date/time			
Bala	5-19.14 840	Ban	5/19/14 9:30			
Relinguished by	Date/time	Accepted by	Date/time			

SHIP TO: QROS
420 Raleigh Street Suite E
Wilmington, NC 28412
Rachel Menoher- rachelm@grosllc.com
910-520-2902

APPENDIX D

PACE ANALYTICAL SERVICES, INC. GROUNDWATER SAMPLE ANALYTICAL RESULTS AND CHAIN OF CUSTODY DOCUMENTATION



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

June 04, 2014

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

RE: Project: NCDOT FRONT ST. WBS17BP.3.R.28 Pace Project No.: 92202425

Dear Chemical Engineer:

Enclosed are the analytical results for sample(s) received by the laboratory on May 22, 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,

angela M. Baioni

Angela Baioni angela.baioni@pacelabs.com Project Manager

Enclosures

cc: Ben Ashba, NCDOT South East Sean O'Neil, NCDOT South East





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

CERTIFICATIONS

Project: NCDOT FRONT ST. WBS17BP.3.R.28

Pace Project No.: 92202425

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 SUMMARY

Project: NCDOT FRONT ST. WBS17BP.3.R.28

Pace Project No.: 92202425

Lab ID	Sample ID	Matrix	Date Collected	Date Received
92202425001	3-08	Water	05/16/14 12:25	05/22/14 09:30



SAMPLE ANALYTE COUNT

Project:NCDOT FRONT ST. WBS17BP.3.R.28Pace Project No.:92202425

Lab ID	Sample ID	Method	Analysts	Analytes Reported	Laboratory
92202425001	3-08	EPA 625	RES	58	PASI-C
		SM 6200B	CAH	63	PASI-C



ANALYTICAL RESULTS

Project: NCDOT FRONT ST. WBS17BP.3.R.28

Pace Project No.: 92202425

Sample: 3-08	Lab ID:	92202425001	Collected:	05/16/1	4 12:25	Received: 05/	22/14 09:30 Ma	atrix: Water	
Parameters	Results	Units	Report Limit	MDL	DF	Prepared	Analyzed	CAS No.	Qual
625 MSSV	Analytical	Method: EPA 6	25 Preparat	ion Metho	d: EPA (625			
Acenaphthene	ND u	ıg/L	5.0	0.25	1	05/23/14 11:30	06/02/14 22:36	83-32-9	
Acenaphthylene	ND u	ıg/L	5.0	0.21	1	05/23/14 11:30	06/02/14 22:36	208-96-8	
Anthracene	ND u	ig/L	5.0	0.14	1	05/23/14 11:30	06/02/14 22:36	120-12-7	
Benzo(a)anthracene	ND u	ig/L	5.0	0.33	1	05/23/14 11:30	06/02/14 22:36	56-55-3	
Benzo(a)pyrene	ND u	-	5.0	0.30	1	05/23/14 11:30	06/02/14 22:36	50-32-8	
Benzo(b)fluoranthene	ND u	ig/L	5.0	0.28	1	05/23/14 11:30	06/02/14 22:36	205-99-2	
Benzo(g,h,i)perylene	ND u	-	5.0	0.38	1	05/23/14 11:30	06/02/14 22:36	191-24-2	
Benzo(k)fluoranthene	ND u	0	5.0	0.43	1	05/23/14 11:30	06/02/14 22:36	207-08-9	
4-Bromophenylphenyl ether	ND u	-	5.0	0.82	1	05/23/14 11:30	06/02/14 22:36	101-55-3	
Butylbenzylphthalate	ND u	-	5.0	0.79	1	05/23/14 11:30	06/02/14 22:36	85-68-7	
4-Chloro-3-methylphenol	ND u	-	5.0	3.7	1	05/23/14 11:30	06/02/14 22:36		
bis(2-Chloroethoxy)methane	ND u	-	10.0	0.92	1	05/23/14 11:30	06/02/14 22:36		
bis(2-Chloroethyl) ether	ND u	-	5.0	1.0	1	05/23/14 11:30	06/02/14 22:36		
bis(2-Chloroisopropyl) ether	ND u		5.0	0.95	1	05/23/14 11:30	06/02/14 22:36		
2-Chloronaphthalene	ND u	-	5.0	0.98	1	05/23/14 11:30	06/02/14 22:36		
2-Chlorophenol	ND u	0	5.0	1.3	1	05/23/14 11:30	06/02/14 22:36		
4-Chlorophenylphenyl ether	ND u	-	5.0	0.87	1	05/23/14 11:30	06/02/14 22:36		
Chrysene	ND u		5.0	0.21	1	05/23/14 11:30	06/02/14 22:36		
Dibenz(a,h)anthracene	ND u	-	5.0	0.55	1	05/23/14 11:30	06/02/14 22:36		
3,3'-Dichlorobenzidine	ND u	-	25.0	2.1	1	05/23/14 11:30	06/02/14 22:36		
2,4-Dichlorophenol	ND u	0	5.0	1.7	1	05/23/14 11:30	06/02/14 22:36		
Diethylphthalate	ND u	-	5.0 5.0	0.58	1	05/23/14 11:30	06/02/14 22:30		
2,4-Dimethylphenol	ND u	0	10.0	1.2	1	05/23/14 11:30	06/02/14 22:30		
Dimethylphthalate	ND u		5.0	0.76	1	05/23/14 11:30	06/02/14 22:30		
Di-n-butylphthalate	ND u	-	5.0 5.0	0.75	1	05/23/14 11:30	06/02/14 22:30		
4,6-Dinitro-2-methylphenol		-	20.0	2.6	1	05/23/14 11:30	06/02/14 22:36		
	ND u ND u	-	20.0 50.0	2.0 9.0	1	05/23/14 11:30	06/02/14 22:36		
2,4-Dinitrophenol 2,4-Dinitrotoluene		0	5.0	9.0 0.90	1	05/23/14 11:30	06/02/14 22:36		
	ND u								
2,6-Dinitrotoluene	ND u	-	5.0 5.0	0.98	1	05/23/14 11:30	06/02/14 22:36		
Di-n-octylphthalate	ND u	-		0.66	1	05/23/14 11:30	06/02/14 22:36		
bis(2-Ethylhexyl)phthalate	ND u	-	5.0	0.79	1	05/23/14 11:30	06/02/14 22:36		
Fluoranthene	ND u	0	5.0	0.21	1	05/23/14 11:30	06/02/14 22:36		
Fluorene	ND u		5.0	0.21	1	05/23/14 11:30	06/02/14 22:36		
Hexachloro-1,3-butadiene	ND u	-	5.0	0.94	1	05/23/14 11:30	06/02/14 22:36		
Hexachlorobenzene	ND u	•	5.0	0.72	1	05/23/14 11:30	06/02/14 22:36		
Hexachlorocyclopentadiene	ND u	-	10.0	0.88	1	05/23/14 11:30	06/02/14 22:36		
Hexachloroethane	ND u	-	5.0	1.1	1	05/23/14 11:30	06/02/14 22:36		
Indeno(1,2,3-cd)pyrene	ND u		5.0	0.29	1	05/23/14 11:30	06/02/14 22:36		
Isophorone	ND u	-	10.0	0.89	1	05/23/14 11:30	06/02/14 22:36		
Naphthalene	ND u	-	5.0	0.34	1	05/23/14 11:30	06/02/14 22:36		
Nitrobenzene	ND u	-	5.0	1.1	1	05/23/14 11:30	06/02/14 22:36		
2-Nitrophenol	ND u	-	5.0	0.91	1	05/23/14 11:30	06/02/14 22:36		
4-Nitrophenol	ND u	-	50.0	4.1	1	05/23/14 11:30	06/02/14 22:36		
N-Nitrosodimethylamine	ND u	-	5.0	0.91	1	05/23/14 11:30	06/02/14 22:36		
N-Nitroso-di-n-propylamine	ND u	-	5.0	0.99	1	05/23/14 11:30	06/02/14 22:36	621-64-7	
N-Nitrosodiphenylamine	ND u	ıg/L	10.0	1.0	1	05/23/14 11:30	06/02/14 22:36	86-30-6	



ANALYTICAL RESULTS

Project: NCDOT FRONT ST. WBS17BP.3.R.28

Pace Project No.:

