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www.catlinusa.com

December 31, 2008

NAVFAC Mid-Atlantic
Marine Corps North Carolina IPT
Environmental Business Line
Code: OPCEV3MA
Attn: Mr. Melvin Acree
6506 Hampton Boulevard
Building C, Room 314
Norfolk, VA 23508-1278

Re: **Site BB-9**
Groundwater Sampling
Marine Corps Base
Camp Lejeune, North Carolina
Navy Contract No. N62470-05-D-6200
Delivery Order No. 0016
CATLIN Project No. 205-077

Dear Mr. Acree:

In September of 2006 a Report of Findings concerning this site was submitted to the Wilmington Regional Office (WiRO) of the North Carolina Department of Environment and Natural Resource (NCDENR) detailing recent soil sampling results and requesting the site be considered for No Further Action (NFA). In an e-mail dated November 1, 2006, Mr. Bruce Reed of the WiRO requested that a groundwater sample be collected from monitoring well BB9-3 in order to justify the No Further Action request. Based on this request CATLIN installed a temporary well (BB9 TW01) on November 20, 2006, because the monitoring well BB9-3 could not be located. A groundwater sample was obtained and analyzed per Environmental Protection Agency (EPA) Method 625 and Massachusetts Department of Environmental Protection (MADEP) VPH and EPH. Results of the laboratory analysis are summarized in Tables 1 through 3. Laboratory results indicated three Method 625 compounds, Benzo[k]fluoranthene, Chrysene, and 1,4-Dichlorobenzene and one MADEP VPH/EPH hydrocarbon fraction, C9-C22 Aromatics above North Carolina Administrative Code (NCAC) 2L Groundwater Quality Standards (GWQS) or Gross Contaminant Levels (GCLs) that prevented the site from being considered for NFA. It was recommended that the groundwater be allowed to naturally attenuate and the site be re-sampled in two years.

CATLIN mobilized to the site on November 14, 2008 to install a permanent monitoring well to replace previously destroyed BB9-3 and collect a representative groundwater sample. Using a trailer mounted Diedrich D-50 drilling rig, CATLIN installed a boring and set a new monitoring well (USTBB9-MW03R). The monitoring well was developed on the same day and later sampled on November 17, 2008. A Well Construction Record and Well Log were submitted to the State on November 24, 2008 and have also been

included with this letter. Monitoring well USTBB9-MW03R was installed in the location of previous well BB9-3 (see attached Figure 1). The groundwater sample was placed in appropriate laboratory provided glassware and transported under proper Chain-of-Custody to SGS Environmental Services, Inc. of Wilmington, North Carolina for analysis by EPA Method 625 and MADEP VPH and EPH. Results of the November 17, 2008 laboratory analyses are summarized in Tables 1-3 and as follows:

EPA METHOD 625

Several EPA Method 625 constituents were detected at low concentrations in excess of laboratory method detection limits. However only Benzo[a]anthracene (23.4 ug/L), Benzo[a]pyrene (14.4 ug/L), Benzo[b]fluoranthene (20.6 ug/L), Benzo[k]fluoranthene (9.5 ug/L), and Chrysene (24.0 ug/L) were above applicable 2L GWQSs and Benzo[a]anthracene, Benzo[b]fluoranthene, and Benzo[k]fluoranthene were above the GCLs. See attached laboratory summary on Table 1. Table 1 also includes laboratory results from the groundwater sample collected from temporary monitoring well BB-9 TW01 in December 2006.

MADEP VPH/EPH

Massachusetts Department of Environmental Protection (MADEP) VPH and EPH fractions were also detected in the USTBB9-MW03R groundwater sample. Only the C9-C22 Aromatic concentration of <4,610 ug/L was in excess of the 2L GWQS of 210 ug/L. There are no established GCLs for the MADEP fractions. See attached laboratory summary on Tables 2 and 3. Tables 2 and 3 also include laboratory results from the groundwater sample collected from temporary monitoring well BB-9 TW01 in December 2006.

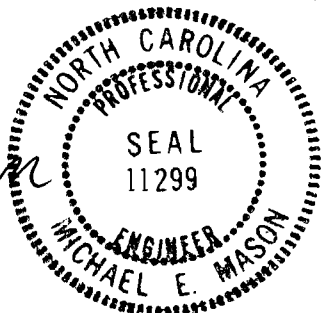
Laboratory reports from both sampling events have been attached for your reference along with Chain of Custody documentation.

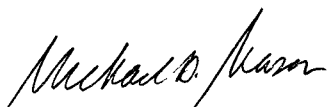
Based on a review of the laboratory results, there are contaminant concentrations in excess of the GCLs that prevent the site from being considered for No Further Action at this time. It is recommended that the groundwater at the site be allowed to naturally attenuate and the site be re-sampled in two years to determine if the contaminant concentrations have been reduced to below acceptable levels.

CATLIN Engineers and Scientists appreciate the opportunity to continue to provide services to NAVFAC Mid-Atlantic and the MCB on your environmental projects.

