

NOR-00973

February 2, 2011

Mr. Stephen Scharf New York Department of Environmental Conservation Division of Environmental Remediation Bureau of Remedial Action A 625 Broadway, 11th Floor Albany, New York 12233-7015

Reference: CLEAN Contract No. N62472-03-0057 Contract Task Order 66

Subject: BPOW 1-3 Outpost Monitoring Wells Repair and Sampling Summary NWIRP Bethpage, New York

Dear Mr. Scharf:

On behalf of the Navy, please find enclosed a copy of the subject document. This document provides a summary of activities to repair and sample outpost monitoring well BPOW 1-3 at Naval Weapons Industrial Reserve Plant (NWIRP) Bethpage, New York.

If you have any questions please contact Ms. Lora Fly, NAVFAC Mid-LANT, at (757) 341-2012.

Sincerely

David D. Brayack, P.È. Project Manager

Enclosure: (1) BPOW 1-3 Outpost Monitoring Well Repair and Sampling Summary NWIRP Bethpage, New York

Distribution: Mid-Lant, Lora Fly NYSDEC (Albany), Henry Wilkie (email) NYSDOH (Troy), Steve Karpinski (email) NAVAIR, Richard Smith USEPA, Carol Stein (email) NGC, Kent Smith (email) South Farmingdale Water District Tetra Tech NUS, Dave Brayack ECOR Solutions, Al Taormina Administrative Record Project File

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BPOW 1-3 OUTPOST MONTORING WELLS REPAIR AND SAMPLING SUMMARY NWIRP BETHPAGE, NEW YORK

INTRODUCTION

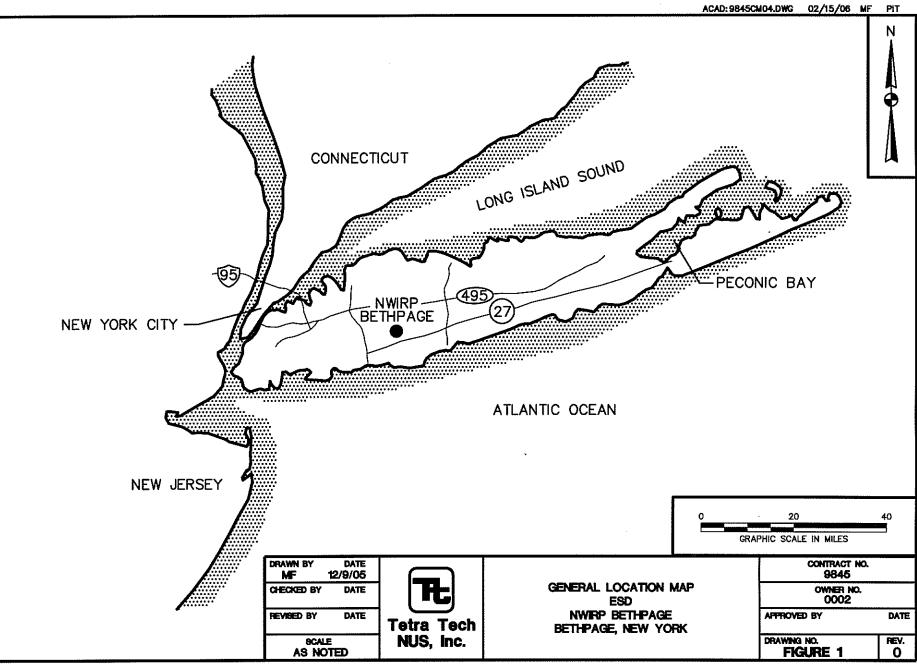
This document summarizes activities conducted to repair and sample Outpost Monitoring Well BPOW 1-3 at Naval Weapons Industrial Reserve Plant (NWIRP) Bethpage (Figure 1).