o.: 92202425

Parameters Results Units Imit MDL DF Prepared Analyzed CAS No. Qual 625 MSV Analytical Method: EPA 625 Preparation Method: EPA 625	Sample: 3-08	Lab ID: 92	202425001	Collected:	05/16/14	12:25	Received: 05/	22/14 09:30 Ma	atrix: Water	
G25 MSSV Analytical Method: EPA 625 Preparation Method: EPA 625 Pentachlorophenol ND ug/L 5.0 0.22 1 06/23/14 11:30 06/02/14 22:36 87-86-5 Phenol ND ug/L 5.0 0.22 1 06/23/14 11:30 06/02/14 22:36 108-95-2 Pyrene ND ug/L 5.0 0.19 1 05/23/14 11:30 06/02/14 22:36 12.96-17 2,46-Trichlorophenol ND ug/L 5.0 0.98 1 05/23/14 11:30 06/02/14 22:36 12.96-22 Surrogates NU 10.120 1 05/23/14 11:30 06/02/14 22:36 18-06-2 Surrogates NU 10.120 1 05/23/14 11:30 06/02/14 22:36 18-16-0 Phenol-d6 (S) 74 % 15-120 1 05/23/14 11:30 06/02/14 22:36 1312-88-3 2-Fluorophenol (S) 34 % 10-120 1 05/23/14 11:30 06/02/14 22:36 1312-88-3 2-Ale-Tribromophenol (S) 34 % 10-120 1 05/23/14 11:			R	eport						
Pentachlorophenol ND ug/L 10. 4.6 1 05/23/14 11:30 06/02/14 22:36 87-86-5 Phenol ND ug/L 5.0 0.22 1 05/23/14 11:30 06/02/14 22:36 10.89-22 Pyrene ND ug/L 5.0 0.98 1 05/23/14 11:30 06/02/14 22:36 12.9-0-0 1,2,4-Trichlorobenzene ND ug/L 5.0 0.98 1 05/23/14 11:30 06/02/14 22:36 12.9-0-0 2,4-Grichlorophenol ND ug/L 1.0 1.3 1 05/23/14 11:30 06/02/14 22:36 88-06-2 Surrogates NU 10.120 1 05/23/14 11:30 06/02/14 22:36 321-60-8 TerphenyL'G14 (S) 71 % 11-131 1 05/23/14 11:30 06/02/14 22:36 321-60-8 2-Fluorobphenol (S) 34 % 10-120 1 05/23/14 11:30 06/02/14 22:36 367-12-4 2-Fluorobphenol (S) 73 % 10-120 1 05/23/14 11:30 06/02/14 22:36 367-12-4 2-Fluorobphenol (S) 73 % <	Parameters	Results	Units L	_imit	MDL	DF	Prepared	Analyzed	CAS No.	Qual
Phenalthrene ND ug/L 5.0 0.22 1 05/02/14 11:30 06/02/14 22:36 129-00-0 Pyene ND ug/L 5.0 0.19 1 05/23/14 11:30 06/02/14 22:36 120-00-0 1.2.4-Trichlorobenzene ND ug/L 5.0 0.98 1 05/23/14 11:30 06/02/14 22:36 120-08-1 2.4-G.Trichlorobenzene ND ug/L 1.0 1 05/23/14 11:30 06/02/14 22:36 120-82-1 2.4-G.Trichlorobenzene ND ug/L 1.0 1 05/23/14 11:30 06/02/14 22:36 120-82-1 2.4-G.Trichlorobenzene/G(S) 74 % 15-120 1 05/23/14 11:30 06/02/14 22:36 171-85-10 Phenold G(S) 2.6 % 10-120 1 05/23/14 11:30 06/02/14 22:36 137-78-3 2-Fluorobjenenl (S) 74 % 11-131 1 05/23/14 11:30 06/02/14 22:36 137-78-3 2-Fluorobjenenl (S) 73 % 10-130 1 05/23/14 11:30 06/02/14 22:36 137-78-3 2-Fluorobjenenl (S) 73 % 10-150 </td <td>625 MSSV</td> <td>Analytical Me</td> <td>thod: EPA 625</td> <td>Preparati</td> <td>on Metho</td> <td>d: EPA 6</td> <td>625</td> <td></td> <td></td> <td></td>	625 MSSV	Analytical Me	thod: EPA 625	Preparati	on Metho	d: EPA 6	625			
Phenol ND ug/L 5.0 1.9 1 05/23/14 11:30 06/02/14 22:36 108-95-2 Pyrene ND ug/L 5.0 0.19 1 05/23/14 11:30 06/02/14 22:36 129-00-0 2.4.6 Trichlorophenol ND ug/L 10.0 1.3 1 05/23/14 11:30 06/02/14 22:36 82-08-1 Surrogates 1 05/23/14 11:30 06/02/14 22:36 16-60-0 2-Fluorobiphenyl (S) 74 % 15-120 1 05/23/14 11:30 06/02/14 22:36 1316-51-0 Prenovle14 (S) 71 % 11-131 1 05/23/14 11:30 06/02/14 22:36 1316-51-0 Phenole16 (S) 26 % 10-120 1 05/23/14 11:30 06/02/14 22:36 1316-51-0 Phenole16 (S) 26 % 10-120 1 05/23/14 11:30 06/02/14 22:36 1316-51-0 Phenole16 (S) 26 % 10-120 1 05/23/14 11:30 06/02/14 22:36 1316-51-0 2.4.6-Tribromophenol (S) 26 % 10-120 1 05/23/14 11:30 06/02/14 22:30 74-32	Pentachlorophenol	ND ug/L		10.0	4.6	1	05/23/14 11:30	06/02/14 22:36	87-86-5	
PyreneND ug/L5.00.19105/23/14 11:3006/02/14 22:36129-00-1.2,4-TichlorobenolND ug/L5.00.98105/23/14 11:3006/02/14 22:3680-62Surrogates105/23/14 11:3006/02/14 22:368165-60-02-Fluorobiphenyl (S)74 %15-120105/23/14 11:3006/02/14 22:36816-68-02-Fluorobiphenyl (S)74 %15-120105/23/14 11:3006/02/14 22:36321-60-8Terphenyl-d14 (S)71 %11-131105/23/14 11:3006/02/14 22:36312-78-8-32-Fluorophenol (S)34 %10-120105/23/14 11:3006/02/14 22:36312-78-8-32-Fluorophenol (S)34 %10-120105/23/14 11:3006/02/14 22:36312-78-8-32-Fluorophenol (S)34 %10-120105/23/14 11:3006/02/14 22:3636-71-242,4,6-Tirbiromophenol (S)73 %10-137105/23/14 11:3006/02/14 22:36131-79-6EoromobenzeneND ug/L0.500.25105/24/14 22:0971-43-2BromochicoromethaneND ug/L0.500.25105/24/14 22:0971-43-2BromochichoromethaneND ug/L0.500.25105/24/14 22:0975-27-4BromochichoromethaneND ug/L0.500.25105/24/14 22:0975-27-4BromochichoromethaneND ug/L0.500.25105/24/14 22:0975-27-4Bromochichoro	Phenanthrene	ND ug/L		5.0	0.22	1	05/23/14 11:30	06/02/14 22:36	85-01-8	
1,2,4-Trichlorobenzene ND ug/L 5.0 0.98 1 05/23/14 11:30 06/02/14 22:36 120-82:1 2,4,6-Trichlorophenol ND ug/L 10.0 1.3 1 05/23/14 11:30 06/02/14 22:36 88-06-2 Nitrobenzene-d5 (S) 70 % 10-120 1 05/23/14 11:30 06/02/14 22:36 216-06-0 2-Fluorobjehenyl (S) 71 % 11-131 1 05/23/14 11:30 06/02/14 22:36 3127-88-3 2-Fluorobjehenyl (S) 34 % 10-120 1 05/23/14 11:30 06/02/14 22:36 3127-88-3 2-Fluorobjehenol (S) 34 % 10-120 1 05/23/14 11:30 06/02/14 22:36 3127-88-3 2-Fluorobjehenol (S) 73 % 10-137 1 05/23/14 11:30 06/02/14 22:30 71-8-2 2-Fluorobjehenol (S) 73 % 10-137 1 05/23/14 11:30 06/02/14 22:00 71-8-2 Eoromolichloromethane ND ug/L 0.50 0.25 1 05/24/14 22:00 74-8-3 Bromochichoromethane ND ug/L 0.50 0.25 1 05/24/14 22:00 75-25-2 Bromochichorome	Phenol	ND ug/L		5.0	1.9	1	05/23/14 11:30	06/02/14 22:36	108-95-2	
2,4,6.Trichlorophenol ND ug/L 10.0 1.3 1 05/23/14 11:30 06/02/14 22:36 88-06-2 Surrogates 0 05/23/14 11:30 06/02/14 22:36 1415-00 2-Fluorobiphenyl (S) 74 % 15-120 1 05/23/14 11:30 06/02/14 22:36 1312-88-3 Terphenyl-d14 (S) 74 % 10-120 1 05/23/14 11:30 06/02/14 22:36 1312-88-3 2-Fluorophenol (S) 34 % 10-120 1 05/23/14 11:30 06/02/14 22:36 1312-88-3 2-Fluorophenol (S) 34 % 10-120 1 05/23/14 11:30 06/02/14 22:36 1312-88-3 2-Fluorophenol (S) 34 % 10-120 1 05/23/14 11:30 06/02/14 22:36 1312-88-3 2-Fluorophenol (S) 34 % 10-120 1 05/23/14 11:30 06/02/14 22:36 1312-88-3 2-Fluorophenol (S) 73 % 10-137 1 05/24/14 22:09 71-43-2 Bromocharone ND ug/L 0.50 0.25 1 05/24/14 22:09 74-32 <t< td=""><td>Pyrene</td><td>ND ug/L</td><td></td><td>5.0</td><td>0.19</td><td>1</td><td>05/23/14 11:30</td><td>06/02/14 22:36</td><td>129-00-0</td><td></td></t<>	Pyrene	ND ug/L		5.0	0.19	1	05/23/14 11:30	06/02/14 22:36	129-00-0	
Surrogates Normal Surrogates Nitrobenzene-d5 (S) 70 % 10-120 1 05/23/14 11:30 06/02/14 22:36 321-60-8 2-Fluorobiphenyl (S) 71 % 11-131 1 05/23/14 11:30 06/02/14 22:36 13127-88-3 2-Fluorobiphenol (S) 28 % 10-120 1 05/23/14 11:30 06/02/14 22:36 367-12-4 2,4,6-Tribromophenol (S) 34 % 10-120 1 05/23/14 11:30 06/02/14 22:36 367-12-4 2,4,6-Tribromophenol (S) 73 % 10-137 1 05/23/14 11:30 06/02/14 22:36 367-12-4 2,4,6-Tribromophenol (S) 73 % 10-137 1 05/23/14 11:30 06/02/14 22:30 367-12-4 2,4,6-Tribromophenol (S) 73 % 10-137 1 05/24/14 22:09 71-43-2 Bromobenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 74-3-2 Bromodichoromethane ND ug/L 0.50 0.25 1 05/24/14 22:09 75-27-4 Bromodichloromethane ND ug/L 0.50	1,2,4-Trichlorobenzene	ND ug/L		5.0	0.98	1	05/23/14 11:30	06/02/14 22:36	120-82-1	
Nitrobizenen-d5 (S) 70 % 10-120 1 05/23/14 11:30 06/02/14 22:36 1465-60-0 2-Fluorobiphenyl (S) 74 % 15-120 1 05/23/14 11:30 06/02/14 22:36 731-81-10 Phenol-d6 (S) 26 % 10-120 1 05/23/14 11:30 06/02/14 22:36 171-88-3 2-Fluorophenol (S) 34 % 10-120 1 05/23/14 11:30 06/02/14 22:36 171-88-3 2-Fluorophenol (S) 34 % 10-120 1 05/23/14 11:30 06/02/14 22:36 171-84-3 2-Fluorophenol (S) 73 % 10-137 1 05/23/14 11:30 06/02/14 22:09 161-72-4 2.4.6-Tibromphenol (S) 73 % 10-137 1 05/23/14 11:30 06/02/14 22:09 161-72-4 2.4.6-Tibromphenol (S) 73 % 10-137 1 05/24/14 22:09 174-32 Bromodichloromethane ND ug/L 0.50 0.25 1 05/24/14 22:09 168-86-1 Bromodichloromethane ND ug/L 0.50 0.25 1 05/24/14 22:09 75-25-2<	2,4,6-Trichlorophenol	ND ug/L		10.0	1.3	1	05/23/14 11:30	06/02/14 22:36	88-06-2	
2-Fluorobiphenyl (S) 74 % 15-120 1 05/23/14 11:30 06/02/14 22:36 321-60-8 Terphenyl-df4 (S) 71 % 11-131 1 05/23/14 11:30 06/02/14 22:36 1718-51-0 Phenol-d6 (S) 26 % 10-120 1 05/23/14 11:30 06/02/14 22:36 1718-78-3 2-Fluorophenol (S) 34 % 10-120 1 05/23/14 11:30 06/02/14 22:36 187-78-8-3 2-Fluorophenol (S) 73 % 10-137 1 05/23/14 11:30 06/02/14 22:36 187-76 6200B MSV Analytical Method: SM 6200F 1 05/23/14 12:09 71-43-2 Bromochloromethane ND ug/L 0.50 0.25 1 05/24/14 22:09 74-87-5 Bromochloromethane ND ug/L 0.50 0.25 1 05/24/14 22:09 74-83-9 Bromochloromethane ND ug/L 0.50 0.25 1 05/24/14 22:09 74-83-9 Bromochloromethane ND ug/L 0.50 0.25 1 05/24/14 22:09 75-2-2	Surrogates									
Terphenyl-d14 (S) 71 % 11-131 1 05/23/14 11:30 06/02/14 22:36 1718-51-0 Phenol-d6 (S) 26 % 10-120 1 05/23/14 11:30 06/02/14 22:36 13127-88-3 2.Fluorophenol (S) 73 % 10-137 1 05/23/14 11:30 06/02/14 22:36 367-12-4 2.4.6-Tribromophenol (S) 73 % 10-137 1 05/23/14 11:30 06/02/14 22:36 371-24 2.4.6-Tribromophenol (S) 73 % 10-137 1 05/23/14 11:30 06/02/14 22:36 371-24 2.4.6-Tribromophenol (S) 73 % 10-137 1 05/23/14 11:30 06/02/14 22:09 71-43-2 Bromobenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 71-43-2 Bromochloromethane ND ug/L 0.50 0.25 1 05/24/14 22:09 75-27-4 Bromomethane ND ug/L 0.50 0.25 1 05/24/14 22:09 74-83-9 n-Butylbenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 74-83-9 sco-Butylbenzene ND ug/L 0.50 0.25 1	Nitrobenzene-d5 (S)	70 %	1	10-120		1		06/02/14 22:36	4165-60-0	
Phenol-d6 (S) 26 % 10-120 1 05/23/14 11:30 06/02/14 22:36 13127-88-3 2-Fluorophenol (S) 34 % 10-120 1 05/23/14 11:30 06/02/14 22:36 367-12-4 2,4,6-Tribromophenol (S) 73 % 10-137 1 05/23/14 11:30 06/02/14 22:36 367-12-4 6200B MSV Analytical Method: SM 6200E E E E Benzene ND ug/L 0.50 0.25 1 05/24/14 22:09 71-43-2 Bromobenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 74-97-5 Bromodichloromethane ND ug/L 0.50 0.25 1 05/24/14 22:09 75-27-4 Bromodichloromethane ND ug/L 0.50 0.25 1 05/24/14 22:09 74-83-9 Bromotom ND ug/L 0.50 0.25 1 05/24/14 22:09 74-83-9 Bromotom ND ug/L 0.50 0.25 1 05/24/14 22:09 74-83-9 Bromotom ND ug/L 0.50 0.	2-Fluorobiphenyl (S)		1	15-120		1	05/23/14 11:30	06/02/14 22:36	321-60-8	
2-Fluorophenol (S) 34 % 10-120 1 05/23/14 11:30 06/02/14 22:36 367-12-4 2,4,6-Tribromophenol (S) 73 % 10-137 1 05/23/14 11:30 06/02/14 22:36 367-12-4 6200B MSV Analytical Method: SM 6200E 5 1 05/24/14 22:09 71-43-2 Bromobenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 71-43-2 Bromobenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 75-27-4 Bromoformomthane ND ug/L 0.50 0.25 1 05/24/14 22:09 75-27-4 Bromoform ND ug/L 0.50 0.25 1 05/24/14 22:09 75-27-4 Bromoform ND ug/L 0.50 0.25 1 05/24/14 22:09 75-25-2 Bromothane ND ug/L 0.50 0.25 1 05/24/14 22:09 14-36-3 sec-Butylbenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 14-36-3 Carbon tetrachloride ND ug/L 0.50 0.25 1 05/24/14 22:09 16-3-3		71 %		11-131		1	05/23/14 11:30	06/02/14 22:36	1718-51-0	
2,4,6-Tribromophenol (S) 73 % 10-137 1 05/23/14 11:30 06/02/14 22:36 118-79-6 6200B MSV Analytical Method: SM 62008 Benzene ND ug/L 0.50 0.25 1 05/24/14 22:09 71-3-2 Bromobenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 74-3-2 Bromodichloromethane ND ug/L 0.50 0.25 1 05/24/14 22:09 74-37-5 Bromodichloromethane ND ug/L 0.50 0.25 1 05/24/14 22:09 75-27-4 Bromodichloromethane ND ug/L 0.50 0.25 1 05/24/14 22:09 75-27-4 Bromodichloromethane ND ug/L 0.50 0.25 1 05/24/14 22:09 74-83-9 n-Butylbenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 74-83-9 c-Tarburghenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 74-83-9 c-Tarburghenzene ND ug/L 0.50 0.25 1 05/24/14 22	Phenol-d6 (S)		1	10-120		1	05/23/14 11:30	06/02/14 22:36	13127-88-3	
6200B MSV Analytical Method: SM 62008 Benzene ND ug/L 0.50 0.25 1 05/24/14 22:09 71-43-2 Bromobenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 74-97-5 Bromodichloromethane ND ug/L 0.50 0.25 1 05/24/14 22:09 74-97-5 Bromodichloromethane ND ug/L 0.50 0.25 1 05/24/14 22:09 75-27-4 Bromoform ND ug/L 0.50 0.25 1 05/24/14 22:09 75-27-8 Bromoform ND ug/L 0.50 0.25 1 05/24/14 22:09 74-83-9 n-Butylbenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 74-83-9 n-Butylbenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 74-83-9 carbon tetrachloride ND ug/L 0.50 0.25 1 05/24/14 22:09 76-82-35 Chlorobenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 76-0-3 <	2-Fluorophenol (S)	34 %	1	10-120		1	05/23/14 11:30	06/02/14 22:36	367-12-4	
Benzene ND ug/L 0.50 0.25 1 05/24/14 22:09 71-43-2 Bromobenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 71-43-2 Bromochloromethane ND ug/L 0.50 0.25 1 05/24/14 22:09 74-97-5 Bromodichloromethane ND ug/L 0.50 0.25 1 05/24/14 22:09 75-27-4 Bromodorm ND ug/L 0.50 0.25 1 05/24/14 22:09 75-25-2 Bromomethane ND ug/L 0.50 0.25 1 05/24/14 22:09 74-83-9 n-Butylbenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 104-51-8 sec-Butylbenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 98-06-6 Carbon tetrachloride ND ug/L 0.50 0.25 1 05/24/14 22:09 96-63 Chlorobenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 76-63-3 Chloroform ND ug/L 0.50	2,4,6-Tribromophenol (S)	73 %	Î	10-137		1	05/23/14 11:30	06/02/14 22:36	118-79-6	
Bromobenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 108-86-1 Bromochloromethane ND ug/L 0.50 0.25 1 05/24/14 22:09 74-97-5 Bromodichloromethane ND ug/L 0.50 0.25 1 05/24/14 22:09 75-27-4 Bromoform ND ug/L 0.50 0.25 1 05/24/14 22:09 75-25-2 Bromoethane ND ug/L 0.50 0.25 1 05/24/14 22:09 74-83-9 n-Butylbenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 135-98-8 tert-Butylbenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 98-06-6 Carbon tetrachloride ND ug/L 0.50 0.25 1 05/24/14 22:09 96-03-5 Chlorobenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 96-63-3 Chlorotomethane ND ug/L 0.50 0.25 1 05/24/14 22:09 96-63-3 Chlorotomethane ND ug/L	6200B MSV	Analytical Me	thod: SM 6200	B						