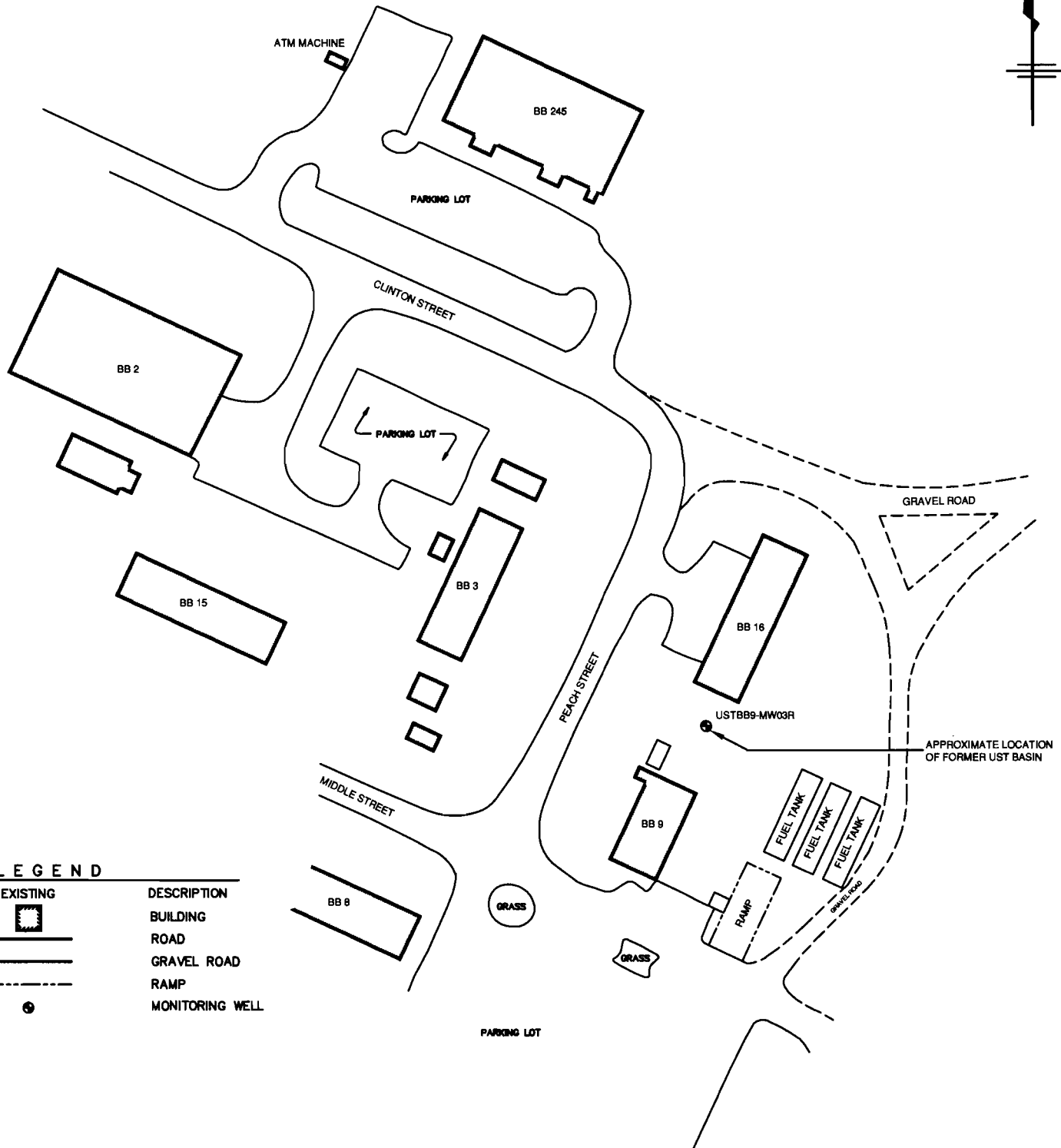
Sincerely,


Michael E. Mason, P.E.
Program Manager




Michael D. Mason
Project Engineer

Attachments



LEGEND

EXISTING	DESCRIPTION
	BUILDING
	ROAD
	GRAVEL ROAD
	RAMP
	MONITORING WELL

NOTE:
 1). FIGURE MODIFIED FROM BAKER ENVIRONMENTAL INC. SITE PLAN
 2). GROUNDWATER MONITORING WELL LOCATION IS APPROXIMATE.



 220 Old Dairy Road Wilmington, NC 28405	PROJECT BB-9 GROUNDWATER SAMPLING MARINE CORP. BASE CAMP LEJEUNE, NC	TITLE BB-9 SITE PLAN WITH MONITORING WELL LOCATION		FIGURE 1
	JOB NO: 205-077	DATE: DEC 2008	SCALE: 1" = 100'	DRAWN BY: MDM

TABLE 1
SUMMARY OF GROUNDWATER LABORATORY RESULTS
EPA METHOD 625

Incident Name and No.: Site BB-9

Well ID	Contaminant of Concern →		Acenaphthene	Anthracene	Benzo[a]anthracene	Benzo[a]pyrene	Benzo[b]fluoranthene	Benzo[g,h,i]perylene	Benzo[k]fluoranthene	Chrysene
	Sample ID	Date Collected								
GCL			2,120	2,100	22	15	0.6	210	0.47	0.8
2L GWQS			80	2,100	0.0479	0.00479	0.0479	210	0.479	4.79
BB-9 TW01	BB-9 TW01	12/8/2006	<1.22	<1.75	<1.36	<1.27	<1.43	<4.57	1.30 J	1.50 J
USTBB9-MW03R	USTBB9-MW03R	11/17/2008	38.6	21.2	23.4	14.4	20.6	9.23	9.5	24.0

BQL = Below Quantitation Limit

All results in micrograms per Liter (ug/L).

NE = Not Established

Bold results indicate concentrations above 2L GWQS or GCL.

< = less than method detection limit

J = Estimated concentration, below calibration range and above Method Detection Limit

NA = Not Analyzed

TABLE 1
SUMMARY OF GROUNDWATER LABORATORY RESI
EPA METHOD 625

Incident Name and No.: Site BB-9

Well ID	Contaminant of Concern →		1,2-Dichlorobenzene	1,4-Dichlorobenzene	Fluoranthene	Fluorene	Phenanthrene	Pyrene	All Other 625 Compounds
	Sample ID	Date Collected							
GCL			24,000	39,500	280	950	410	210	Varies
2L GWQS			24	1.4	280	280	210	210	Varies
BB-9 TW01	BB-9 TW01	12/8/2006	4.40 J	4.20 J	1.70 J	<1.22	1.80 J	2.90 J	BMDL
USTBB9-MW03R	USTBB9-MW03R	11/17/2008	NA	NA	75.2	12.3	12.1	64.7	BMDL

BQL = Below Quantitation Limit

All results in micrograms per Liter (ug/L).

NE = Not Established

Bold results indicate concentrations above 2L GWQS or GCL.

< = less than method detection limit

J = Estimated concentration, below calibration range and above Metho

NA = Not Analyzed

TABLE 2
SUMMARY OF GROUNDWATER LABORATORY RESULTS
MADEP METHOD VPH/EPH

Incident Name and No.: Site BB-9

Well ID	Hydrocarbon Fraction		VPH			EPH		
			C5-C8 Aliphatics	C9-C12 Aliphatics	C9-C10 Aromatics	C9-C18 Aliphatics	C19-C36 Aliphatics	C11-C22 Aromatics
	Sample ID	Date Collected						
BB-9 TW01	BB-9 TW01	12/8/2006	<100	<100	<100	420	1,500	240
USTBB9-MW03R	USTBB9-MW03R	11/17/2008	<100	<100	<100	833	2,870	4,510

All results in micrograms per Liter (ug/L).

TABLE 3
SUMMARY OF GROUNDWATER LABORATORY RESULTS
MADEP METHOD VPH/EPH AS COMPARED TO NCDENR 2L GWQS

Incident Name and No.: Site BB-9

Well ID	Hydrocarbon Fraction of Concern →		C5-C8 Aliphatics	C9-C18 Aliphatics	C19-C36 Aliphatics	C9-C22 Aromatics
	Sample ID	Date Collected				
GCL 2L GWQS			NE 420	NE 4,200	NE 42,000	NE 210
BB-9 TW01	BB-9 TW01	12/8/2006	<100	<520	1,500	<340*
USTBB9-MW03R	USTBB9-MW03R	11/17/2008	<100	<933*	2,870	<4,610*

All results in micrograms per Liter (ug/L).

NE = Not Established

Bold results indicate concentrations above 2L GWQS or GCL.

* = The value represents the sum of the reported practical quantitation limit of one fraction and the detected concentration of the other fraction.

WELL LOG

CATLIN




ENGINEERS and SCIENTISTS
205-077
Wilmington, NC

SHEET 1 OF 1

PROJECT NO.: 205-077	STATE: NC	COUNTY: Onslow	LOCATION: Camp Lejeune
PROJECT NAME: BB-9		LOGGED BY: Tom Stetler	WELL ID: USTBB9-MW03R
NORTHING:		DRILLER: Bobbie D. Fowler	
EASTING:		CREW: John Wood	
SYSTEM:		BORING LOCATION: Proximity of former well BB9-3.	T.O.C. ELEV.:
DRILL MACHINE: Diedrich D-50	METHOD: HSA	0 HOUR DTW: 5.5	TOTAL DEPTH: 9.0
START DATE: 11/14/08	FINISH DATE: 11/14/08	24 HOUR DTW: 5.3	WELL DEPTH: 9.0

DEPTH	BLOW COUNT				OVA (ppm)	LAB.	U S C S	L O G	SOIL AND ROCK DESCRIPTION	WELL DETAIL
	6in	6in	6in	6in						
0.0									LAND SURFACE	0.0
1.5									White limestone gravel in auger cuttings	2" Sch. 40 PVC
3.0	1	1	1	4	0.0		SM		Dark brown, SILTY vf. SAND. Moist. Wet at base of interval. Tr. mod. rounded limestone gravel. Fill.	2" Sch. 40 PVC
5.0										2" Slot 010 Sch. 40 PVC
8.0	6	50/5"			1884		SM		Same as above. Refusal at 9.0' BLS on concrete. Dark brown, SILTY vf. SAND w/ tr. gravel. Strong HCO and petroleum sheen on sample. Saturated.	2" Slot 010 Sch. 40 PVC
9.0									Boring Terminated at Depth 9.0 ft on concrete.	9.0

CATLIN BORING LOG 205-077_BB-9.GPJ CATLIN.GDT 12/16/08

 Portland Cement
  Bentonite Pellets
  #2 Medium Sand



Mr. Mike E. Mason
Richard Catlin & Associates
P.O. Box 10279
Wilmington NC 28404-0279

Report Number: G128-1872

Client Project: 205-077 (BB-9)

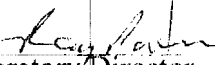
Dear Mr. Mason:

Enclosed are the results of the analytical services performed under the referenced project. The samples are certified to meet the requirements of the National Environmental Laboratory Accreditation Conference Standards. Copies of this report and supporting data will be retained in our files for a period of five years in the event they are required for future reference. Any samples submitted to our laboratory will be retained for a maximum of thirty (30) days from the date of this report unless other arrangements are requested.