BACKGROUND

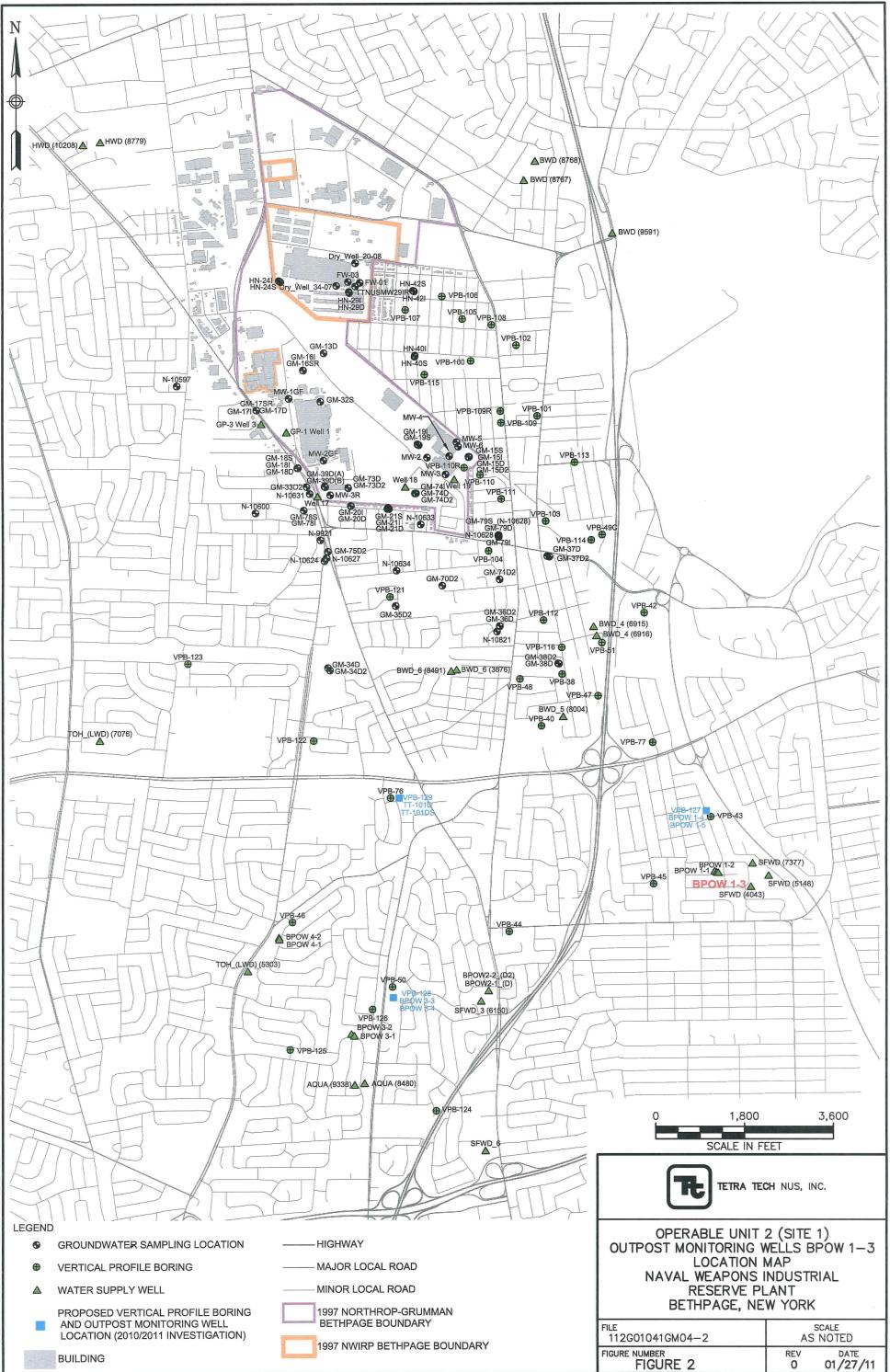
Groundwater samples collected during quarterly sampling events from outpost monitoring wells BPOW 1-1 (screened 200 to 235 feet bgs) and 1-3 (screened 375 to 400 feet bgs) consistently contained total VOCs of approximately 3 and 10 µg/L, respectively; however VOCs were not detected in groundwater samples collected from outpost monitoring well BPOW 1-2 (screened 310 to 340 feet bgs) that is screened between the BPOW 1-1 and 1-3 screen intervals. Based on a review of groundwater results and evaluation of the condition of BPOW 1-3, it was determined that the well casing may be cracked and has allowed shallower contaminated groundwater to infiltrate the well and be detected in BPOW 1-3. Therefore, a decision was made to seal the well by installing a 2-inch monitoring well inside the existing 4-inch casing.