Bromochloromethane ND ug/L 0.50 0.25 1 05/24/14 22:09 74-97-5 Bromodichloromethane ND ug/L 0.50 0.25 1 05/24/14 22:09 75-27-4 Bromodichloromethane ND ug/L 0.50 0.25 1 05/24/14 22:09 75-25-2 Bromomethane ND ug/L 0.50 0.25 1 05/24/14 22:09 74-83-9 n-Butylbenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 135-88-8 tert-Butylbenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 98-06-6 Carbon tetrachloride ND ug/L 0.50 0.25 1 05/24/14 22:09 98-08-6 Chlorobenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 98-08-6 Chlorobenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 96-03-5 Chlorobenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 95-03-3 Chlorotofure ND ug/L	Benzene	ND ug/L		0.50	0.25	1		05/24/14 22:09	71-43-2	
Bromodichloromethane ND ug/L 0.50 0.25 1 05/24/14 22:09 75-27-4 Bromoform ND ug/L 0.50 0.25 1 05/24/14 22:09 75-25-2 Bromomethane ND ug/L 0.50 0.50 1 05/24/14 22:09 74-83-9 n-Butylbenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 104-51-8 sec-Butylbenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 98-06-6 Carbon tetrachloride ND ug/L 0.50 0.25 1 05/24/14 22:09 98-06-6 Chlorobenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 98-06-6 Chlorobenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 96-07-3 Chlorobenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 75-0-3 Chloroform ND ug/L 0.50 0.25 1 05/24/14 22:09 96-49-8 2-Chlorotoluene ND ug/L 0.50<	Bromobenzene	ND ug/L		0.50	0.25	1		05/24/14 22:09	108-86-1	
BromoformND ug/L0.500.25105/24/14 22:0975-25-2BromomethaneND ug/L5.00.50105/24/14 22:0974-83-9n-ButylbenzeneND ug/L0.500.25105/24/14 22:09104-51-8sec-ButylbenzeneND ug/L0.500.25105/24/14 22:09135-98-8tert-ButylbenzeneND ug/L0.500.25105/24/14 22:0998-06-6Carbon tetrachlorideND ug/L0.500.25105/24/14 22:0996-63ChlorobenzeneND ug/L0.500.25105/24/14 22:0975-00-3ChloroformND ug/L0.500.25105/24/14 22:0975-00-3ChloroformND ug/L0.500.25105/24/14 22:0974-87-32-ChlorotolueneND ug/L0.500.25105/24/14 22:0974-87-32-ChlorotolueneND ug/L0.500.25105/24/14 22:0974-87-32-ChlorotolueneND ug/L0.500.25105/24/14 22:0995-49-84-ChlorotolueneND ug/L0.500.25105/24/14 22:0996-12-8DibromochloromethaneND ug/L0.500.25105/24/14 22:0996-12-81,2-Dibromo-thane (EDB)ND ug/L0.500.25105/24/14 22:0974-85-31,2-Dibromoethane (EDB)ND ug/L0.500.25105/24/14 22:0974-95-31,2-DibromoethaneND ug/L <t< td=""><td>Bromochloromethane</td><td>ND ug/L</td><td></td><td>0.50</td><td>0.25</td><td>1</td><td></td><td>05/24/14 22:09</td><td>74-97-5</td><td></td></t<>	Bromochloromethane	ND ug/L		0.50	0.25	1		05/24/14 22:09	74-97-5	
BromomethaneND ug/L5.00.50105/24/14 22:0974-83-9n-ButylbenzeneND ug/L0.500.25105/24/14 22:09104-51-8sec-ButylbenzeneND ug/L0.500.25105/24/14 22:09135-98-8tert-ButylbenzeneND ug/L0.500.25105/24/14 22:0998-06-6Carbon tetrachlorideND ug/L0.500.25105/24/14 22:0956-23-5ChlorobenzeneND ug/L0.500.25105/24/14 22:0975-00-3ChlorothaneND ug/L1.00.50105/24/14 22:0975-00-3ChlorothaneND ug/L0.500.25105/24/14 22:0976-6-3ChlorothueneND ug/L1.00.50105/24/14 22:0974-87-32-ChlorothueneND ug/L0.500.25105/24/14 22:0974-87-32-ChlorothueneND ug/L0.500.25105/24/14 22:0974-87-32-ChlorothueneND ug/L0.500.25105/24/14 22:0974-87-32-ChlorothueneND ug/L0.500.25105/24/14 22:0974-87-31,2-Dibromo-3-chloropropaneND ug/L0.500.25105/24/14 22:0974-87-31,2-Dibromo-thane (EDB)ND ug/L0.500.25105/24/14 22:0974-95-31,2-DibromoethaneND ug/L0.500.25105/24/14 22:0974-95-31,2-DibromoethaneND ug/L<	Bromodichloromethane	ND ug/L		0.50	0.25	1		05/24/14 22:09	75-27-4	
n-ButylbenzeneND ug/L0.500.25105/24/14 22:09104-51-8sec-ButylbenzeneND ug/L0.500.25105/24/14 22:09135-98-8tert-ButylbenzeneND ug/L0.500.25105/24/14 22:0998-06-6Carbon tetrachlorideND ug/L0.500.25105/24/14 22:0956-23-5ChlorobenzeneND ug/L0.500.25105/24/14 22:0975-00-3ChloroformND ug/L0.500.25105/24/14 22:0975-00-3ChloroformND ug/L0.500.25105/24/14 22:0976-66-3ChlorotolueneND ug/L0.500.25105/24/14 22:0995-49-84-ChlorotolueneND ug/L0.500.25105/24/14 22:0995-49-81,2-Dibromo-3-chloropropaneND ug/L0.500.25105/24/14 22:0996-12-8DibromochloromethaneKD ug/L0.500.25105/24/14 22:0996-12-8DibromochloromethaneKD ug/L0.500.25105/24/14 22:0996-12-8DibromochloromethaneKD ug/L0.500.25105/24/14 22:09106-43-4DibromochloromethaneKD ug/L0.500.25105/24/14 22:09106-93-4DibromochloromethaneKD ug/L0.500.25105/24/14 22:0974-95-31,2-DichlorobenzeneND ug/L0.500.25105/24/14 22:0995-50-11,3-Dich	Bromoform	ND ug/L		0.50	0.25	1		05/24/14 22:09	75-25-2	
sec-ButylbenzeneND ug/L0.500.25105/24/14 22:09135-98-8tert-ButylbenzeneND ug/L0.500.25105/24/14 22:0998-06-6Carbon tetrachlorideND ug/L0.500.25105/24/14 22:0956-23-5ChlorobenzeneND ug/L0.500.25105/24/14 22:0975-00-3ChlorothaneND ug/L1.00.50105/24/14 22:0975-00-3ChlorothaneND ug/L0.500.25105/24/14 22:0975-00-3ChlorothaneND ug/L0.500.25105/24/14 22:0974-87-32-ChlorotolueneND ug/L0.500.25105/24/14 22:0995-49-84-ChlorotolueneND ug/L0.500.25105/24/14 22:0996-12-8Dibromo-3-chloropropaneND ug/L0.500.25105/24/14 22:0996-12-8Dibromoethane (EDB)ND ug/L0.500.25105/24/14 22:09124-48-11,2-DibromoethaneND ug/L0.500.25105/24/14 22:0914-95-31,2-DichlorobenzeneND ug/L0.500.25105/24/14 22:0974-95-31,2-DichlorobenzeneND ug/L0.500.25105/24/14 22:0995-50-11,3-DichlorobenzeneND ug/L0.500.25105/24/14 22:0995-50-11,3-DichlorobenzeneND ug/L0.500.25105/24/14 22:0995-50-1	Bromomethane	ND ug/L		5.0	0.50	1		05/24/14 22:09	74-83-9	
tert-ButylbenzeneND ug/L0.500.25105/24/14 22:0998-06-6Carbon tetrachlorideND ug/L0.500.25105/24/14 22:0956-23-5ChlorobenzeneND ug/L0.500.25105/24/14 22:09108-90-7ChlorothaneND ug/L1.00.50105/24/14 22:0975-00-3ChlorothaneND ug/L0.500.25105/24/14 22:0967-66-3ChlorothaneND ug/L1.00.50105/24/14 22:0974-87-32-ChlorotolueneND ug/L0.500.25105/24/14 22:0995-49-84-ChlorotolueneND ug/L0.500.25105/24/14 22:0996-12-81,2-Dibromo-3-chloropropaneND ug/L1.00.50105/24/14 22:09124-48-11,2-Dibromoethane (EDB)ND ug/L0.500.25105/24/14 22:09106-93-41,2-DibromoethaneND ug/L0.500.25105/24/14 22:0974-95-31,2-DichlorobenzeneND ug/L0.500.25105/24/14 22:0974-95-31,2-DichlorobenzeneND ug/L0.500.25105/24/14 22:0995-50-11,3-DichlorobenzeneND ug/L0.500.25105/24/14 22:0995-50-11,3-DichlorobenzeneND ug/L0.500.25105/24/14 22:0995-50-1	n-Butylbenzene	ND ug/L		0.50	0.25	1		05/24/14 22:09	104-51-8	
Carbon tetrachlorideND ug/L0.500.25105/24/14 22:0956-23-5ChlorobenzeneND ug/L0.500.25105/24/14 22:09108-90-7ChloroethaneND ug/L1.00.50105/24/14 22:0975-00-3ChloroformND ug/L0.500.25105/24/14 22:0967-66-3ChloroethaneND ug/L1.00.50105/24/14 22:0974-87-32-ChlorotolueneND ug/L0.500.25105/24/14 22:0995-49-84-ChlorotolueneND ug/L0.500.25105/24/14 22:0996-12-81,2-Dibromo-3-chloropropaneND ug/L0.500.25105/24/14 22:09124-48-11,2-Dibromoethane (EDB)ND ug/L0.500.25105/24/14 22:09124-48-11,2-DichlorobenzeneND ug/L0.500.25105/24/14 22:09124-48-11,2-DichlorobenzeneND ug/L0.500.25105/24/14 22:09124-48-11,3-DichlorobenzeneND ug/L0.500.25105/24/14 22:0974-95-31,3-DichlorobenzeneND ug/L0.500.25105/24/14 22:0974-95-31,3-DichlorobenzeneND ug/L0.500.25105/24/14 22:0974-95-31,3-DichlorobenzeneND ug/L0.500.25105/24/14 22:0995-50-11,3-DichlorobenzeneND ug/L0.500.25105/24/14 22:09541-73-1 </td <td>sec-Butylbenzene</td> <td>ND ug/L</td> <td></td> <td>0.50</td> <td>0.25</td> <td>1</td> <td></td> <td>05/24/14 22:09</td> <td>135-98-8</td> <td></td>	sec-Butylbenzene	ND ug/L		0.50	0.25	1		05/24/14 22:09	135-98-8	
ChlorobenzeneND ug/L0.500.25105/24/14 22:09108-90-7ChloroethaneND ug/L1.00.50105/24/14 22:0975-00-3ChloroformND ug/L0.500.25105/24/14 22:0967-66-3ChloromethaneND ug/L1.00.50105/24/14 22:0974-87-32-ChlorotolueneND ug/L0.500.25105/24/14 22:0995-49-84-ChlorotolueneND ug/L0.500.25105/24/14 22:0996-12-81,2-Dibromo-3-chloropropaneND ug/L0.500.25105/24/14 22:0996-12-8Dibromoethane (EDB)ND ug/L0.500.25105/24/14 22:09124-48-11,2-DibromoethaneND ug/L0.500.25105/24/14 22:09124-48-11,2-DibromoethaneND ug/L0.500.25105/24/14 22:09124-48-11,2-DibromoethaneND ug/L0.500.25105/24/14 22:09124-48-11,2-DibromoethaneND ug/L0.500.25105/24/14 22:0974-95-31,2-DichlorobenzeneND ug/L0.500.25105/24/14 22:0995-50-11,3-DichlorobenzeneND ug/L0.500.25105/24/14 22:09541-73-1	tert-Butylbenzene	ND ug/L		0.50	0.25	1		05/24/14 22:09	98-06-6	
ChloroethaneND ug/L1.00.50105/24/14 22:0975-00-3ChloroformND ug/L0.500.25105/24/14 22:0967-66-3ChloromethaneND ug/L1.00.50105/24/14 22:0974-87-32-ChlorotolueneND ug/L0.500.25105/24/14 22:0995-49-84-ChlorotolueneND ug/L0.500.25105/24/14 22:0996-43-41,2-Dibromo-3-chloropropaneND ug/L1.00.50105/24/14 22:0996-12-8DibromochloromethaneND ug/L0.500.25105/24/14 22:09124-48-11,2-Dibromoethane (EDB)ND ug/L0.500.25105/24/14 22:09106-93-4DibromoethaneND ug/L0.500.25105/24/14 22:0974-95-31,2-DichlorobenzeneND ug/L0.500.25105/24/14 22:0995-50-11,3-DichlorobenzeneND ug/L0.500.25105/24/14 22:0995-50-11,3-DichlorobenzeneND ug/L0.500.25105/24/14 22:0995-50-1	Carbon tetrachloride	ND ug/L		0.50	0.25	1		05/24/14 22:09	56-23-5	
ChloroformND ug/L0.500.25105/24/14 22:0967-66-3ChloromethaneND ug/L1.00.50105/24/14 22:0974-87-32-ChlorotolueneND ug/L0.500.25105/24/14 22:0995-49-84-ChlorotolueneND ug/L0.500.25105/24/14 22:09106-43-41,2-Dibromo-3-chloropropaneND ug/L1.00.50105/24/14 22:0996-12-8DibromochloromethaneND ug/L0.500.25105/24/14 22:09124-48-11,2-Dibromoethane (EDB)ND ug/L0.500.25105/24/14 22:09106-93-4DibromomethaneND ug/L0.500.25105/24/14 22:0974-95-31,2-DichlorobenzeneND ug/L0.500.25105/24/14 22:0995-50-11,3-DichlorobenzeneND ug/L0.500.25105/24/14 22:0995-50-11,3-DichlorobenzeneND ug/L0.500.25105/24/14 22:0995-1-1	Chlorobenzene	ND ug/L		0.50	0.25	1		05/24/14 22:09	108-90-7	
ChloromethaneND ug/L1.00.50105/24/14 22:0974-87-32-ChlorotolueneND ug/L0.500.25105/24/14 22:0995-49-84-ChlorotolueneND ug/L0.500.25105/24/14 22:09106-43-41,2-Dibromo-3-chloropropaneND ug/L1.00.50105/24/14 22:0996-12-8DibromochloromethaneND ug/L0.500.25105/24/14 22:09124-48-11,2-Dibromoethane (EDB)ND ug/L0.500.25105/24/14 22:09106-93-4DibromomethaneND ug/L0.500.25105/24/14 22:0974-95-31,2-DichlorobenzeneND ug/L0.500.25105/24/14 22:0995-50-11,3-DichlorobenzeneND ug/L0.500.25105/24/14 22:0995-50-11,3-DichlorobenzeneND ug/L0.500.25105/24/14 22:09541-73-1	Chloroethane	ND ug/L		1.0	0.50	1		05/24/14 22:09	75-00-3	
2-ChlorotolueneND ug/L0.500.25105/24/14 22:0995-49-84-ChlorotolueneND ug/L0.500.25105/24/14 22:09106-43-41,2-Dibromo-3-chloropropaneND ug/L1.00.50105/24/14 22:0996-12-8DibromochloromethaneND ug/L0.500.25105/24/14 22:09124-48-11,2-Dibromoethane (EDB)ND ug/L0.500.25105/24/14 22:09106-93-4DibromoethaneND ug/L0.500.25105/24/14 22:0974-95-31,2-DichlorobenzeneND ug/L0.500.25105/24/14 22:0995-50-11,3-DichlorobenzeneND ug/L0.500.25105/24/14 22:09541-73-1	Chloroform	ND ug/L		0.50	0.25	1		05/24/14 22:09	67-66-3	
4-ChlorotolueneND ug/L0.500.25105/24/14 22:09106-43-41,2-Dibromo-3-chloropropaneND ug/L1.00.50105/24/14 22:0996-12-8DibromochloromethaneND ug/L0.500.25105/24/14 22:09124-48-11,2-Dibromoethane (EDB)ND ug/L0.500.25105/24/14 22:09106-93-4DibromomethaneND ug/L0.500.25105/24/14 22:0974-95-31,2-DichlorobenzeneND ug/L0.500.25105/24/14 22:0995-50-11,3-DichlorobenzeneND ug/L0.500.25105/24/14 22:09541-73-1	Chloromethane	ND ug/L		1.0	0.50	1		05/24/14 22:09	74-87-3	
1,2-Dibromo-3-chloropropaneND ug/L1.00.50105/24/14 22:0996-12-8DibromochloromethaneND ug/L0.500.25105/24/14 22:09124-48-11,2-Dibromoethane (EDB)ND ug/L0.500.25105/24/14 22:09106-93-4DibromomethaneND ug/L0.500.25105/24/14 22:0974-95-31,2-DichlorobenzeneND ug/L0.500.25105/24/14 22:0995-50-11,3-DichlorobenzeneND ug/L0.500.25105/24/14 22:09541-73-1	2-Chlorotoluene	ND ug/L		0.50	0.25	1		05/24/14 22:09	95-49-8	
DibromochloromethaneND ug/L0.500.25105/24/14 22:09124-48-11,2-Dibromoethane (EDB)ND ug/L0.500.25105/24/14 22:09106-93-4DibromomethaneND ug/L0.500.25105/24/14 22:0974-95-31,2-DichlorobenzeneND ug/L0.500.25105/24/14 22:0995-50-11,3-DichlorobenzeneND ug/L0.500.25105/24/14 22:09541-73-1	4-Chlorotoluene	ND ug/L		0.50	0.25	1		05/24/14 22:09	106-43-4	
1,2-Dibromoethane (EDB)ND ug/L0.500.25105/24/14 22:09106-93-4DibromomethaneND ug/L0.500.25105/24/14 22:0974-95-31,2-DichlorobenzeneND ug/L0.500.25105/24/14 22:0995-50-11,3-DichlorobenzeneND ug/L0.500.25105/24/14 22:09541-73-1	1,2-Dibromo-3-chloropropane	ND ug/L		1.0	0.50	1		05/24/14 22:09	96-12-8	
Dibromomethane ND ug/L 0.50 0.25 1 05/24/14 22:09 74-95-3 1,2-Dichlorobenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 95-50-1 1,3-Dichlorobenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 541-73-1	Dibromochloromethane	ND ug/L		0.50	0.25	1		05/24/14 22:09	124-48-1	
1,2-DichlorobenzeneND ug/L0.500.25105/24/14 22:0995-50-11,3-DichlorobenzeneND ug/L0.500.25105/24/14 22:09541-73-1	1,2-Dibromoethane (EDB)	ND ug/L		0.50	0.25	1		05/24/14 22:09	106-93-4	
1,3-Dichlorobenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 541-73-1	Dibromomethane	ND ug/L		0.50	0.25	1		05/24/14 22:09	74-95-3	
	1,2-Dichlorobenzene	ND ug/L		0.50	0.25	1		05/24/14 22:09	95-50-1	
1,4-Dichlorobenzene ND ug/L 0.50 0.25 1 05/24/14 22:09 106-46-7	1,3-Dichlorobenzene	ND ug/L		0.50	0.25	1		05/24/14 22:09	541-73-1	
	1,4-Dichlorobenzene	ND ug/L		0.50	0.25	1		05/24/14 22:09	106-46-7	
Dichlorodifluoromethane ND ug/L 0.50 0.25 1 05/24/14 22:09 75-71-8	Dichlorodifluoromethane	ND ug/L		0.50	0.25	1		05/24/14 22:09	75-71-8	
1,1-Dichloroethane ND ug/L 0.50 0.25 1 05/24/14 22:09 75-34-3	1,1-Dichloroethane	ND ug/L		0.50	0.25	1		05/24/14 22:09	75-34-3	
1,2-Dichloroethane ND ug/L 0.50 0.25 1 05/24/14 22:09 107-06-2	1,2-Dichloroethane	ND ug/L		0.50	0.25	1		05/24/14 22:09	107-06-2	
1,1-Dichloroethene ND ug/L 0.50 0.25 1 05/24/14 22:09 75-35-4	1,1-Dichloroethene	ND ug/L		0.50	0.25	1		05/24/14 22:09	75-35-4	
cis-1,2-Dichloroethene ND ug/L 0.50 0.25 1 05/24/14 22:09 156-59-2	cis-1,2-Dichloroethene	ND ug/L		0.50	0.25	1		05/24/14 22:09	156-59-2	
trans-1,2-Dichloroethene ND ug/L 0.50 0.25 1 05/24/14 22:09 156-60-5	trans-1,2-Dichloroethene	ND ug/L		0.50	0.25	1		05/24/14 22:09	156-60-5	
1,2-Dichloropropane ND ug/L 0.50 0.25 1 05/24/14 22:09 78-87-5	1,2-Dichloropropane	ND ug/L		0.50	0.25	1		05/24/14 22:09	78-87-5	
1,3-Dichloropropane ND ug/L 0.50 0.25 1 05/24/14 22:09 142-28-9	1,3-Dichloropropane	ND ug/L		0.50	0.25	1		05/24/14 22:09	142-28-9	