If there are any questions about the report or the services performed during this project, please call SGS/Paradigm at (910) 350-1903. We will be happy to answer any questions or concerns which you may have.

Thank you for using SGS/Paradigm Analytical Labs for your analytical services. We look forward to working with you again on any additional analytical needs which you may have.

Sincerely,
SGS/Paradigm Analytical Laboratories, Inc.


Laboratory Director
J. Patrick Weaver

12/21/2006
Date



Results for Semivolatiles
by GCMS 625

Client Sample ID: BB-9 TW01
Client Project ID: 205-077 (BB-9)
Lab Sample ID: G128-1872-1E
Lab Project ID: G128-1872

Analyzed By: EAW
Date Collected: 12/8/2006 12:00
Date Received: 12/11/2006
Date Extracted: 12/12/2006
Matrix: Water

Compound	Result ug/L	RL ug/L	MDL ug/L	Dilution Factor	Date Analyzed	Flag
Acenaphthene	BQL	10.0	1.22	1	12/13/2006	
Acenaphthylene	BQL	10.0	1.12	1	12/13/2006	
Anthracene	BQL	10.0	1.75	1	12/13/2006	
Benzo[a]anthracene	BQL	10.0	1.36	1	12/13/2006	
Benzo[a]pyrene	BQL	10.0	1.27	1	12/13/2006	
Benzo[b]fluoranthene	BQL	10.0	1.43	1	12/13/2006	
Benzo[g,h,i]perylene	BQL	10.0	4.57	1	12/13/2006	
Benzo[k]fluoranthene	1.30	10.0	1.09	1	12/13/2006	J
Bis(2-chloroethoxy)methane	BQL	10.0	1.11	1	12/13/2006	
Bis(2-chloroethyl)ether	BQL	10.0	1.09	1	12/13/2006	
Bis(2-chloroisopropyl)ether	BQL	10.0	1.57	1	12/13/2006	
Bis(2-ethylhexyl)phthalate	BQL	10.0	1.33	1	12/13/2006	
4-bromophenyl phenyl ether	BQL	10.0	1.99	1	12/13/2006	
Butylbenzylphthalate	BQL	10.0	1.53	1	12/13/2006	
2-Chloronaphthalene	BQL	10.0	1.25	1	12/13/2006	
2-Chlorophenol	BQL	10.0	4.22	1	12/13/2006	
4-Chloro-3-methylphenol	BQL	10.0	3.26	1	12/13/2006	
4-Chlorophenyl phenyl ether	BQL	10.0	1.42	1	12/13/2006	
Chrysene	1.50	10.0	1.11	1	12/13/2006	J
Dibenzo[a,h]anthracene	BQL	10.0	4.87	1	12/13/2006	
Di-n-Butylphthalate	BQL	10.0	1.65	1	12/13/2006	
1,2-Dichlorobenzene	4.40	10.0	1.25	1	12/13/2006	J
1,3-Dichlorobenzene	BQL	10.0	1.24	1	12/13/2006	
1,4-Dichlorobenzene	4.20	10.0	1.20	1	12/13/2006	J
3,3'-Dichlorobenzidine	BQL	20.0	4.10	1	12/13/2006	
2,4-Dichlorophenol	BQL	10.0	3.75	1	12/13/2006	
Diethylphthalate	BQL	10.0	1.48	1	12/13/2006	
Dimethylphthalate	BQL	10.0	1.04	1	12/13/2006	
2,4-Dimethylphenol	BQL	10.0	9.25	1	12/13/2006	
Di-n-octylphthalate	BQL	10.0	1.16	1	12/13/2006	
4,6-Dinitro-2-methylphenol	BQL	50.0	3.71	1	12/13/2006	
2,4-Dinitrophenol	BQL	50.0	4.20	1	12/13/2006	
2,4-Dinitrotoluene	BQL	10.0	1.52	1	12/13/2006	
2,6-Dinitrotoluene	BQL	10.0	1.41	1	12/13/2006	
Diphenylamine *	BQL	10.0	1.53	1	12/13/2006	
Fluoranthene	1.70	10.0	1.41	1	12/13/2006	J
Fluorene	BQL	10.0	1.22	1	12/13/2006	
Hexachlorobenzene	BQL	10.0	1.22	1	12/13/2006	
Hexachlorobutadiene	BQL	10.0	1.58	1	12/13/2006	
Hexachlorocyclopentadiene	BQL	20.0	20.0	1	12/13/2006	
Hexachloroethane	BQL	10.0	1.58	1	12/13/2006	
Indeno(1,2,3-c,d)pyrene	BQL	10.0	4.57	1	12/13/2006	
Isophorone	BQL	10.0	1.27	1	12/13/2006	
Naphthalene	BQL	10.0	1.08	1	12/13/2006	
Nitrobenzene	BQL	10.0	1.32	1	12/13/2006	
2-Nitrophenol	BQL	10.0	3.52	1	12/13/2006	
4-Nitrophenol	BQL	50.0	3.17	1	12/13/2006	
N-Nitrosodi-n-propylamine	BQL	10.0	1.87	1	12/13/2006	
Pentachlorophenol	BQL	50.0	2.83	1	12/13/2006	
Phenanthrene	1.80	10.0	1.38	1	12/13/2006	J



Results for Semivolatiles
by GCMS 625

Client Sample ID: BB-9 TW01
Client Project ID: 205-077 (BB-9)
Lab Sample ID: G128-1872-1E
Lab Project ID: G128-1872

Analyzed By: EAW
Date Collected: 12/8/2006 12:00
Date Received: 12/11/2006
Date Extracted: 12/12/2006
Matrix: Water

Compound	Result ug/L	RL ug/L	MDL ug/L	Dilution Factor	Date Analyzed	Flag
Phenol	BQL	10.0	3.38	1	12/13/2006	
Pyrene	2.90	10.0	2.08	1	12/13/2006	J
1,2,4-Trichlorobenzene	BQL	10.0	1.33	1	12/13/2006	
2,4,6-Trichlorophenol	BQL	10.0	2.92	1	12/13/2006	
		Spike Added	Spike Result	Percent Recovered		
2-Fluorobiphenyl		10	6.6	66		
2-Fluorophenol		10	8.7	87		
Nitrobenzene-d5		10	8.9	89		
Phenol-d6		10	9	90		
2,4,6-Tribromophenol		10	8.8	88		
4-Terphenyl-d14		10	5.4	54		

Comments:

* N-Nitrosodiphenylamine is reported as the breakdown product Diphenylamine.

Flags:

BQL = Below Quantitation Limits.
J = Detected below the quantitation limit.

Reviewed By: full



VPH (Aliphatics/Aromatics) Laboratory Reporting Form

Client Name: Richard Catlin & Associates

Project Name: 205-077 (BB-9)

Sample Information and Analytical Results	
Sample Identification	BB-9 TW01
Sample Matrix	Water
Collection Option (for Soil)*	
Date Collected	12/08/06
Date Received	12/11/06
Date Extracted	12/11/06
Date Analyzed	12/11/06
Dry Weight	
Dilution Factor	1
C ₅ -C ₈ Aliphatics**	< 100 (µg/L)
C ₉ -C ₁₂ Aliphatics**	< 100 (µg/L)
C ₉ -C ₁₀ Aromatics**	< 100 (µg/L)
Surrogate % Recovery - PID	110
Surrogate % Recovery - FID	100

* = Option 1 = Established fill line on vial, Option 2 = Sampling Device/Brand, or Option 3 = Field weight of soil.