OUTPOST MONITORING WELL BPOW 1-3

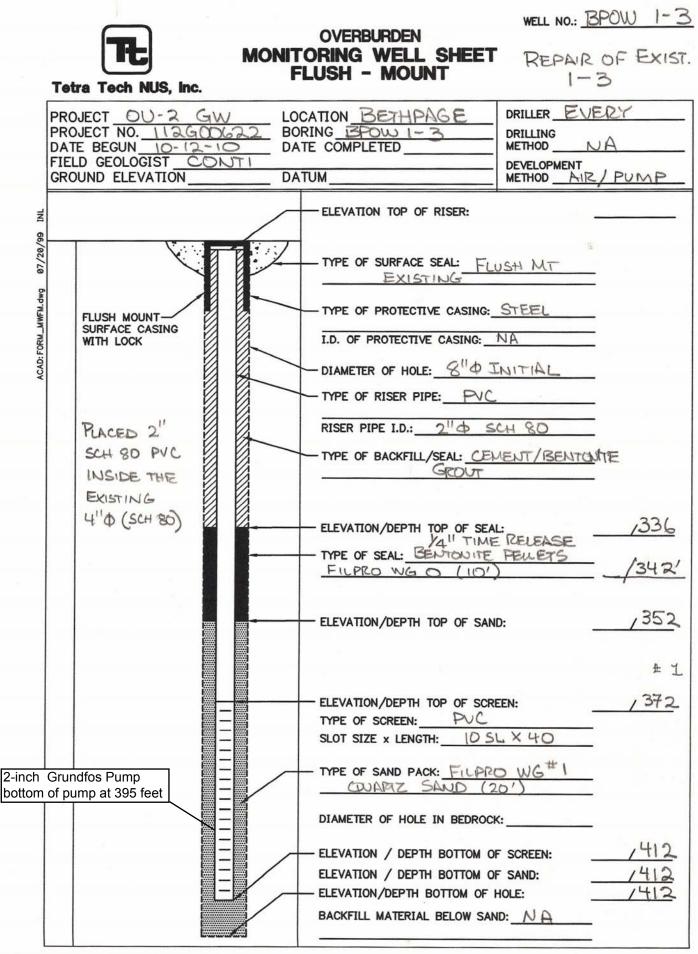
- BPOW 1-3 was repaired on October 12, 2010. The monitoring well repair construction log is provided in Attachment 1.
- BPOW 1-3 was re-developed on October 18, 2010. The monitoring well development log is provided in Attachment 2.
- A new dedicated 2-inch submersible pump (Grundfos[®]) was installed on December 7, 2010.
- Groundwater samples were collected from BPOW 2-1 on October 18, 2010 and December 8, 2010. The sample collected on October 18, 2010 was collected at the end of development to provide an initial evaluation of groundwater quality and are considered to be screening level quality. The sample collected on December 7, 2010 is representative of stabilized groundwater conditions and is considered a high level quality sample.
- Samples were analyzed for Target Compound List (TCL) VOCs. Sample logs sheets documenting the collection of these samples are provided in Attachment 3. Sample Chain of Custody forms are provided in Attachment 4.
- Site-related VOCs were not detected in groundwater samples collected from BPOW 1-3 during the October and December 2010 sampling events.
- Data validation reports are provided in Attachment 5.



FORM CADD NO. TTNUS-AH.DWG - REV 1 -9/10/98



ATTACHMENT 1 MONITORING WELL REPAIR CONSTRUCTION LOG



Pump installed on 12/8/2010.

ATTACHMENT 2 MONITORING WELL DEVELOPMENT LOG



Tetra Tech NUS, Inc.

MONITORING WELL DEVELOPMENT RECORD

Page ____ of ____

Dev. Metho	oped: <u>10 18</u> od: <u>PUMP</u> e: <u>2" GRUN</u>		Screen Length (ft. Specific Capacity: Casing ID (in.):			Project Number	112G0062: 7 GPM	(2"\$ 11	NSIDE OF EXIS 4" SCH 80. RED WELL.)
Time	Estimated Sediment Thickness (Ft.)	Cumulative Water Volume (Gal.)	Water Level Readings (Ft. below TOC)	Temperature (Degrees C)	pН	Specific Conductance (Units)	Turbidity (NTU)	R	emarks , color, etc.)
1050	1								
1100		70	31.60	14.81	4.21	. 234	42.8	SL. GLOUD	٢
1110	_	140	31.60	13.97	4.22	.123	2.9	CLEAR.	PH PAPER: 5
1120		210	31-60	13.50	4.15	-118	0.6	11	1
1130		280	31.60	13.25	4.13	.118	1.2		
1140	-	350	31.60	13.09	4.11	.116	0.4	ų	
1150	-	420	31.60	13.10	4.11	•115	1.1	U	
1210	-	560	31.60	13.17	4.11	-116	1.0		
1220		630	31.60	13.05	4.14	.121	0.8	11	★ 5→6
1240	-	770	31.60	13.23	4.13	-118	1.1	u	
1250		840	31.60	13.05	4.11	• 119	1.5	L C	
1300	-	910	31.60	-	/	/	/	11	
1310	-	980	31.60	13.2	4.13	-120	1.4	11	