ANALYTICAL RESULTS

Project: NCDOT FRONT ST. WBS17BP.3.R.28

Pace Project No.: 9

92202425

Sample: 3-08	Lab ID: 92202425001	Collecte	d: 05/16/14	12:25	Received: 05/22/14 09:30 Matrix: Water	
		Report				
Parameters	Results Units	Limit	MDL	DF	Prepared Analyzed CAS No.	Qual
6200B MSV	Analytical Method: SM 62	200B				
2,2-Dichloropropane	ND ug/L	0.50	0.25	1	05/24/14 22:09 594-20-7	
1,1-Dichloropropene	ND ug/L	0.50	0.25	1	05/24/14 22:09 563-58-6	
cis-1,3-Dichloropropene	ND ug/L	0.50	0.25	1	05/24/14 22:09 10061-01-5	
trans-1,3-Dichloropropene	ND ug/L	0.50	0.25	1	05/24/14 22:09 10061-02-6	
Diisopropyl ether	ND ug/L	0.50	0.25	1	05/24/14 22:09 108-20-3	
Ethylbenzene	ND ug/L	0.50	0.25	1	05/24/14 22:09 100-41-4	
Hexachloro-1,3-butadiene	ND ug/L	2.0	1.0	1	05/24/14 22:09 87-68-3	
Isopropylbenzene (Cumene)	ND ug/L	0.50	0.25	1	05/24/14 22:09 98-82-8	
Methylene Chloride	ND ug/L	2.0	1.0	1	05/24/14 22:09 75-09-2	
Methyl-tert-butyl ether	ND ug/L	0.50	0.25	1	05/24/14 22:09 1634-04-4	
Naphthalene	ND ug/L	2.0	1.0	1	05/24/14 22:09 91-20-3	
n-Propylbenzene	ND ug/L	0.50	0.25	1	05/24/14 22:09 103-65-1	
Styrene	ND ug/L	0.50	0.25	1	05/24/14 22:09 100-42-5	
1,1,1,2-Tetrachloroethane	ND ug/L	0.50	0.25	1	05/24/14 22:09 630-20-6	
1,1,2,2-Tetrachloroethane	ND ug/L	0.50	0.25	1	05/24/14 22:09 79-34-5	
Tetrachloroethene	ND ug/L	0.50	0.25	1	05/24/14 22:09 127-18-4	
Toluene	ND ug/L	0.50	0.25	1	05/24/14 22:09 108-88-3	
1,2,3-Trichlorobenzene	ND ug/L	2.0	1.0	1	05/24/14 22:09 87-61-6	
1,2,4-Trichlorobenzene	ND ug/L	2.0	1.0	1	05/24/14 22:09 120-82-1	
1,1,1-Trichloroethane	ND ug/L	0.50	0.25	1	05/24/14 22:09 71-55-6	
1,1,2-Trichloroethane	ND ug/L	0.50	0.25	1	05/24/14 22:09 79-00-5	
Trichloroethene	ND ug/L	0.50	0.25	1	05/24/14 22:09 79-01-6	
Trichlorofluoromethane	ND ug/L	1.0	0.50	1	05/24/14 22:09 75-69-4	
1,2,3-Trichloropropane	ND ug/L	0.50	0.25	1	05/24/14 22:09 96-18-4	
1,2,4-Trimethylbenzene	ND ug/L	0.50	0.25	1	05/24/14 22:09 95-63-6	
1,3,5-Trimethylbenzene	ND ug/L	0.50	0.25	1	05/24/14 22:09 108-67-8	
Vinyl chloride	ND ug/L	1.0	0.25	1	05/24/14 22:09 75-01-4	
m&p-Xylene	ND ug/L	1.0	0.50	1	05/24/14 22:09 179601-23-1	
o-Xylene	ND ug/L	0.50	0.25	1	05/24/14 22:09 95-47-6	
Surrogates	···· ••• ••• •••	5.00	0.20	-		
1,2-Dichloroethane-d4 (S)	98 %	70-130		1	05/24/14 22:09 17060-07-0	
4-Bromofluorobenzene (S)	95 %	70-130		1	05/24/14 22:09 460-00-4	
Toluene-d8 (S)	101 %	70-130		1	05/24/14 22:09 2037-26-5	