** = Excludes any surrogates or internal standards.

Lab Info: g128-1872-1b

Reviewed By: me



Attachment 2

VPH Laboratory Reporting Form

Calibration and QA/QC Information

FID Initial Calibration Date: 11/21/06 PID Initial Calibration Date: 11/21/06

Calibration Ranges and Limits

Range	MDL (07/15/2004) (µg/L)	ML (µg/L)	RL	
			(µg/L)	(mg/Kg)
C ₅ -C ₈ Aliphatics	4.4	14	100	10
C ₉ -C ₁₂ Aliphatics	3.4	11	100	10
C ₉ -C ₁₀ Aromatics	0.13	0.41	100	10

Calibration Concentration Levels

Range	Levels (µg/L)	%RSD or CCC	Method of Quantitation
C ₅ -C ₈ Aliphatics	40	21.8	Calibration Factor
	1000		
	2000		
	3000		
	4000		
C ₉ -C ₁₂ Aliphatics	10	24.00	Calibration Factor
	250		
	500		
	750		
	1000		
C ₉ -C ₁₀ Aromatics	10	0.99	Linear Regression
	250		
	500		
	750		
	1000		

Calibration Check Date: 12/11/06

Calibration Check

Range	Levels (µg/L) (mg/Kg)		RPD
	C ₅ -C ₈ Aliphatics	2000	
C ₉ -C ₁₂ Aliphatics	500	50	1.1
C ₉ -C ₁₀ Aromatics	500	50	-12.1

MDL = Method Detection Limit
ML = Minimum Limit
RL = Reportable Limit

RPD = Relative Percent Difference
%RSD = Percent Relative Standard Deviation
CCC = Correlation Coefficient of Curve



EPH (Aliphatics/Aromatics) Results

by MDEP-EPH

Client Name: Richard Catlin & Associates

Project Name: 205-077 (BB-9)

Sample Information and Analytical Results	
Sample Identification	BB-9 TW01
Sample Matrix	Water
Date Collected	12/08/06
Date Received	12/11/06
Date Extracted	12/14/06
Date Analyzed	12/18/06
Dry Weight	
Dilution Factor	1:1
C ₉ -C ₁₈ Aliphatics*	420 (ug/L)
C ₁₉ -C ₃₆ Aliphatics*	1500 (ug/L)
C ₁₁ -C ₂₂ Aromatics*	240 (ug/L)
Aliphatic Surrogate % Recovery	46
Aromatic Surrogate % Recovery	40
Fractionation Surrogate 1 % Recovery	85

Comments:

* = Excludes any surrogates or internal standards.

Lab info: G128-1872-1F

Reviewed By: RM



EPH Laboratory Reporting Form

Calibration and QA/QC Information
--

Initial Calibration Date: 12/11/06**Calibration Ranges and Limits**

Range	MDL (2/2004) ($\mu\text{g/L}$)	ML ($\mu\text{g/L}$)	RL ($\mu\text{g/L}$)	RL (mg/Kg)
C ₉ -C ₁₈ Aliphatics	3.84	12.2	100	10
C ₁₉ -C ₃₆ Aliphatics	0.57	1.8	100	10
C ₁₁ -C ₂₂ Aromatics	4.54	14.4	100	10

Calibration Concentration Levels

Range	Levels ($\mu\text{g/mL}$)	%RSD or CCC	Method of Quantitation
C ₉ -C ₁₈ Aliphatics	6	1.00	Linear Regression
	30		
	60		
	120		
	240		
C ₁₉ -C ₃₆ Aliphatics	8	1.0	Linear Regression
	40		
	80		
	160		
	320		
C ₁₁ -C ₂₂ Aromatics	17	1	Linear Regression
	85		
	170		
	340		
	680		

Calibration Check Date: 12/18/06**Calibration Check**

Range	Levels ($\mu\text{g/mL}$)	RPD
C ₉ -C ₁₈ Aliphatics	120	-17.5
C ₁₉ -C ₃₆ Aliphatics	160	-15.8
C ₁₁ -C ₂₂ Aromatics	340	-0.2

MDL = Method Detection Limit

ML = Minimum Limit

RL = Reportable Limit

RPD = Relative Percent Difference

%RSD = Percent Relative Standard Deviation

CCC = Correlation Coefficient of Curve



List of Reporting Abbreviations and Data Qualifiers

B = Compound also detected in batch blank

BQL = Below Quantitation Limit (RL or MDL)

DF = Dilution Factor

Dup = Duplicate

D = Detected, but RPD is > 40% between results in dual column method.

E = Estimated concentration, exceeds calibration range.

J = Estimated concentration, below calibration range and above MDL

LCS(D) = Laboratory Control Spike (Duplicate)

MDL = Method Detection Limit

MS(D) = Matrix Spike (Duplicate)

PQL = Practical Quantitation Limit

RL = Reporting Limit

RPD = Relative Percent Difference

mg/kg = milligram per kilogram, ppm, parts per million

ug/kg = micrograms per kilogram, ppb, parts per billion

mg/L = milligram per liter, ppm, parts per million

ug/L = micrograms per liter, ppb, parts per billion

% Rec = Percent Recovery

% solids = Percent Solids

Special Notes:

- 1) Metals and mercury samples are digested with a hot block, see the standard operating procedure document for details.
- 2) Uncertainty for all reported data is less than or equal to 30 percent.



CHAIN OF CUSTODY RECORD
SGS Environmental Services Inc.

Locations Nationwide
 • Alaska • Hawaii
 • Louisiana • Maryland
 • New Jersey • North Carolina
 • West Virginia

www.us.sgs.com 0668



CLIENT: **DOD Lejuene**

CONTACT: **Mike E. Mason** PHONE NO.: **(910) 352-7741**

PROJECT: **206-077 (BB-9)** SITE/PWSID#:

REPORTS TO: **Mike E. Mason** E-MAIL:

VOICE TO: **Catlin** QUOTE #:

P.O. NUMBER **261211-3**

SGS Reference: **G128-1872**

PAGE _____ OF _____

LAB NO.	SAMPLE IDENTIFICATION	DATE	TIME	MATRIX	CONTAINERS	SAMPLE TYPE	Preservatives Used										REMARKS							
							Analysis Required	C=COMP	G=GRAB	3	625	EPH	VPH	HCL	HCL									
	BB-9 TW01	12/8/06	1200	G/W	4	G	1	1	2															

Collected/Relinquished By: (1) **Justin Hester** Date: **12/11** Time: **10:55** Received By: **John Flinn**

Shipping Carrier: _____ Samples Received Cold? (Circle) YES NO

Shipping Ticket No: _____ Temperature (C): **5.8°C**

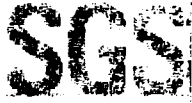
Relinquished By: (2) _____ Date: _____ Time: _____ Received By: _____

Special Deliverable Requirements: _____ Chain of Custody Seal: (Circle) INTACT BROKEN **ABSENT**

Relinquished By: (3) _____ Date: _____ Time: _____ Received By: _____

Requested Turnaround Time and Special Instructions: **Standard**

Relinquished By: (4) _____ Date: _____ Time: _____ Received By: _____



Mike D. Mason
Richard Catlin & Associates
220 Old Dairy Rd.
Wilmington, NC 28405

Report Number: G128-2277

Client Project: Site BB-9 Groundwater Sampling

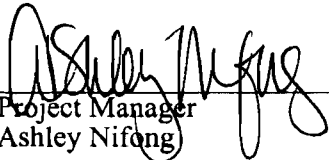
Dear Mike D. Mason,

Enclosed are the results of the analytical services performed under the referenced project. The samples are certified to meet the requirements of the National Environmental Laboratory Accreditation Conference Standards. Copies of this report and supporting data will be retained in our files for a period of five years in the event they are required for future reference. Any samples submitted to our laboratory will be retained for a maximum of thirty (30) days from the date of this report unless other arrangements are requested.