ATTACHMENT 3 GROUNDWATER SAMPLE LOG

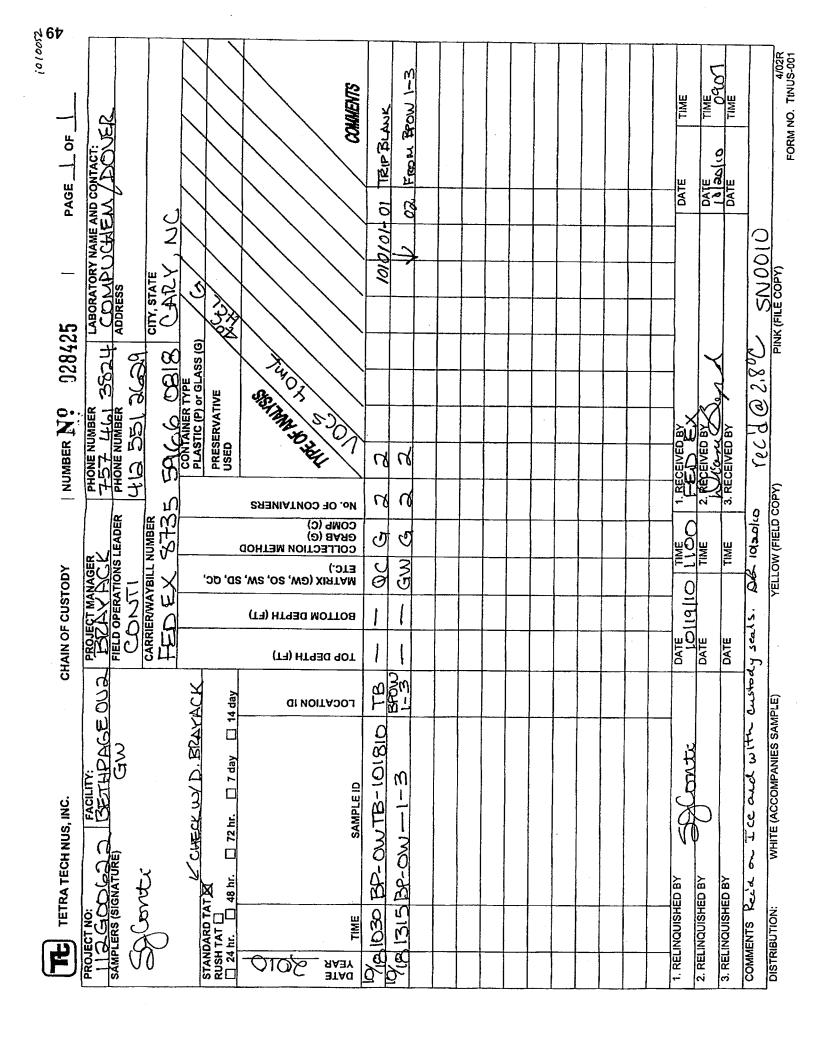
\square	
H	Tetra Tech NUS, Inc.