Project: NCDOT FRONT ST. WBS17BP.3.R.28

Project: NCDOT FF	RONT ST. WBS17BP.3.R.	28			
Pace Project No.: 92202425					
QC Batch: MSV/269	40	Analysis Meth	od: SN	/ 6200B	
QC Batch Method: SM 6200	-	Analysis Desc		00B MSV	
		Analysis Desc	1.001. 02		
Associated Lab Samples: 92	202425001				
METHOD BLANK: 1206081		Matrix: \	Nater		
Associated Lab Samples: 92	202425001				
		Blank	Reporting		
Parameter	Units	Result	Limit	Analyzed	Qualifiers
1,1,1,2-Tetrachloroethane	ug/L	ND	0.50	05/24/14 18:18	
1,1,1-Trichloroethane	ug/L	ND	0.50	05/24/14 18:18	
1,1,2,2-Tetrachloroethane	ug/L	ND	0.50	05/24/14 18:18	
1,1,2-Trichloroethane	ug/L	ND	0.50	05/24/14 18:18	
1,1-Dichloroethane	ug/L	ND	0.50	05/24/14 18:18	
1,1-Dichloroethene	ug/L	ND	0.50	05/24/14 18:18	
1,1-Dichloropropene	ug/L	ND	0.50	05/24/14 18:18	
1,2,3-Trichlorobenzene	ug/L	ND	2.0	05/24/14 18:18	
1,2,3-Trichloropropane	ug/L	ND	0.50	05/24/14 18:18	
1,2,4-Trichlorobenzene	ug/L	ND	2.0	05/24/14 18:18	
1,2,4-Trimethylbenzene	ug/L	ND	0.50	05/24/14 18:18	
1,2-Dibromo-3-chloropropane	ug/L	ND	1.0	05/24/14 18:18	
1,2-Dibromoethane (EDB)	ug/L	ND	0.50	05/24/14 18:18	
1,2-Dichlorobenzene	ug/L	ND	0.50	05/24/14 18:18	
1,2-Dichloroethane	ug/L	ND	0.50	05/24/14 18:18	
1,2-Dichloropropane	ug/L	ND	0.50	05/24/14 18:18	
1,3,5-Trimethylbenzene	ug/L	ND	0.50	05/24/14 18:18	
1,3-Dichlorobenzene	ug/L	ND	0.50	05/24/14 18:18	
1,3-Dichloropropane	ug/L	ND	0.50	05/24/14 18:18	
1,4-Dichlorobenzene	ug/L	ND	0.50	05/24/14 18:18	
2,2-Dichloropropane	ug/L	ND	0.50	05/24/14 18:18	
2-Chlorotoluene	ug/L	ND	0.50	05/24/14 18:18	
4-Chlorotoluene	ug/L	ND	0.50	05/24/14 18:18	
Benzene	ug/L	ND	0.50	05/24/14 18:18	
Bromobenzene	ug/L	ND	0.50	05/24/14 18:18	
Bromochloromethane	ug/L	ND	0.50	05/24/14 18:18	
Bromodichloromethane	ug/L	ND	0.50	05/24/14 18:18	
Bromoform	ug/L	ND	0.50	05/24/14 18:18	
Bromomethane	ug/L	ND	5.0	05/24/14 18:18	

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.

ND

0.50 05/24/14 18:18

0.50 05/24/14 18:18

1.0 05/24/14 18:18

0.50 05/24/14 18:18

1.0 05/24/14 18:18

0.50 05/24/14 18:18

0.50 05/24/14 18:18

0.50 05/24/14 18:18

0.50 05/24/14 18:18

0.50 05/24/14 18:18

0.50 05/24/14 18:18

0.50 05/24/14 18:18

REPORT OF LABORATORY ANALYSIS

Carbon tetrachloride

cis-1,2-Dichloroethene

cis-1,3-Dichloropropene

Dibromochloromethane

Dichlorodifluoromethane

Chlorobenzene

Chloromethane

Dibromomethane

Diisopropyl ether

Ethylbenzene

Chloroethane

Chloroform

ug/L



Matrix: Water

Project: NCDOT FRONT ST. WBS17BP.3.R.28

Pace Project No.: 92202425

METHOD BLANK: 1206081

Associated Lab Samples: 92202425001

Parameter	Units	Blank Result	Reporting Limit	Analyzed	Qualifiers
Hexachloro-1,3-butadiene	ug/L	ND	2.0	05/24/14 18:18	
Isopropylbenzene (Cumene)	ug/L	ND	0.50	05/24/14 18:18	
m&p-Xylene	ug/L	ND	1.0	05/24/14 18:18	
Methyl-tert-butyl ether	ug/L	ND	0.50	05/24/14 18:18	
Methylene Chloride	ug/L	ND	2.0	05/24/14 18:18	
n-Butylbenzene	ug/L	ND	0.50	05/24/14 18:18	
n-Propylbenzene	ug/L	ND	0.50	05/24/14 18:18	
Naphthalene	ug/L	ND	2.0	05/24/14 18:18	
o-Xylene	ug/L	ND	0.50	05/24/14 18:18	
sec-Butylbenzene	ug/L	ND	0.50	05/24/14 18:18	
Styrene	ug/L	ND	0.50	05/24/14 18:18	
ert-Butylbenzene	ug/L	ND	0.50	05/24/14 18:18	
Tetrachloroethene	ug/L	ND	0.50	05/24/14 18:18	
Toluene	ug/L	ND	0.50	05/24/14 18:18	
rans-1,2-Dichloroethene	ug/L	ND	0.50	05/24/14 18:18	
rans-1,3-Dichloropropene	ug/L	ND	0.50	05/24/14 18:18	
Frichloroethene	ug/L	ND	0.50	05/24/14 18:18	
Trichlorofluoromethane	ug/L	ND	1.0	05/24/14 18:18	
/inyl chloride	ug/L	ND	1.0	05/24/14 18:18	
I,2-Dichloroethane-d4 (S)	%	97	70-130	05/24/14 18:18	
I-Bromofluorobenzene (S)	%	96	70-130	05/24/14 18:18	
Toluene-d8 (S)	%	100	70-130	05/24/14 18:18	

LABORATORY CONTROL SAMPLE: 1206082

		Spike	LCS	LCS	% Rec	
Parameter	Units	Conc.	Result	% Rec	Limits	Qualifiers
1,1,1,2-Tetrachloroethane	ug/L	50	57.3	115	60-140	
1,1,1-Trichloroethane	ug/L	50	48.8	98	60-140	
1,1,2,2-Tetrachloroethane	ug/L	50	54.0	108	60-140	
1,1,2-Trichloroethane	ug/L	50	52.7	105	60-140	
1,1-Dichloroethane	ug/L	50	47.5	95	60-140	
1,1-Dichloroethene	ug/L	50	50.9	102	60-140	
1,1-Dichloropropene	ug/L	50	51.0	102	60-140	
1,2,3-Trichlorobenzene	ug/L	50	57.5	115	60-140	
1,2,3-Trichloropropane	ug/L	50	52.7	105	60-140	
1,2,4-Trichlorobenzene	ug/L	50	55.2	110	60-140	
1,2,4-Trimethylbenzene	ug/L	50	57.9	116	60-140	
1,2-Dibromo-3-chloropropane	ug/L	50	55.7	111	60-140	
1,2-Dibromoethane (EDB)	ug/L	50	54.6	109	60-140	
1,2-Dichlorobenzene	ug/L	50	54.9	110	60-140	
1,2-Dichloroethane	ug/L	50	48.2	96	60-140	
1,2-Dichloropropane	ug/L	50	49.9	100	60-140	
1,3,5-Trimethylbenzene	ug/L	50	56.4	113	60-140	
1,3-Dichlorobenzene	ug/L	50	54.7	109	60-140	

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: NCDOT FRONT ST. WBS17BP.3.R.28