If there are any questions about the report or the services performed during this project, please call SGS Environmental Services at (910) 350-1903. We will be happy to answer any questions or concerns which you may have.

Thank you for using SGS Environmental Services for your analytical services. We look forward to working with you again on any additional analytical needs which you may have.

Sincerely,
SGS Environmental Services, Inc.

 12/1/08
Project Manager Date
Ashley Nifong

Case Narrative
SGS Project: **G128-2277**
Project Name: **Site BB-9 GW Sampling**

For Method: EPH

- The submitted sample was accepted into the lab on November 18th, 2008 and analyzed on November 19th and 23rd, 2008 by method EPH.
- The LCS/LCSD associated with batch 12995 had reported recoveries for the C₉-C₁₈ aliphatic that are outside of the method's QC limits. The sample was reextracted in batch 13033. The LCS/LCSD from this batch also has reported recoveries for the C₉-C₁₈ aliphatic that are outside of the method's QC limits. Both sets of data are reported in this SDG.



Craig R. Tronzo
Data Validation

List of Reporting Abbreviations
And Data Qualifiers

B = Compound also detected in batch blank

BQL = Below Quantification Limit (RL or MDL)

DF = Dilution Factor

Dup = Duplicate

P/D = Detected, but RPD is > 25/40% between results in dual column method.

E = Estimated concentration, exceeds calibration range.

J = Estimated concentration, below calibration range and above MDL

LCS(D) = Laboratory Control Spike (Duplicate)

MDL = Method Detection Limit

MS(D) = Matrix Spike (Duplicate)

PQL = Practical Quantitation Limit

RL/CL = Reporting Limit / Control Limit

RPD = Relative Percent Difference

mg/kg = milligram per kilogram, ppm, parts per million

ug/kg = micrograms per kilogram, ppb, parts per billion

mg/L = milligram per liter, ppm, parts per million

ug/L = micrograms per liter, ppb, parts per billion

% Rec = Percent Recovery

% solids = Percent Solids

Special Notes:

- 1) Metals and mercury samples are digested with a hot block, see the standard operating procedure document for details.
- 2) Uncertainty for all reported data is less than or equal to 30 percent.

**Results for Semivolatiles
by GCMS 625**

Client Sample ID: USTBB9-MW03R
 Client Project ID: Site BB-9 Groundwater Sampling
 Lab Sample ID: G128-2277-11
 Lab Project ID: G128-2277

Analyzed By: DES
 Date Collected: 11/17/2008 11:00
 Date Received: 11/18/2008
 Date Extracted: 11/21/2008
 Matrix: Water

Initial/Final Amt: 905 mL / 5.0 mL

Compound	Result ug/L	RL ug/L	MDL ug/L	Dilution Factor	Date Analyzed	Flag
Acenaphthene	38.6	5.52	0.823	1	11/26/2008	
Acenaphthylene	BQL	5.52	0.823	1	11/26/2008	
Anthracene	21.2	5.52	0.967	1	11/26/2008	
Benzo[a]anthracene	23.4	5.52	0.751	1	11/26/2008	
Benzo[a]pyrene	14.4	5.52	0.702	1	11/26/2008	
Benzo[b]fluoranthene	20.6	5.52	0.790	1	11/26/2008	
Benzo[g,h,i]perylene	9.23	5.52	0.680	1	11/26/2008	
Benzo[k]fluoranthene	9.50	5.52	0.608	1	11/26/2008	
Bis(2-chloroethoxy)methane	BQL	5.52	1.14	1	11/26/2008	
Bis(2-chloroethyl)ether	BQL	5.52	1.15	1	11/26/2008	
Bis(2-chloroisopropyl)ether	BQL	5.52	1.08	1	11/26/2008	
Bis(2-ethylhexyl)phthalate	BQL	5.52	0.453	1	11/26/2008	
4-bromophenyl phenyl ether	BQL	5.52	0.862	1	11/26/2008	
Butylbenzylphthalate	BQL	5.52	0.492	1	11/26/2008	
2-Chloronaphthalene	BQL	5.52	0.956	1	11/26/2008	
2-Chlorophenol	BQL	5.52	1.29	1	11/26/2008	
4-Chloro-3-methylphenol	BQL	5.52	0.878	1	11/26/2008	
4-Chlorophenyl phenyl ether	BQL	5.52	3.60	1	11/26/2008	
Chrysene	24.0	5.52	0.613	1	11/26/2008	
Dibenzo[a,h]anthracene	BQL	5.52	0.486	1	11/26/2008	
Di-n-Butylphthalate	BQL	5.52	0.912	1	11/26/2008	
3,3'-Dichlorobenzidine	BQL	11.0	1.35	1	11/26/2008	
2,4-Dichlorophenol	BQL	5.52	1.24	1	11/26/2008	
Diethylphthalate	BQL	5.52	0.818	1	11/26/2008	
Dimethylphthalate	BQL	5.52	0.613	1	11/26/2008	
2,4-Dimethylphenol	BQL	5.52	1.79	1	11/26/2008	
Di-n-octylphthalate	BQL	5.52	0.641	1	11/26/2008	
4,6-Dinitro-2-methylphenol	BQL	27.6	0.608	1	11/26/2008	
2,4-Dinitrophenol	BQL	27.6	0.707	1	11/26/2008	
2,4-Dinitrotoluene	BQL	5.52	0.591	1	11/26/2008	
2,6-Dinitrotoluene	BQL	5.52	0.718	1	11/26/2008	
Diphenylamine *	BQL	5.52	0.630	1	11/26/2008	
Fluoranthene	75.2	5.52	0.779	1	11/26/2008	
Fluorene	12.3	5.52	0.801	1	11/26/2008	
Hexachlorobenzene	BQL	5.52	0.558	1	11/26/2008	
Hexachlorobutadiene	BQL	5.52	0.840	1	11/26/2008	
Hexachlorocyclopentadiene	BQL	11.0	11.0	1	11/26/2008	
Hexachloroethane	BQL	5.52	0.823	1	11/26/2008	
Indeno(1,2,3-c,d)pyrene	BQL	5.52	2.52	1	11/26/2008	
Isophorone	BQL	5.52	0.978	1	11/26/2008	
Naphthalene	BQL	5.52	1.01	1	11/26/2008	
Nitrobenzene	BQL	5.52	1.16	1	11/26/2008	
2-Nitrophenol	BQL	5.52	1.36	1	11/26/2008	
4-Nitrophenol	BQL	27.6	1.19	1	11/26/2008	
N-Nitrosodi-n-propylamine	BQL	5.52	1.66	1	11/26/2008	
Pentachlorophenol	BQL	27.6	1.56	1	11/26/2008	
Phenanthrene	12.1	5.52	0.492	1	11/26/2008	

**Results for Semivolatiles
by GCMS 625**

Client Sample ID: USTBB9-MW03R
 Client Project ID: Site BB-9 Groundwater Sampling
 Lab Sample ID: G128-2277-11
 Lab Project ID: G128-2277

Analyzed By: DES
 Date Collected: 11/17/2008 11:00
 Date Received: 11/18/2008
 Date Extracted: 11/21/2008
 Matrix: Water


Initial/Final Amt: 905 mL / 5.0 mL

Compound	Result ug/L	RL ug/L	MDL ug/L	Dilution Factor	Date Analyzed	Flag
Phenol	BQL	5.52	1.17	1	11/26/2008	
Pyrene	64.7	5.52	2.28	1	11/26/2008	
1,2,4-Trichlorobenzene	BQL	5.52	0.796	1	11/26/2008	
2,4,6-Trichlorophenol	BQL	5.52	1.02	1	11/26/2008	
		Spike Added	Spike Result	Percent Recovered		
2-Fluorobiphenyl		10	6.3	63		
2-Fluorophenol		10	8.2	82		
Nitrobenzene-d5		10	9.8	98		
Phenol-d6		10	9	90		
2,4,6-Tribromophenol		10	7.5	75		
4-Terphenyl-d14		10	3	30		

Comments:

Flags:

BQL = Below Quantitation Limits.
 J = Detected below the quantitation limit.