Pace Project No.: 92202425

LABORATORY CONTROL SAMPLE: 1206082

Dickhorobenzene ug/L 50 52.7 105 60-140 Dichloropropane ug/L 50 54.6 109 60-140 Dichlorobulene ug/L 50 53.1 106 60-140 nzene ug/L 50 53.8 106 60-140 mochloromethane ug/L 50 51.2 102 60-140 mochloromethane ug/L 50 51.2 102 60-140 mochloromethane ug/L 50 48.1 96 60-140 mochloromethane ug/L 50 51.7 103 60-140 iorobenzene ug/L 50 52.5 105 60-140 lorobenzene ug/L 50 52.5 101 60-140 lorobenzene ug/L 50 54.7 109 60-140 lorobenzene ug/L 50 51.1 102 60-140 lorobenzene ug/L 50 51.6 1014 60-1	Parameter	Units	Spike Conc.	LCS Result	LCS % Rec	% Rec Limits	Qualifiers
Dickhorobenzene ug/L 50 52.7 105 60-140 Dichloropropane ug/L 50 53.4 109 60-140 Dichlorobulene ug/L 50 53.1 106 60-140 Increne ug/L 50 53.9 108 60-140 Increne ug/L 50 51.2 102 60-140 Increne ug/L 50 51.2 102 60-140 Increne ug/L 50 56.6 113 60-140 Increnteritane ug/L 50 54.7 109 60-140 Increnteritane ug/L 50 52.5 103 60-140 Increnteritane ug/L 50 52.5 105 60-140 Increnteritane ug/L 50 56.6 113 60-140 Increnteritane ug/L 50 51.7 103 60-140 Increnteritane ug/L 50 51.7 109 60-140	1,3-Dichloropropane	ug/L		52.4	105	60-140	
Chlorotoluene ug/L 50 54.6 109 60-140 Chlorotoluene ug/L 50 53.1 106 60-140 Dronobnormethane ug/L 50 53.9 108 60-140 Dronobnormethane ug/L 50 51.2 102 60-140 Dronomethane ug/L 50 54.6 113 60-140 Dronomethane ug/L 50 48.1 96 60-140 Dronomethane ug/L 50 54.7 109 60-140 Iorobenzene ug/L 50 54.7 109 60-140 Iorobenzene ug/L 50 52.5 105 60-140 Iorobenzene ug/L 50 52.6 101 60-140 Iorobenzene ug/L 50 54.7 109 60-140 Iorobenzene ug/L 50 54.7 109 60-140 Iorobenzene ug/L 50 51.7 103 60-140 <td>1,4-Dichlorobenzene</td> <td></td> <td>50</td> <td>52.7</td> <td>105</td> <td>60-140</td> <td></td>	1,4-Dichlorobenzene		50	52.7	105	60-140	
biorotoluene ug/L 50 53.1 106 60-140 nzene ug/L 50 52.8 106 60-140 pmobenzene ug/L 50 51.2 102 60-140 pmochloromethane ug/L 50 56.6 113 60-140 pmodinbromethane ug/L 50 48.6 99 60-140 pmodinbromethane ug/L 50 48.7 109 60-140 pmoderbane ug/L 50 54.7 103 60-140 lorobenzene ug/L 50 51.7 103 60-140 lorobenzene ug/L 50 52.5 105 60-140 loroform ug/L 50 51.1 102 60-140 loromethane ug/L 50 54.7 109 60-140 rycomachicromethane ug/L 50 54.7 109 60-140 rycomachicromethane ug/L 50 52.5 105 60-14	2,2-Dichloropropane	ug/L	50	47.3	95	60-140	
Shlorobluene ug/L 50 53.1 106 60-140 nzene ug/L 50 52.8 108 60-140 mochoromethane ug/L 50 51.2 102 60-140 mochichoromethane ug/L 50 56.6 113 60-140 morofic ug/L 50 48.1 96 60-140 morofic ug/L 50 48.1 96 60-140 morethane ug/L 50 54.7 109 60-140 lorobenzene ug/L 50 51.7 103 60-140 lorobenzene ug/L 50 52.5 105 60-140 lorobenzene ug/L 50 56.6 113 60-140 lorobenzene ug/L 50 54.7 109 60-140 lorobenzene ug/L 50 54.7 109 60-140 lorobenzene ug/L 50 52.5 105 60-140	2-Chlorotoluene	ug/L	50	54.6	109	60-140	
nzene ug/L 50 52.8 106 60-140 mochenzene ug/L 50 53.9 108 60-140 modeichloromethane ug/L 50 56.6 113 60-140 modoichloromethane ug/L 50 48.1 96 60-140 morenthane ug/L 50 48.1 96 60-140 rbon tetrachloride ug/L 50 54.7 109 60-140 lorobenzene ug/L 50 52.5 105 60-140 lorobratene ug/L 50 52.5 105 60-140 loroform ug/L 50 56.6 113 60-140 loroform ug/L 50 54.7 109 60-140 -1_2-Dichloroethene ug/L 50 54.7 109 60-140 vormomethane ug/L 50 52.5 105 60-140 vormomethane ug/L 50 57.6 115 60-140 <td>-Chlorotoluene</td> <td>-</td> <td>50</td> <td>53.1</td> <td>106</td> <td>60-140</td> <td></td>	-Chlorotoluene	-	50	53.1	106	60-140	
nmochoramethane ug/L 50 53.9 108 60-140 nmochhoromethane ug/L 50 51.2 102 60-140 nmochhoromethane ug/L 50 48.1 96 60-140 nmorethane ug/L 50 48.1 96 60-140 nmorethane ug/L 50 51.7 103 60-140 lorobenzene ug/L 50 52.5 105 60-140 lorobenzene ug/L 50 56.6 113 60-140 loromethane ug/L 50 56.6 113 60-140 loromethane ug/L 50 51.1 102 60-140 loromethane ug/L 50 54.7 109 60-140 romochloropropene ug/L 50 52.5 105 60-140 romochloromethane ug/L 50 52.5 105 60-140 romochloromethane ug/L 50 57.6 115 <	senzene	-	50	52.8	106	60-140	
amodiloromethane ug/L 50 51.2 102 60-140 amodichloromethane ug/L 50 56.6 113 60-140 amodichloromethane ug/L 50 49.6 99 60-140 bronotethane ug/L 50 54.7 109 60-140 lorobenzene ug/L 50 52.5 105 60-140 lorobenzene ug/L 50 52.5 105 60-140 lorobenzene ug/L 50 56.6 113 60-140 lorobenzene ug/L 50 51.1 102 60-140 loromethane ug/L 50 54.7 109 60-140 romodichloromethane ug/L 50 47.9 96 60-140 romodichloromethane ug/L 50 52.5 105 60-140 romodichloromethane ug/L 50 52.5 105 60-140 romodichloromethane ug/L 50 52.5	Bromobenzene	-	50	53.9	108	60-140	
pmodichloromethane ug/L 50 56.6 113 60-140 pmoform ug/L 50 48.1 96 60-140 pmomethane ug/L 50 54.7 109 60-140 lorobenzene ug/L 50 51.7 103 60-140 lorobenzene ug/L 50 52.5 105 60-140 lorobenzene ug/L 50 56.6 113 60-140 lorobentane ug/L 50 56.6 113 60-140 loromethane ug/L 50 54.7 109 60-140 -1,2-Dichloroethene ug/L 50 47.9 96 60-140 voromethane ug/L 50 47.9 96 60-140 voromethane ug/L 50 52.5 105 60-140 voromethane ug/L 50 57.6 115 60-140 voromethane ug/L 50 57.6 115 60-140 </td <td>Bromochloromethane</td> <td>-</td> <td>50</td> <td>51.2</td> <td>102</td> <td>60-140</td> <td></td>	Bromochloromethane	-	50	51.2	102	60-140	
nmonorm ug/L 50 48.1 96 60-140 mmonethane ug/L 50 49.6 99 60-140 bon tetrachloride ug/L 50 54.7 103 60-140 lorobenzene ug/L 50 51.7 103 60-140 lorobenzene ug/L 50 50.6 101 60-140 lorobenzene ug/L 50 50.6 113 60-140 lorobenzene ug/L 50 51.1 102 60-140 lorobenzene ug/L 50 54.7 109 60-140 -1,2-Dichloroptene ug/L 50 47.9 96 60-140 vormonethane ug/L 50 52.5 105 60-140 vormonethane ug/L 50 52.5 105 60-140 vormonethane ug/L 50 57.6 115 60-140 vormonethane ug/L 50 57.6 115 60-140	romodichloromethane		50	56.6	113	60-140	
pmomethane ug/L 50 49.6 99 60-140 hon tetrachloride ug/L 50 54.7 109 60-140 lorobenzene ug/L 50 51.7 103 60-140 lorobentane ug/L 50 52.5 105 60-140 lorobentane ug/L 50 56.6 113 60-140 lorobentane ug/L 50 51.1 102 60-140 1,2-Dichloroptene ug/L 50 54.7 109 60-140 romonethane ug/L 50 47.9 96 60-140 romonethane ug/L 50 47.9 96 60-140 romonethane ug/L 50 52.5 105 60-140 schoro-1,3-butadiene ug/L 50 57.2 114 60-140 up/L-pr/Sylene ug/L 50 51.6 103 60-140 up/L-pr/Sylence ug/L 50 51.6 103 6	romoform		50	48.1	96	60-140	
rbon tetrachloride ug/L 50 54.7 109 60-140 lorobenzene ug/L 50 51.7 103 60-140 lorobenzene ug/L 50 52.5 105 60-140 lorobentane ug/L 50 56.6 113 60-140 loromethane ug/L 50 51.1 102 60-140 romochloromethane ug/L 50 54.7 109 60-140 romochloromethane ug/L 50 54.7 109 60-140 romochloromethane ug/L 50 47.9 96 60-140 romochloromethane ug/L 50 62.3 125 60-140 romochloromethane ug/L 50 52.5 105 60-140 vjbenzene ug/L 50 57.2 114 60-140 vjbenzene ug/L 50 57.6 115 60-140 vjbenzene ug/L 50 57.6 115	romomethane		50	49.6	99	60-140	
borobenzene ug/L 50 51.7 103 60-140 loroethane ug/L 50 52.5 105 60-140 loroethane ug/L 50 50.6 101 60-140 loroethane ug/L 50 56.6 113 60-140 -1,2-Dichloroethene ug/L 50 54.7 109 60-140 -1,3-Dichloropropene ug/L 50 47.9 96 60-140 oromochloromethane ug/L 50 47.9 96 60-140 oromochloromethane ug/L 50 62.3 125 60-140 oromochloromethane ug/L 50 52.5 105 60-140 ylbenzene ug/L 50 57.2 114 60-140 yub/serzene ug/L 50 57.6 115 60-140 yub/serzene ug/L 50 57.6 115 60-140 yub/serzene ug/L 50 56.1 112	arbon tetrachloride	-					
boroethane ug/L 50 52.5 105 60-140 loroform ug/L 50 50.6 101 60-140 loromethane ug/L 50 56.6 113 60-140 -1,2-Dichloroethene ug/L 50 51.1 102 60-140 -1,3-Dichloropropene ug/L 50 47.9 96 60-140 oromomethane ug/L 50 47.9 96 60-140 oromomethane ug/L 50 49.9 100 60-140 oromomethane ug/L 50 52.5 105 60-140 oromomethane ug/L 50 52.5 105 60-140 holorofil/uromethane ug/L 50 57.6 115 60-140 hylbenzene ug/L 50 57.6 115 60-140 hylbenzene ug/L 50 51.6 103 60-140 hylbenzene ug/L 50 51.6 103 60-1	hlorobenzene					60-140	
boroform ug/L 50 50.6 101 60-140 loromethane ug/L 50 56.6 113 60-140 -1,2-Dichloroethene ug/L 50 51.1 102 60-140 -1,3-Dichloropropene ug/L 50 47.9 96 60-140 oromochloromethane ug/L 50 47.9 96 60-140 oromochloromethane ug/L 50 49.9 100 60-140 oromochloromethane ug/L 50 52.5 105 60-140 ylbenzene ug/L 50 57.2 114 60-140 ylbenzene ug/L 50 57.6 115 60-140 ypoylbenzene (Cumene) ug/L 50 57.6 115 60-140 thylere ther ug/L 50 57.6 115 60-140 thylere ther ug/L 50 51.6 103 60-140 thylere ther ug/L 50 51.6 103<	hloroethane	-					
boromethane ug/L 50 56.6 113 60-140 -1,2-Dichloroethene ug/L 50 51.1 102 60-140 -1,3-Dichloropropene ug/L 50 54.7 109 60-140 promochloromethane ug/L 50 49.9 100 60-140 promothane ug/L 50 49.9 100 60-140 promothane ug/L 50 49.9 100 60-140 schoro-1,3-butadiene ug/L 50 52.5 105 60-140 yblenzene ug/L 50 57.2 114 60-140 propyletner ug/L 50 57.6 115 60-140 propylenzene (Cumene) ug/L 50 57.6 115 60-140 thylene Aug/L 50 51.6 103 60-140 thylene ug/L 50 51.3 103 60-140 thylene Aug/L 50 57.1 114 60-140	hloroform	-					
-1,2-Dichloroethene ug/L 50 51.1 102 60-140 -1,3-Dichloropropene ug/L 50 54.7 109 60-140 oromochloromethane ug/L 50 47.9 96 60-140 oromochloromethane ug/L 50 49.9 100 60-140 oromomethane ug/L 50 62.3 125 60-140 sopropyl ether ug/L 50 52.5 105 60-140 viblorodifluoromethane ug/L 50 57.2 114 60-140 propylbenzene (Cumene) ug/L 50 57.6 115 60-140 propylbenzene (Cumene) ug/L 50 57.6 115 60-140 propylbenzene (Cumene) ug/L 50 51.6 103 60-140 thyl-tert-butyl ether ug/L 50 51.6 103 60-140 thyl-tert-butyl ether ug/L 50 51.3 103 60-140 viblenzene ug/L 50 56.1 112 60-140 viblenzene ug/L <td>hloromethane</td> <td>0</td> <td>50</td> <td>56.6</td> <td>113</td> <td>60-140</td> <td></td>	hloromethane	0	50	56.6	113	60-140	
-1,3-Dichloropropene ug/L 50 54.7 109 60-140 promochloromethane ug/L 50 47.9 96 60-140 promomethane ug/L 50 49.9 100 60-140 promomethane ug/L 50 62.3 125 60-140 sopropylether ug/L 50 52.5 105 60-140 nylbenzene ug/L 50 52.5 105 60-140 nylbenzene ug/L 50 57.2 114 60-140 propylbenzene (Cumene) ug/L 50 57.6 115 60-140 proylbenzene ug/L 100 108 108 60-140 thyl-tert-butyl ether ug/L 50 51.6 103 60-140 thyl-tert-butyl ether ug/L 50 51.3 103 60-140 thyl-tert-butyl ether ug/L 50 51.3 103 60-140 thyl-tert-butyl ether ug/L 50 57.1 114 60-140 thylenzene ug/L 50 <	s-1.2-Dichloroethene	0					
bromochloromethane ug/L 50 47.9 96 60-140 promochloromethane ug/L 50 49.9 100 60-140 promochloromethane ug/L 50 62.3 125 60-140 sopropyl ether ug/L 50 52.5 105 60-140 nylbenzene ug/L 50 57.2 114 60-140 propylbenzene (Cumene) ug/L 50 57.2 114 60-140 propylbenzene (Cumene) ug/L 50 57.6 115 60-140 thyl-tert-butyl ether ug/L 50 57.6 115 60-140 thyl-tert-butyl ether ug/L 50 51.6 103 60-140 propylBenzene ug/L 50 56.1 112 60-140 propylBenzene ug/L 50 56.1 112 60-140 propylBenzene ug/L 50 57.1 114 60-140 propylBenzene ug/L 50		-				60-140	
orromomethane ug/L 50 49.9 100 60-140 chlorodifluoromethane ug/L 50 62.3 125 60-140 sopropyl ether ug/L 50 52.5 105 60-140 ylbenzene ug/L 50 52.5 105 60-140 yachloro-1,3-butadiene ug/L 50 57.2 114 60-140 propylbenzene (Cumene) ug/L 50 57.6 115 60-140 yb-Xylene ug/L 50 51.6 103 60-140 thyl-tert-butyl ether ug/L 50 51.6 103 60-140 stylbenzene ug/L 50 51.6 103 60-140 stylbenzene ug/L 50 51.3 103 60-140 verylbenzene ug/L 50 51.3 103 60-140 verylbenzene ug/L 50 57.1 114 60-140 verne ug/L 50 57.0 114 <td></td> <td>-</td> <td></td> <td></td> <td></td> <td></td> <td></td>		-					
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		•					
P-Dichloroethane-d4 (S) % 94 70-130	2-Dichloroethane-d4 (S)	%	50	02.1			
	Bromofluorobenzene (S)						
	bluene-d8 (S)						

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: NCDOT FRONT ST. WBS17BP.3.R.28

Pace Project No.: 92202425

MATRIX SPIKE & MATRIX SPI	KE DUPLICATI	E: 12072			1207286							
			MS	MSD					04 F			
Parameter	922 Units	02441006 Result	Spike Conc.	Spike Conc.	MS Result	MSD Result	MS % Rec	MSD % Rec	% Rec Limits	RPD	Max RPD	Qual
1,1,1,2-Tetrachloroethane	ug/L	ND	20	20	23.8	20.6	119	103	60-140	15	30	
1,1,1-Trichloroethane	ug/L	ND	20	20	22.3	19.5	111	98	60-140	13	30	
1,1,2,2-Tetrachloroethane	ug/L	ND	20	20	22.2	19.5	111	97	60-140	13	30	
1,1,2-Trichloroethane	ug/L	ND	20	20	22.2	19.3	111	97	60-140	14	30	
1,1-Dichloroethane	ug/L	ND	20	20	20.8	18.5	104	93	60-140	11	30	
1,1-Dichloroethene	ug/L	ND	20	20	23.0	20.5	115	103	60-140	11	30	
1,1-Dichloropropene	ug/L	ND	20	20	23.1	20.5	115	103	60-140	12	30	
1,2,3-Trichlorobenzene	ug/L	ND	20	20	21.7	19.2	109	96	60-140	12	30	
1,2,3-Trichloropropane	ug/L	ND	20	20	22.1	19.6	111	98	60-140	12	30	
1,2,4-Trichlorobenzene	ug/L	ND	20	20	21.0	18.5	105	92	60-140	13	30	
1,2,4-Trimethylbenzene	ug/L	ND	20	20	23.2	20.7	116	103	60-140	12	30	
1,2-Dibromo-3-chloropropane	ug/L	ND	20	20	23.1	19.6	115	98	60-140	16	30	
1,2-Dibromoethane (EDB)	ug/L	ND	20	20	22.6	20.3	113	101	60-140	11	30	
1,2-Dichlorobenzene	ug/L	ND	20	20	22.2	19.4	111	97	60-140	14	30	
1,2-Dichloroethane	ug/L	ND	20	20	20.8	18.5	104	92	60-140	12	30	
1,2-Dichloropropane	ug/L	ND	20	20	21.3	19.2	107	96	60-140	11	30	
1,3,5-Trimethylbenzene	ug/L	ND	20	20	22.8	20.6	114	103	60-140	10	30	
1,3-Dichlorobenzene	ug/L	ND	20	20	22.3	19.4	111	97	60-140	14	30	
1,3-Dichloropropane	ug/L	ND	20	20	21.7	19.5	109	97	60-140	11	30	
1,4-Dichlorobenzene	ug/L	ND	20	20	21.7	18.7	105	93	60-140	13	30	
2,2-Dichloropropane	ug/L	ND	20	20	20.5	18.2	100	91	60-140	12	30	
2-Chlorotoluene	ug/L	ND	20	20	20.3	10.2	102	99	60-140	12	30	
4-Chlorotoluene	ug/L	ND	20	20 20	22.3	19.8	107	99 96	60-140 60-140	11	30	
Benzene	ug/L	ND	20	20	21.3	20.2	107	101	60-140 60-140	12	30	
Bromobenzene	ug/L ug/L	ND	20	20 20	22.0	20.2 19.5	114	97	60-140 60-140	13	30	
Bromochloromethane		ND	20	20 20	22.2	20.5	116	103	60-140 60-140	12	30	
	ug/L					20.5	115				30	
Bromodichloromethane	ug/L	ND ND	20	20	22.9			103	60-140	10		
Bromoform	ug/L		20	20	20.4	18.2	102 117	91 96	60-140	11	30 30	
Bromomethane	ug/L	ND ND	20	20	23.3	19.3			60-140	19	30 30	
Carbon tetrachloride	ug/L	ND	20	20 20	24.0	21.3	120 108	106	60-140	12	30 30	
Chlorobenzene	ug/L		20		21.5	19.2		96	60-140	11		
Chloroethane	ug/L	ND	20	20	23.8	22.8	119	114	60-140	4	30	
Chloroform	ug/L	ND	20	20	22.2	19.7	111	99	60-140	12	30	
Chloromethane	ug/L	ND	20	20	24.0	21.9	120	109	60-140	9	30	
cis-1,2-Dichloroethene	ug/L	ND	20	20	22.6	19.6	113	98	60-140	15	30	
cis-1,3-Dichloropropene	ug/L	ND	20	20	22.0	19.5	110	98	60-140	12	30	
Dibromochloromethane	ug/L	ND	20	20	20.5	17.7	103	89	60-140	14	30	
Dibromomethane	ug/L	ND	20	20	21.0	18.7	105	94	60-140		30	
Dichlorodifluoromethane	ug/L	ND	20	20	26.9	24.0	135	120	60-140		30	
Diisopropyl ether	ug/L	ND	20	20	22.6	20.4	113	102	60-140		30	
Ethylbenzene	ug/L	ND	20	20	21.9	19.7	110	99	60-140		30	
Hexachloro-1,3-butadiene	ug/L	ND	20	20	22.6	20.4	113	102	60-140		30	
Isopropylbenzene (Cumene)	ug/L	ND	20	20	23.5	21.3	118	107	60-140		30	
m&p-Xylene	ug/L	ND	40	40	45.2	40.4	113	101	60-140		30	
Methyl-tert-butyl ether	ug/L	0.74	20	20	23.2	20.6	112	99	60-140		30	
Methylene Chloride	ug/L	ND	20	20	21.4	18.5	107	93	60-140	14	30	

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: NCDOT FRONT ST. WBS17BP.3.R.28

Pace Project No.: 92202425

MATRIX SPIKE & MATRIX SP	IKE DUPLICAT	E: 12072	85		1207286							
			MS	MSD								
	922	202441006	Spike	Spike	MS	MSD	MS	MSD	% Rec		Max	
Parameter	Units	Result	Conc.	Conc.	Result	Result	% Rec	% Rec	Limits	RPD	RPD	Qual
n-Butylbenzene	ug/L	ND	20	20	22.8	20.6	114	103	60-140	10	30	
n-Propylbenzene	ug/L	ND	20	20	22.7	20.6	114	103	60-140	10	30	
Naphthalene	ug/L	ND	20	20	22.2	19.1	111	96	60-140	15	30	
o-Xylene	ug/L	ND	20	20	22.9	20.5	115	102	60-140	11	30	
sec-Butylbenzene	ug/L	ND	20	20	23.4	20.9	117	105	60-140	11	30	
Styrene	ug/L	ND	20	20	23.9	21.0	120	105	60-140	13	30	
tert-Butylbenzene	ug/L	ND	20	20	23.4	21.0	117	105	60-140	11	30	
Tetrachloroethene	ug/L	ND	20	20	21.9	19.7	109	99	60-140	10	30	
Toluene	ug/L	ND	20	20	21.8	19.6	109	98	60-140	11	30	
trans-1,2-Dichloroethene	ug/L	ND	20	20	22.1	19.9	110	99	60-140	11	30	
trans-1,3-Dichloropropene	ug/L	ND	20	20	19.5	17.5	98	87	60-140	11	30	
Trichloroethene	ug/L	ND	20	20	22.3	20.1	111	100	60-140	10	30	
Trichlorofluoromethane	ug/L	ND	20	20	24.8	22.2	124	111	60-140	11	30	
Vinyl chloride	ug/L	ND	20	20	28.4	25.3	142	126	60-140	12	30	MO
1,2-Dichloroethane-d4 (S)	%						97	97	70-130			
4-Bromofluorobenzene (S)	%						100	100	70-130			
Toluene-d8 (S)	%						99	100	70-130			

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



Analysis Method:

Project: NCDOT FRONT ST. WBS17BP.3.R.28

Pace Project No.: 92202425

QC Batch:OEXT/27878QC Batch Method:EPA 625

Analysis Description: 625 MSS

EPA 625

Associated Lab Samples: 92202425001

METHOD BLANK: 1205691 Matrix: Water Associated Lab Samples: 92202425001 Blank Reporting Parameter Units Result Limit Qualifiers Analyzed 1,2,4-Trichlorobenzene ug/L ND 5.0 06/02/14 22:08 2,4,6-Trichlorophenol ug/L ND 10.0 06/02/14 22:08 ND 2,4-Dichlorophenol ug/L 5.0 06/02/14 22:08 2,4-Dimethylphenol ug/L ND 06/02/14 22:08 10.0 2,4-Dinitrophenol ug/L ND 50.0 06/02/14 22:08 2,4-Dinitrotoluene ug/L ND 5.0 06/02/14 22:08 2,6-Dinitrotoluene ug/L ND 5.0 06/02/14 22:08 2-Chloronaphthalene ug/L ND 5.0 06/02/14 22:08 2-Chlorophenol ug/L ND 5.0 06/02/14 22:08 2-Nitrophenol ug/L ND 5.0 06/02/14 22:08 3,3'-Dichlorobenzidine ND 25.0 06/02/14 22:08 ug/L ug/L 4.6-Dinitro-2-methylphenol ND 20.0 06/02/14 22:08 ND 06/02/14 22:08 4-Bromophenylphenyl ether ug/L 5.0 ND 06/02/14 22:08 4-Chloro-3-methylphenol ug/L 5.0 4-Chlorophenylphenyl ether ug/L ND 5.0 06/02/14 22:08 4-Nitrophenol ug/L ND 50.0 06/02/14 22:08 Acenaphthene ug/L ND 5.0 06/02/14 22:08 Acenaphthylene ug/L ND 5.0 06/02/14 22:08 Anthracene ND 5.0 06/02/14 22:08 ug/L ND 06/02/14 22:08 Benzo(a)anthracene ug/L 5.0 Benzo(a)pyrene ug/L ND 5.0 06/02/14 22:08 Benzo(b)fluoranthene ug/L ND 5.0 06/02/14 22:08 Benzo(g,h,i)perylene ug/L ND 5.0 06/02/14 22:08 Benzo(k)fluoranthene ND 50 06/02/14 22:08 ug/L bis(2-Chloroethoxy)methane ug/L ND 10.0 06/02/14 22:08 bis(2-Chloroethyl) ether ug/L ND 5.0 06/02/14 22:08 ND bis(2-Chloroisopropyl) ether ug/L 5.0 06/02/14 22:08 bis(2-Ethylhexyl)phthalate ug/L ND 5.0 06/02/14 22:08 Butylbenzylphthalate ug/L ND 5.0 06/02/14 22:08 Chrysene ND 06/02/14 22:08 ug/L 5.0 Di-n-butylphthalate ug/L ND 50 06/02/14 22:08 Di-n-octylphthalate ug/L ND 50 06/02/14 22:08 Dibenz(a,h)anthracene ug/L ND 5.0 06/02/14 22:08 Diethylphthalate ug/L ND 50 06/02/14 22:08 Dimethylphthalate ug/L ND 5.0 06/02/14 22:08 Fluoranthene ug/L ND 5.0 06/02/14 22:08 Fluorene ug/L ND 5.0 06/02/14 22:08 Hexachloro-1,3-butadiene ug/L ND 06/02/14 22:08 5.0 Hexachlorobenzene ug/L ND 5.0 06/02/14 22:08

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.

10.0

5.0

06/02/14 22:08

06/02/14 22:08

ND

ND

REPORT OF LABORATORY ANALYSIS

Hexachlorocyclopentadiene

Hexachloroethane

ug/L

ug/L



Matrix: Water

Project: NCDOT FRONT ST. WBS17BP.3.R.28

Pace Project No.: 92202425

METHOD BLANK: 1205691

Associated Lab Samples: 92202425001

Blank Reporting Parameter Units Result Limit Analyzed Qualifiers ND Indeno(1,2,3-cd)pyrene ug/L 5.0 06/02/14 22:08 Isophorone ug/L ND 10.0 06/02/14 22:08 N-Nitroso-di-n-propylamine ND 5.0 06/02/14 22:08 ug/L 5.0 06/02/14 22:08 ND N-Nitrosodimethylamine ug/L N-Nitrosodiphenylamine ug/L ND 10.0 06/02/14 22:08 Naphthalene ug/L ND 5.0 06/02/14 22:08 Nitrobenzene ug/L ND 5.0 06/02/14 22:08 Pentachlorophenol ug/L ND 10.0 06/02/14 22:08 ug/L Phenanthrene ND 5.0 06/02/14 22:08 Phenol ND 5.0 06/02/14 22:08 ug/L ND Pyrene ug/L 5.0 06/02/14 22:08 2,4,6-Tribromophenol (S) % 64 10-137 06/02/14 22:08 2-Fluorobiphenyl (S) % 65 15-120 06/02/14 22:08 2-Fluorophenol (S) % 34 10-120 06/02/14 22:08 % Nitrobenzene-d5 (S) 64 10-120 06/02/14 22:08 Phenol-d6 (S) % 27 10-120 06/02/14 22:08 % Terphenyl-d14 (S) 80 11-131 06/02/14 22:08

LABORATORY CONTROL SAMPLE: 1205692

		Spike	LCS	LCS	% Rec	
Parameter	Units	Conc.	Result	% Rec	Limits	Qualifiers
1,2,4-Trichlorobenzene	ug/L		19.9	40	44-142	L2
2,4,6-Trichlorophenol	ug/L	50	20.9	42	37-144	
2,4-Dichlorophenol	ug/L	50	18.2	36	1-191	
2,4-Dimethylphenol	ug/L	50	18.2	36	32-119	
2,4-Dinitrophenol	ug/L	250	171	69	1-181	
2,4-Dinitrotoluene	ug/L	50	55.1	110	39-139	
2,6-Dinitrotoluene	ug/L	50	45.2	90	50-158	
-Chloronaphthalene	ug/L	50	19.5	39	60-118 I	L2
-Chlorophenol	ug/L	50	19.5	39	23-134	
-Nitrophenol	ug/L	50	18.9	38	29-182	
,3'-Dichlorobenzidine	ug/L	100	88.7	89	1-262	
,6-Dinitro-2-methylphenol	ug/L	100	83.2	83	1-181	
-Bromophenylphenyl ether	ug/L	50	34.2	68	53-127	
-Chloro-3-methylphenol	ug/L	100	48.2	48	22-147	
-Chlorophenylphenyl ether	ug/L	50	31.8	64	25-158	
-Nitrophenol	ug/L	250	111	44	1-132	
cenaphthene	ug/L	50	23.8	48	47-145	
cenaphthylene	ug/L	50	23.3	47	33-145	
Inthracene	ug/L	50	39.7	79	1-166	
enzo(a)anthracene	ug/L	50	42.2	84	33-143	
Benzo(a)pyrene	ug/L	50	43.2	86	17-163	
Benzo(b)fluoranthene	ug/L	50	41.5	83	24-159	
3enzo(g,h,i)perylene	ug/L	50	42.6	85	1-219	

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: NCDOT FRONT ST. WBS17BP.3.R.28

Pace Project No.: 92202425

LABORATORY CONTROL SAMPLE: 1205692

		Spike	LCS	LCS	% Rec	
Parameter	Units	Conc.	Result	% Rec	Limits	Qualifiers
Benzo(k)fluoranthene	ug/L	50	42.5	85	11-162	
bis(2-Chloroethoxy)methane	ug/L	50	19.9	40	33-184	
bis(2-Chloroethyl) ether	ug/L	50	21.2	42	12-158	
bis(2-Chloroisopropyl) ether	ug/L	50	20.0	40	36-166	
bis(2-Ethylhexyl)phthalate	ug/L	50	47.0	94	8-158	
Butylbenzylphthalate	ug/L	50	47.6	95	1-152	
Chrysene	ug/L	50	43.3	87	17-168	
Di-n-butylphthalate	ug/L	50	44.8	90	1-118	
Di-n-octylphthalate	ug/L	50	45.8	92	4-146	
Dibenz(a,h)anthracene	ug/L	50	44.2	88	1-227	
Diethylphthalate	ug/L	50	46.0	92	1-114	
Dimethylphthalate	ug/L	50	39.7	79	1-112	
Fluoranthene	ug/L	50	41.9	84	26-137	
Fluorene	ug/L	50	33.5	67	59-121	
Hexachloro-1,3-butadiene	ug/L	50	19.8	40	24-116	
Hexachlorobenzene	ug/L	50	35.2	70	1-152	
Hexachlorocyclopentadiene	ug/L	50	14.5	29	25-150	
Hexachloroethane	ug/L	50	19.3	39	40-113 L	.2
Indeno(1,2,3-cd)pyrene	ug/L	50	34.9	70	1-171	
Isophorone	ug/L	50	23.0	46	21-196	
N-Nitroso-di-n-propylamine	ug/L	50	18.7	37	1-230	
N-Nitrosodimethylamine	ug/L	50	12.6	25	25-150	
N-Nitrosodiphenylamine	ug/L	50	36.8	74	25-150	
Naphthalene	ug/L	50	20.4	41	21-133	
Nitrobenzene	ug/L	50	25.2	50	35-180	
Pentachlorophenol	ug/L	100	80.8	81	14-176	
Phenanthrene	ug/L	50	38.0	76	54-120	
Phenol	ug/L	50	9.5	19	5-112	
Pyrene	ug/L	50	43.6	87	52-115	
2,4,6-Tribromophenol (S)	%			72	10-137	
2-Fluorobiphenyl (S)	%			36	15-120	
2-Fluorophenol (S)	%			23	10-120	
Nitrobenzene-d5 (S)	%			40	10-120	
Phenol-d6 (S)	%			16	10-120	
Terphenyl-d14 (S)	%			88	11-131	