Reviewed By:  _____

Results of Library Search for Semivolatile Compounds by GCMS

Client Sample ID: USTBB9-MW03R
 Client Project ID: Site BB-9 Groundwater Sampling
 Lab Sample ID: G128-2277-1I
 Lab Project ID: G128-2277
 Sample Wt/Vol: 905 ML
 Dilution: 1

Analyzed By: DES
 Date Collected: 11/17/2008 11:00
 Date Received: 11/18/2008
 Date Extracted: 11/21/2008
 Date Analyzed: 11/26/2008
 Matrix: Water

No.	Compound	Retention Time	CAS#	Match Probability	Result (ug/L)
1	Unknown	9.56			23.6
2	Unknown	5.24			17.3
3	Naphthalene, 2,6-dimethyl-	6.89	000581-42-0	96	12.8
4	9H-Fluoren-9-ol	8.68	001689-64-1	91	12.6
5	Naphthalene, 2,6-dimethyl-	6.96	000581-42-0	98	11.4
6	Unknown	8.29			8.97
7	Naphthalene, 1,6-dimethyl-	6.99	000575-43-9	96	7.81
8	Unknown	8.81			7.71
9	Pentadecane, 2,6,10,14-tetramethyl-	8.32	001921-70-6	72	7.44
10					

Comment:

Tentatively Identified Compound (TIC) refers to substances which are not present in the list of target compounds. Therefore, not all TICs are identified and quantitated using individual standards. TIC listings are prepared utilizing a computerized library search of electron impact mass spectral data and evaluation of the relevant data by a mass spectral data specialist.

Quantitation is accomplished by relative peak area of the compound compared to that of the nearest internal standard from the total ion chromatogram. TICs are identified and quantitated only if the peak area is equal to or greater than 10% of that of the nearest internal standard. Quantitation provided is an estimate.

Reviewed by: _____



EPH (Aliphatics/Aromatics) Laboratory Reporting Form

Client Name: Richard Catlin & Associates

Project Name: Site BB-9 Groundwater Sampling

Sample Information	
Sample Identification	USTBB9-MW03R
Sample Matrix	WATER
Date Collected	11/17/08
Date Received	11/18/08
Date Extracted	11/19/08
Date Analyzed	11/21/08 00:26 - 11/21/08 00:54
Dry Weight	NA
Dilution Factor	1 - 1
Initial Volume (mL)	903
Final Volume (mL)	5.0

Analytical Results			
Analytes**	Result µg/L	Report Limit µg/L	Flags
C9-C18 Aliphatics	833	100	
C19-C36 Aliphatics	2870	100	
C11-C22 Aromatics	4510	100	

Surrogates	Percent Recovery	Flags	Limits	
			Lower	Upper
Aliphatic (tricosane)	57.1		40	140
Aromatic (ortho-terphenyl)	56.2		40	140
Fractionation 1 (2-bromonaphthalene)	115		40	140
Fractionation 2 (2-fluorobiphenyl)	103		40	140

** = Excludes any surrogates or internal standards and are unadjusted for individual analytes.

Lab Info: G128-2277-1H	Lab Info: G128-2277-1H
Aliphatic: EP112008/021F1901.D	Aromatic: EP112008/022F2001.D

Reviewed By: 

EPH (Aliphatics/Aromatics) Laboratory Reporting Form

Client Name: Richard Catlin & Associates

Project Name: Site BB-9 Groundwater Sampling

Sample Information	
Sample Identification	USTBB9-MW03R
Sample Matrix	WATER
Date Collected	11/17/08
Date Received	11/18/08
Date Extracted	11/23/08
Date Analyzed	11/27/08 00:04 - 11/27/08 00:32
Dry Weight	NA
Dilution Factor	1 - 1
Initial Volume (mL)	965
Final Volume (mL)	5.0

Analytical Results			
Analytes**	Result µg/L	Report Limit µg/L	Flags
C9-C18 Aliphatics	1050	100	
C19-C36 Aliphatics	3470	100	
C11-C22 Aromatics	5260	100	

Surrogates	Percent Recovery	Flags	Limits	
			Lower	Upper
Aliphatic (tricosane)	45.7		40	140
Aromatic (ortho-terphenyl)	49.8		40	140
Fractionation 1 (2-bromonaphthalene)	119		40	140
Fractionation 2 (2-fluorobiphenyl)	106		40	140

** = Excludes any surrogates or internal standards and are unadjusted for individual analytes.

Lab Info: G128-2277-1K	Lab Info: G128-2277-1K
Aliphatic: EP112608/028F2001.D	Aromatic: EP112608/029F2101.D

Reviewed By: *MLC*

Attachment 3

EPH Laboratory Reporting Form

Calibration and QA/QC Information

Initial Calibration Date: 10/22/08

Calibration Ranges and Limits

Range	MDL		ML		RL	
	(02/15/08) (µg/L)	(02/11/08) (mg/Kg)	(µg/L)	(mg/Kg)	(µg/L)	(mg/Kg)
C9-C18 Aliphatics	1.66	0.274	5.28	0.871	100	10
C19-C36 Aliphatics	2.79	0.201	8.87	0.639	100	10
C11-C22 Aromatics	2.64	0.110	8.40	0.350	100	10

Calibration Concentration Levels

Range	Levels (µg/L)	Levels (mg/Kg)	%RSD if CF r if LR	Method of Quantitation
C ₉ -C ₁₈ Aliphatics	200	33.3	17.73	Calibration Factor
	100	16.7		
	50	8.33		
	25	4.17		
	5	0.833		
C ₁₉ -C ₃₆ Aliphatics	200	33.3	10.65	Calibration Factor
	100	16.7		
	50	8.33		
	25	4.17		
	5	0.833		
C ₁₁ -C ₂₂ Aromatics	200	33.3	10.39	Calibration Factor
	100	16.7		
	50	8.33		
	25	4.17		
	5	0.833		

Calibration Check Date: 11/20/08
11/20/08

Filenames: ep112008/001f0101.d
ep112008/002f0201.d

Calibration Check

Range	Levels (µg/L)	Levels (mg/Kg)	%Difference if CF %Drift if LR	Limits
C9-C18 Aliphatics	100	16.7	13.3	±25%
C19-C36 Aliphatics	100	16.7	8.2	±25%
C11-C22 Aromatics	100	16.7	3.1	±25%

MDL = Method Detection Limit
ML = Minimum Limit
RL = Reportable Limit

RPD = Relative Percent Difference
%RSD = Percent Relative Standard Deviation
CCC = Correlation Coefficient of Curve

Attachment 3

EPH Laboratory Reporting Form

Calibration and QA/QC Information

Initial Calibration Date: 10/22/08

Calibration Ranges and Limits

Range	MDL		ML		RL	
	(02/15/08) (µg/L)	(02/11/08) (mg/Kg)	(µg/L)	(mg/Kg)	(µg/L)	(mg/Kg)
C9-C18 Aliphatics	1.66	0.274	5.28	0.871	100	10
C19-C36 Aliphatics	2.79	0.201	8.87	0.639	100	10
C11-C22 Aromatics	2.64	0.110	8.40	0.350	100	10

Calibration Concentration Levels

Range	Levels (µg/L)	Levels (mg/Kg)	%RSD if CF r if LR	Method of Quantitation
C ₉ -C ₁₈ Aliphatics	200	33.3	17.73	Calibration Factor
	100	16.7		
	50	8.33		
	25	4.17		
	5	0.833		
C ₁₉ -C ₃₆ Aliphatics	200	33.3	10.65	Calibration Factor
	100	16.7		
	50	8.33		
	25	4.17		
	5	0.833		
C ₁₁ -C ₂₂ Aromatics	200	33.3	10.39	Calibration Factor
	100	16.7		
	50	8.33		
	25	4.17		
	5	0.833		