MATRIX SPIKE & MATRIX S	PIKE DUPLICAT	E: 12056	93		1205694							
	92	202592004	MS Spike	MSD Spike	MS	MSD	MS	MSD	% Rec		Мах	
Parameter	Units	Result	Conc.	Conc.	Result	Result	% Rec	% Rec	Limits	RPD	RPD	Qual
1,2,4-Trichlorobenzene	ug/L	<5.0	100	100	75.7	68.0	76	68	44-142	11	30	
2,4,6-Trichlorophenol	ug/L	<10.0	100	100	81.4	70.8	81	71	37-144	14	30	
2,4-Dichlorophenol	ug/L	<5.0	100	100	85.7	73.1	86	73	1-191	16	30	
2,4-Dimethylphenol	ug/L	<10.0	100	100	78.4	32.7	78	33	32-119	82	30	R1
2,4-Dinitrophenol	ug/L	<50.0	500	500	426	403	85	81	1-181	5	30	

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REPORT OF LABORATORY ANALYSIS



Project: NCDOT FRONT ST. WBS17BP.3.R.28

Pace Project No.: 92202425

MATRIX SPIKE & MATRIX SP	PIKE DUPLICATE:	12056	93		1205694							
			MS	MSD								
Devenueter		2592004	Spike	Spike	MS	MSD	MS % Data	MSD	% Rec		Max	Qual
Parameter	Units	Result	Conc.	Conc.	Result	Result	% Rec	% Rec	Limits	RPD		Qual
2,4-Dinitrotoluene	ug/L	<5.0	100	100	104	103	104	103	39-139	0	30	
2,6-Dinitrotoluene	ug/L	<5.0	100	100	104	97.2	104	97	50-158	7	30	
2-Chloronaphthalene	ug/L	<5.0	100	100	82.9	73.8	83	74	60-118	12	30	
2-Chlorophenol	ug/L	<5.0	100	100	89.9	70.5	90	71	23-134	24	30	
2-Nitrophenol	ug/L	<5.0	100	100	84.6	70.5	85	71	29-182	18	30	
3,3'-Dichlorobenzidine	ug/L	<25.0	200	200	158	140	79	70	1-262	12	30	
4,6-Dinitro-2-methylphenol	ug/L	<20.0	200	200	180	170	90	85	1-181	6	30	
4-Bromophenylphenyl ether	ug/L	<5.0	100	100	85.7	81.6	86	82	53-127	5	30	
4-Chloro-3-methylphenol	ug/L	<5.0	200	200	181	153	90	76	22-147	17	30	
4-Chlorophenylphenyl ether	ug/L	<5.0	100	100	87.6	84.7	88	85	25-158	3	30	
4-Nitrophenol	ug/L	<50.0	500	500	310	285	62	57	1-132	8	30	
Acenaphthene	ug/L	<5.0	100	100	84.8	77.8	83	76	47-145	9	30	
Acenaphthylene	ug/L	<5.0	100	100	82.2	76.0	82	76	33-145	8	30	
Anthracene	ug/L	<5.0	100	100	85.2	83.1	85	83	1-166	2	30	
Benzo(a)anthracene	ug/L	<5.0	100	100	87.9	81.3	88	81	33-143	8	30	
Benzo(a)pyrene	ug/L	<5.0	100	100	91.5	87.6	91	88	17-163	4	30	
Benzo(b)fluoranthene	ug/L	<5.0	100	100	88.9	86.8	89	87	24-159	2	30	
Benzo(g,h,i)perylene	ug/L	<5.0	100	100	86.0	81.6	86	82	1-219	5	30	
Benzo(k)fluoranthene	ug/L	<5.0	100	100	92.2	87.6	92	88	11-162	5	30	
bis(2-Chloroethoxy)methane	ug/L	<10.0	100	100	86.0	75.6	86	76	33-184	13	30	
bis(2-Chloroethyl) ether	ug/L	<5.0	100	100	93.2	74.7	93	75	12-158	22	30	
bis(2-Chloroisopropyl) ether	ug/L	<5.0	100	100	90.0	73.3	90	73	36-166	21	30	
bis(2-Ethylhexyl)phthalate	ug/L	<5.0	100	100	95.8	90.6	96	91	8-158	6	30	
Butylbenzylphthalate	ug/L	<5.0	100	100	97.7	89.5	98	89	1-152	9	30	
Chrysene	ug/L	<5.0	100	100	88.9	83.5	89	84	17-168	6	30	
Di-n-butylphthalate	ug/L	<5.0	100	100	89.0	90.8	89	91	1-118	2	30	
Di-n-octylphthalate	ug/L	<5.0	100	100	92.4	85.1	92	85	4-146	8	30	
Dibenz(a,h)anthracene	ug/L	<5.0	100	100	88.7	84.1	89	84	1-227	5	30	
Diethylphthalate	ug/L	<5.0	100	100	87.2	88.4	87	88	1-227	1	30	
Dimethylphthalate	•	<5.0	100	100	86.7	81.5	87	81	1-112	6	30	
Fluoranthene	ug/L ug/L	<5.0 <5.0	100	100	83.5	86.1	83	86	26-137	3	30	
Fluorene		<5.0 <5.0	100	100	86.2	84.2	84	82	20-137 59-121	2	30	
	ug/L						04 72					
Hexachloro-1,3-butadiene Hexachlorobenzene	ug/L	<5.0 <5.0	100 100	100 100	71.8 78.4	64.7 74.0	72	65 74	24-116 1-152	10 6	30 30	
	ug/L						80					
Hexachlorocyclopentadiene	ug/L	<10.0	100 100	100 100	80.5 83.3	65.5 72.1	80 83	65 72	25-150 40-113	21 14	30 30	
Hexachloroethane	ug/L	<5.0						72				
Indeno(1,2,3-cd)pyrene	ug/L	<5.0	100	100	68.8	65.2	69	65	1-171	5	30	
Isophorone	ug/L	<10.0	100	100	100	87.8	100	88	21-196	13	30	
N-Nitroso-di-n-propylamine	ug/L	<5.0	100	100	107	79.8	107	80	1-230	29	30	
N-Nitrosodimethylamine	ug/L	<5.0	100	100	64.8	53.6	65	54	25-150	19	30	
N-Nitrosodiphenylamine	ug/L	<10.0	100	100	87.6	81.6	88	82	25-150	7	30	
Naphthalene	ug/L	<5.0	100	100	82.5	74.5	78	70	21-133	10	30	
Nitrobenzene	ug/L	<5.0	100	100	97.9	84.5	98	84	35-180	15	30	
Pentachlorophenol	ug/L	<10.0	200	200	169	161	85	81	14-176	5	30	
Phenanthrene	ug/L	<5.0	100	100	84.4	82.3	82	80	54-120	3	30	
Phenol	ug/L	<5.0	100	100	61.4	43.6	61	44	5-112	34	30	K1

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: NCDOT FRONT ST. WBS17BP.3.R.28

Pace Project No.: 92202425

MATRIX SPIKE & MATRIX SP	PIKE DUPLICAT	E: 12056	93		1205694							
Parameter	92: Units	202592004 Result	MS Spike Conc.	MSD Spike Conc.	MS Result	MSD Result	MS % Rec	MSD % Rec	% Rec Limits	RPD	Max RPD	Qual
Pyrene	ug/L	<5.0	100	100	91.2	82.8	91	83	52-115	10	30	
2,4,6-Tribromophenol (S)	%						82	82	10-137			
2-Fluorobiphenyl (S)	%						81	75	15-120			
2-Fluorophenol (S)	%						58	49	10-120			
Nitrobenzene-d5 (S)	%						77	72	10-120			
Phenol-d6 (S)	%						57	41	10-120			
Terphenyl-d14 (S)	%						86	84	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.



QUALIFIERS

Project: NCDOT FRONT ST. WBS17BP.3.R.28

Pace Project No.: 92202425

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

- L2 Analyte recovery in the laboratory control sample (LCS) was below QC limits. Results for this analyte in associated samples may be biased low.
- 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:NCDOT FRONT ST. WBS17BP.3.R.28Pace Project No.:92202425

Lab ID	Sample ID	QC Batch Method	QC Batch	Analytical Method	Analytical Batch
92202425001	3-08	EPA 625	OEXT/27878	EPA 625	MSSV/9170
92202425001	3-08	SM 6200B	MSV/26940		

	Documen		Document Revised: April 07. 2014
Pace Analytical*	Sample Condition Up Document		Page 1 of 2 Issuing Authority:
	F-CHR-CS-		Pace Huntersville Quality Office
Client Name: Catlin	NCDUT		
	PS Client Comme	ercial Pace Other_	Optional
Custody Seal on Cooler/Box Presen	it: 🗌 yes 🖄 no	Seals intact: yes	no Proj. Due Date: Proj. Name:
Packing Material:	Bubble Bags 🗌 Nor	ne 🗌 Other	
Thermometer Used: IR Gun T1102	T1401 Type of Ice:	Wet Blue None	Samples on ice, cooling process has begun
	: No Correction T1:	301: No Correction	Date and Initials of person examining
Corrected Cooler Temp.: 2.1	°C Biological T	issue is Frozen: Yes No	$_{\text{N/A}}$ contents: $_{\text{A}}$ $_{\text{S}}$ $_{\text{S}}$ $_{\text{C}}$ $_{\text{C}}$ $_{\text{C}}$ $_{\text{C}}$ $_{\text{C}}$
Temp should be above freezing to 6°C		Comments:	0000 51001
Chain of Custody Present:	Yes No	□n/A 1.	
Chain of Custody Filled Out:	Xyes 🗆 No	□n/a 2.	
Chain of Custody Relinquished:	'ØYes □No	□n/a 3.	
Sampler Name & Signature on COC:	XYes □No	□n/a 4.	
Samples Arrived within Hold Time:	XYes □No	□n/a 5.	
Short Hold Time Analysis (<72hr):	□Yes ØÅNo	□n/a 6.	
Rush Turn Around Time Requested	: 🛛 Yes 🕅 No	□n/a 7.	
Sufficient Volume:	⊠Yes □No	□n/a 8.	
Correct Containers Used:	\$\$Yes □No	□n/a 9.	
-Pace Containers Used:	□Yes ¤No	□n/a	
Containers Intact:	`⊠(Yes □No	□n/a 10.	
Filtered volume received for Dissolved	d tests □Yes □No	QN/A 11.	
Sample Labels match COC:		IN/A 12. Samples	From 10 3-08 do nut
-Includes date/time/ID/Analysis	Matrix:	have date	
All containers needing preservation have bee		XIN/A 13.	
All containers needing preservation are fou compliance with EPA recommendation.			
exceptions: VOA, coliform, TOC, O&G, WI-DRO	(water) Øvyes □No		
Samples checked for dechlorination:	Yes DNo	□n/A 14.	
Headspace in VOA Vials (>6mm):	TYes WNO	□n/A 15.	
Trip Blank Present:	□Yes □No	XN/A 16.	
Trip Blank Custody Seals Present	□Yes □No	₩N/A	
Pace Trip Blank Lot # (if purchased):_			
Client Notification/ Resolution:			Field Data Required? Y / N
Person Contacted:		Date/Time:	
Comments/ Resolution:			
SCURF Review: JDB	Date: 52214	1.10#	92202425
SRF Review: AMB	Date: 5-23-14	WOT	
Note: Whenever there is a discrepancy a samples, a copy of this form will be se Certification Office (i.e out of hold, in	ent to the North Carolina DEHM	NR	

incorrect containers)

FIGURATION Note of an operation of a second construction of a second constr							p		12	11	10	9	8	7	6	UI	4	ω	N	-	ITEM #			Req	Pho	S ^m	5	Add.	Corr	Sec	
Inter our construction I	*Important Note: By signing this form you are acc	0					J Values	ADDITIONAL COMMENTS												0	SAMPLE ID (A-Z, 0-9 /,-) Sample IDS MUST BE UNIQUE	7		1		all To: 24. ash ba @ attinusa.com	ton.NC		2	Section A Required Client Information:	Pace Analytical®
Inter our construction I	cepting Pa	ORIGI					Sers										25					Of de				N.a.		Сору	Repo	Sect	
Inter our construction I	ce's NET	NAL				A		-	_		Andreastan				-			-		8		Ś		S S Numb	ZName	hase Or		-	ort To: 1	tion B	
Inter our construction I	30 day p					Z		RELINO	-			-	-				-	-	-	,			-)er:	10	P	4	0	Re	piect Inf	
Inter our construction I	payment terms and a		S/		•	esent fa		UISHED BY / AF													COMPOSIT START DATE			0,	Front	+ Ne	-		Ashba	ormation:	
Inter our construction I	agreeing t		AMPLE			0		FILIATIO											. 1	X	TIME	COLLE			St				-		
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