Calibration Check Date: 11/20/08
11/21/08

Filenames: ep112008/029f2701.d
ep112008/030f2801.d

Calibration Check

Range	Levels (µg/L)	Levels (mg/Kg)	%Difference if CF %Drift if LR	Limits
C9-C18 Aliphatics	100	16.7	18.6	≤±25%
C19-C36 Aliphatics	100	16.7	6.3	≤±25%
C11-C22 Aromatics	100	16.7	-24.0	≤±25%

MDL = Method Detection Limit
ML = Minimum Limit
RL = Reportable Limit

RPD = Relative Percent Difference
%RSD = Percent Relative Standard Deviation
CCC = Correlation Coefficient of Curve

Attachment 3

EPH Laboratory Reporting Form

Calibration and QA/QC Information

Initial Calibration Date: 10/22/08

Calibration Ranges and Limits

Range	MDL		ML		RL	
	(02/15/08) (µg/L)	(02/11/08) (mg/Kg)	(µg/L)	(mg/Kg)	(µg/L)	(mg/Kg)
C9-C18 Aliphatics	1.66	0.274	5.28	0.871	100	10
C19-C36 Aliphatics	2.79	0.201	8.87	0.639	100	10
C11-C22 Aromatics	2.64	0.110	8.40	0.350	100	10

Calibration Concentration Levels

Range	Levels (µg/L)	Levels (mg/Kg)	%RSD if CF r if LR	Method of Quantitation
C ₉ -C ₁₈ Aliphatics	200	33.3	17.73	Calibration Factor
	100	16.7		
	50	8.33		
	25	4.17		
	5	0.833		
C ₁₉ -C ₃₆ Aliphatics	200	33.3	10.65	Calibration Factor
	100	16.7		
	50	8.33		
	25	4.17		
	5	0.833		
C ₁₁ -C ₂₂ Aromatics	200	33.3	10.39	Calibration Factor
	100	16.7		
	50	8.33		
	25	4.17		
	5	0.833		

Calibration Check Date: 11/26/08
11/26/08

FileNames: ep112608/022f1401.d
ep112608/023f1501.d

Calibration Check

Range	Levels (µg/L)	Levels (mg/Kg)	%Difference if CF %Drift if LR	Limits
C9-C18 Aliphatics	100	16.7	16.1	≤±25%
C19-C36 Aliphatics	100	16.7	11.4	≤±25%
C11-C22 Aromatics	100	16.7	4.1	≤±25%

MDL = Method Detection Limit
ML = Minimum Limit
RL = Reportable Limit

RPD = Relative Percent Difference
%RSD = Percent Relative Standard Deviation
CCC = Correlation Coefficient of Curve

Attachment 3

EPH Laboratory Reporting Form

Calibration and QA/QC Information

Initial Calibration Date: 10/22/08

Calibration Ranges and Limits

Range	MDL		ML		RL	
	(02/15/08) (µg/L)	(02/11/08) (mg/Kg)	(µg/L)	(mg/Kg)	(µg/L)	(mg/Kg)
C9-C18 Aliphatics	1.66	0.274	5.28	0.871	100	10
C19-C36 Aliphatics	2.79	0.201	8.87	0.639	100	10
C11-C22 Aromatics	2.64	0.110	8.40	0.350	100	10

Calibration Concentration Levels

Range	Levels (µg/L)	Levels (mg/Kg)	%RSD if CF r if LR	Method of Quantitation
C ₉ -C ₁₈ Aliphatics	200	33.3	17.73	Calibration Factor
	100	16.7		
	50	8.33		
	25	4.17		
	5	0.833		
C ₁₉ -C ₃₆ Aliphatics	200	33.3	10.65	Calibration Factor
	100	16.7		
	50	8.33		
	25	4.17		
	5	0.833		
C ₁₁ -C ₂₂ Aromatics	200	33.3	10.39	Calibration Factor
	100	16.7		
	50	8.33		
	25	4.17		
	5	0.833		

Calibration Check Date: 11/26/08
11/27/08

Filenames: ep112608/034f2601.d
ep112608/035f2701.d

Calibration Check

Range	Levels (µg/L)	Levels (mg/Kg)	%Difference if CF %Drift if LR	Limits
C9-C18 Aliphatics	100	16.7	20.8	≤±25%
C19-C36 Aliphatics	100	16.7	16.6	≤±25%
C11-C22 Aromatics	100	16.7	11.3	≤±25%

MDL = Method Detection Limit
ML = Minimum Limit
RL = Reportable Limit

RPD = Relative Percent Difference
%RSD = Percent Relative Standard Deviation
CCC = Correlation Coefficient of Curve

VPH (Aliphatics/Aromatics) Laboratory Reporting Form

Client Name: Richard Catlin & Associates

Project Name: Site BB-9 Groundwater Sampling

Sample Information	
Sample Identification	USTBB9-MW03R
Sample Matrix	Water
Collection Option (for Soil)*	NA
Date Collected	11/17/08
Date Received	11/18/08
Date Extracted	11/25/08 23:11 - 11/25/08 23:11
Date Analyzed	11/25/08 23:11 - 11/25/08 23:11
Dry Weight	NA
Dilution Factor	1 - 1

Analytical Results				
Analyte	Result µg/L	Report Limit µg/L	Flags	
C ₅ -C ₈ Aliphatics**	BQL	100		
C ₉ -C ₁₂ Aliphatics**	BQL	100		
C ₉ -C ₁₀ Aromatics**	BQL	100		
	Percent Recovery	Flags	Limits Lower Upper	
Surrogate % Recovery - PID	90.0		70	130
Surrogate % Recovery - FID	95.2		70	130

* = Option 1 = Established fill line on vial, Option 2 = Sampling Device/Brand, or Option 3 = Field weight of soil.

** = Excludes any surrogates or internal standards and are unadjusted for individual analytes.

Lab Info: g128-2277-1a	Lab Info: g128-2277-1a
FID Info: VP112508/034F0101.D	PID Info: VP112508/034R0101.D

Reviewed By: 

Attachment 2

VPH Laboratory Reporting Form

Calibration and QA/QC Information

FID Initial Calibration Date: 11/18/08 PID Initial Calibration Date: 11/18/08

Calibration Ranges and Limits

Range	MDL		ML		RL	
	(µg/L)	(mg/Kg)	(µg/L)	(mg/Kg)	(µg/L)	(mg/Kg)
C ₅ -C ₈ Aliphatics	2.02	0.175	6.42	0.557	100	10
C ₉ -C ₁₂ Aliphatics	1.51	0.118	4.80	0.375	100	10
C ₉ -C ₁₀ Aromatics	0.902	0.132	2.87	0.420	100	10

Calibration Concentration Levels

Range	Levels (µg/L)	Levels (mg/Kg)	%RSD if CF r if LR	Method of Quantitation
C ₅ -C ₈ Aliphatics	10	0.8	4.30	Calibration Factor
	50	4		
	100	8		
	200	16		
	500	40		
C ₉ -C ₁₂ Aliphatics	10	0.8	1.00	Linear Regression
	50	4		
	100	8		
	200	16		
	500	40		
C ₉ -C ₁₀ Aromatics	10	0.8	14.21	Calibration Factor
	50	4		
	100	8		
	200	16		
	500	40		

Calibration Check Date: 11/25/08 Filename: VP112508/002F0101.d

Calibration Check

Range	Levels (µg/L)	Levels (mg/Kg)	%Difference if CF %Drift if LR	Limits
C ₅ -C ₈ Aliphatics	200	16	10.7 ✓	±25%
C ₉ -C ₁₂ Aliphatics	200	16	3.1 ✓	±25%
C ₉ -C ₁₀ Aromatics	200	16	4.8 ✓	±25%

MDL = Method Detection Limit
ML = Minimum Limit
RL = Reportable Limit

RPD = Relative Percent Difference
%RSD = Percent Relative Standard Deviation
CCC = Correlation Coefficient of Curve

Attachment 2

VPH Laboratory Reporting Form

Calibration and QA/QC Information

FID Initial Calibration Date: 11/18/08 PID Initial Calibration Date: 11/18/08

Calibration Ranges and Limits

Range	MDL		ML		RL	
	(µg/L)	(mg/Kg)	(µg/L)	(mg/Kg)	(µg/L)	(mg/Kg)
C ₅ -C ₈ Aliphatics	2.02	0.175	6.42	0.557	100	10
C ₉ -C ₁₂ Aliphatics	1.51	0.118	4.80	0.375	100	10
C ₉ -C ₁₀ Aromatics	0.902	0.132	2.87	0.420	100	10

Calibration Concentration Levels

Range	Levels (µg/L)	Levels (mg/Kg)	%RSD if CF r if LR	Method of Quantitation
C ₅ -C ₈ Aliphatics	10	0.8	4.30	Calibration Factor
	50	4		
	100	8		
	200	16		
	500	40		
C ₉ -C ₁₂ Aliphatics	10	0.8	1.00	Linear Regression
	50	4		
	100	8		
	200	16		
	500	40		
C ₉ -C ₁₀ Aromatics	10	0.8	14.21	Calibration Factor
	50	4		
	100	8		
	200	16		
	500	40		

Calibration Check Date: 11/25/08 Filename: VP112508/011F0101.d

Calibration Check

Range	Levels (µg/L)	Levels (mg/Kg)	%Difference if CF %Drift if LR	Limits
C ₅ -C ₈ Aliphatics	200	16	6.6 ✓	±25%
C ₉ -C ₁₂ Aliphatics	200	16	-3.6 ✓	±25%
C ₉ -C ₁₀ Aromatics	200	16	3.9 ✓	±25%

MDL = Method Detection Limit
ML = Minimum Limit
RL = Reportable Limit

RPD = Relative Percent Difference
%RSD = Percent Relative Standard Deviation
CCC = Correlation Coefficient of Curve

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VPH Laboratory Reporting Form

Calibration and QA/QC Information

FID Initial Calibration Date: 11/18/08 PID Initial Calibration Date: 11/18/08

Calibration Ranges and Limits

Range	MDL		ML		RL	
	(µg/L)	(mg/Kg)	(µg/L)	(mg/Kg)	(µg/L)	(mg/Kg)
C ₅ -C ₈ Aliphatics	2.02	0.175	6.42	0.557	100	10
C ₉ -C ₁₂ Aliphatics	1.51	0.118	4.80	0.375	100	10
C ₉ -C ₁₀ Aromatics	0.902	0.132	2.87	0.420	100	10

Calibration Concentration Levels

Range	Levels (µg/L)	Levels (mg/Kg)	%RSD if CF r if LR	Method of Quantitation
C ₅ -C ₈ Aliphatics	10	0.8	4.30	Calibration Factor
	50	4		
	100	8		
	200	16		
	500	40		
C ₉ -C ₁₂ Aliphatics	10	0.8	1.00	Linear Regression
	50	4		
	100	8		
	200	16		
	500	40		
C ₉ -C ₁₀ Aromatics	10	0.8	14.21	Calibration Factor
	50	4		
	100	8		
	200	16		
	500	40		

Calibration Check Date: 11/25/08 Filename: VP112508/043F0101.d

Calibration Check

Range	Levels (µg/L)	Levels (mg/Kg)	%Difference if CF %Drift if LR	Limits
C ₅ -C ₈ Aliphatics	200	16	-3.8 ✓	±25%
C ₉ -C ₁₂ Aliphatics	200	16	-7.4 ✓	±25%
C ₉ -C ₁₀ Aromatics	200	16	1.3 ✓	±25%

MDL = Method Detection Limit
ML = Minimum Limit
RL = Reportable Limit

RPD = Relative Percent Difference
%RSD = Percent Relative Standard Deviation
CCC = Correlation Coefficient of Curve

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VPH Laboratory Reporting Form

Calibration and QA/QC Information

FID Initial Calibration Date: 11/18/08 PID Initial Calibration Date: 11/18/08

Calibration Ranges and Limits

Range	MDL		ML		RL	
	(µg/L)	(mg/Kg)	(µg/L)	(mg/Kg)	(µg/L)	(mg/Kg)
C ₅ -C ₈ Aliphatics	2.02	0.175	6.42	0.557	100	10
C ₉ -C ₁₂ Aliphatics	1.51	0.118	4.80	0.375	100	10
C ₉ -C ₁₀ Aromatics	0.902	0.132	2.87	0.420	100	10

Calibration Concentration Levels

Range	Levels (µg/L)	Levels (mg/Kg)	%RSD if CF r if LR	Method of Quantitation
C ₅ -C ₈ Aliphatics	10	0.8	4.30	Calibration Factor
	50	4		
	100	8		
	200	16		
	500	40		
C ₉ -C ₁₂ Aliphatics	10	0.8	1.00	Linear Regression
	50	4		
	100	8		
	200	16		
	500	40		
C ₉ -C ₁₀ Aromatics	10	0.8	14.21	Calibration Factor
	50	4		
	100	8		
	200	16		
	500	40		

Calibration Check Date: 11/26/08 Filename: VP112608/002F0101.d

Calibration Check

Range	Levels (µg/L)	Levels (mg/Kg)	%Difference if CF %Drift if LR	Limits
C ₅ -C ₈ Aliphatics	200	16	2.7 ✓	±25%
C ₉ -C ₁₂ Aliphatics	200	16	-7.1 /	±25%
C ₉ -C ₁₀ Aromatics	200	16	2.2 ✓	±25%

MDL = Method Detection Limit
ML = Minimum Limit
RL = Reportable Limit

RPD = Relative Percent Difference
%RSD = Percent Relative Standard Deviation
CCC = Correlation Coefficient of Curve

Attachment 2

VPH Laboratory Reporting Form

Calibration and QA/QC Information

FID Initial Calibration Date: 11/18/08 PID Initial Calibration Date: 11/18/08

Calibration Ranges and Limits

Range	MDL		ML		RL	
	(µg/L)	(mg/Kg)	(µg/L)	(mg/Kg)	(µg/L)	(mg/Kg)
C ₅ -C ₈ Aliphatics	2.02	0.175	6.42	0.557	100	10
C ₉ -C ₁₂ Aliphatics	1.51	0.118	4.80	0.375	100	10
C ₉ -C ₁₀ Aromatics	0.902	0.132	2.87	0.420	100	10

Calibration Concentration Levels

Range	Levels (µg/L)	Levels (mg/Kg)	%RSD if CF r if LR	Method of Quantitation
C ₅ -C ₈ Aliphatics	10	0.8	4.30	Calibration Factor
	50	4		
	100	8		
	200	16		
	500	40		
C ₉ -C ₁₂ Aliphatics	10	0.8	1.00	Linear Regression
	50	4		
	100	8		
	200	16		
	500	40		
C ₉ -C ₁₀ Aromatics	10	0.8	14.21	Calibration Factor
	50	4		
	100	8		
	200	16		
	500	40		

Calibration Check Date: 11/26/08 Filename: VP112608/043F0101.d

Calibration Check

Range	Levels (µg/L)	Levels (mg/Kg)	%Difference if CF %Drift if LR	Limits
C ₅ -C ₈ Aliphatics	200	16	-2.0 ✓	±25%
C ₉ -C ₁₂ Aliphatics	200	16	-11.6 ✓	±25%
C ₉ -C ₁₀ Aromatics	200	16	0.1 /	±25%

MDL = Method Detection Limit
ML = Minimum Limit
RL = Reportable Limit

RPD = Relative Percent Difference
%RSD = Percent Relative Standard Deviation
CCC = Correlation Coefficient of Curve