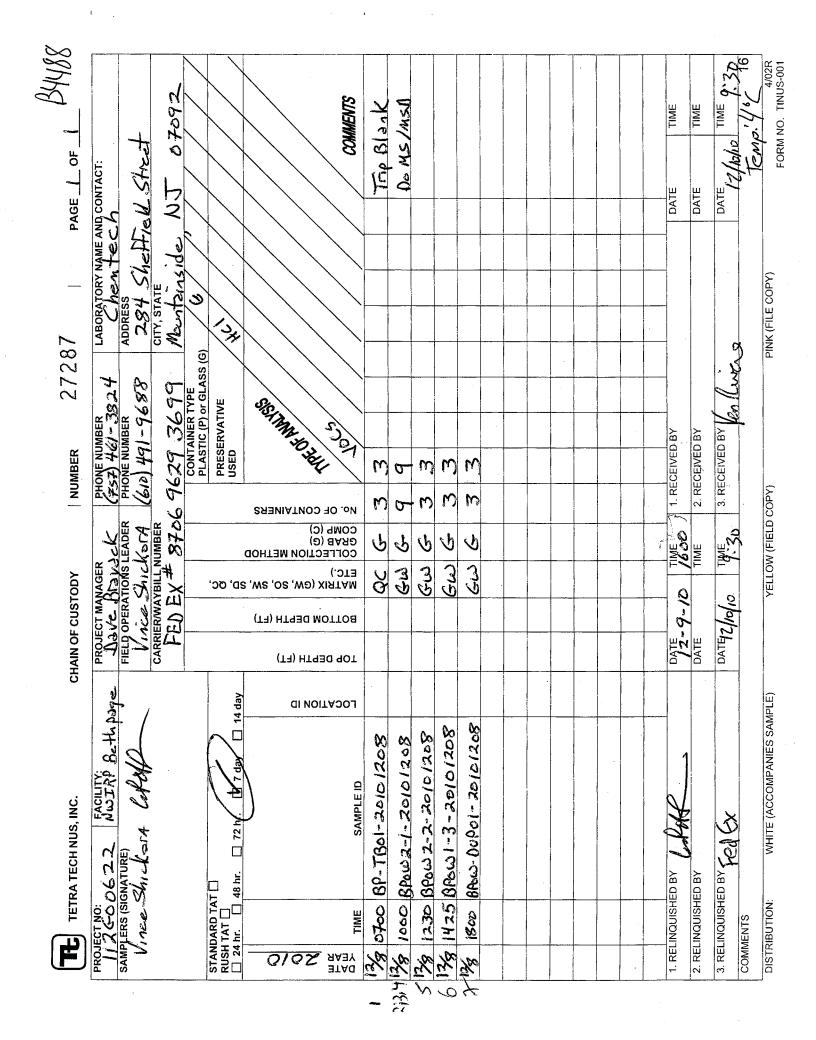
GROUNDWATER SAMPLE LOG SHEET

Project Site Name: Project No.:	BETHPA	22			Sample Sample		BPDW SJC	
[] Domestic Well Data [x] Monitoring Well Data [] Other Well Type: [] QA Sample Type:	a				[X] Lo	No.: Sample: w Concentr gh Concentr		15
SAMPLING DATA:								
Date: 10/18/10	Color	pН	S.C.	Temp.	Turbidity	DO	ORP	Other
Time: 1315	Visual	Standard 4.13	mS/cm	Degrees C 13.21	NTU	mg/l	mV	NA
Method: PUMP PURGE DATA:	CLEAR	14.13	-120	13.2	1.4	NA	NA	
Alighter App			alem Honeston, course	addidaa thataaay turi				
Date: 10/18/10		-	-					
Method: PUMP								-
Monitor Reading (ppm): O		-						-
Well Casing Diameter & Material	-					SHEET	-	
Type: 2" & PVC SCH 80			FOR	- PURG	EDAT	TA.	1	-
Total Well Depth (TD): ~ 412		1	_					
Static Water Level (WL): 30.05			(T)					
One Casing Volume(gal/L):								
Start Purge (hrs): 1050								
End Purge (hrs): 1310			1.					
Total Purge Time (min): 140				1.1.1.1				
Total Vol. Purged (gal/)_): 980								
SAMPLE COLLECTION INFORMA	TION: Strike	thru analy	sis not rec	uired			and the second	
Analysis		Preser	vative		Container R	equirements		Collected
VOCs		HCL/4	DEGC	(2-)	Oml Glass V	lals		V
			A	-				1000
		-						-
		-						-
		-						-
		-						
		1						
DBSERVATIONS / NOTES:								
2" MW = 0.163 gal/ft								
Sample taken at discreet intervals u	sing a hydrop	unch sampl	er unless o	therwise note	ed.			
			-					
Not enough volume for water quality Check box if not enough volume.	parameters	N	A		PURG	ING WA	S PAR	TOF
					DEV	SLOW	ED DI	QAAL
Jsed pH paper instead of water qua	lity meter	NI	A		Dou	N G 1	300 4	ND
Check box if used nH naner			<u>.</u>		Dou	PLED FI	ZOM 4	1
Check box if used pH paper.					SILIM	run r		
Check box if used pH paper.					SHOW	ICN CA.	IDIE S	TACE
Check box if used pH paper.					VAL	LED SAN	UPUE F	PORT
					VALI	UED SAN	UPUE F	2014
Check box if used pH paper. Circle if Applicable: MS/MSD Duplicate ID No.					VALI Signature(s	UED SAN	APUE F	2012

Tetra Tech NUS, Ir	ЪС.	GROU	JNDWA	TER SA	MPLE LO	DG SHEE	т		
9							Page	of	
Project Site Name: Project No.:	NWIR 	A Ber 5006	4page 22		_ Sample	Location:	<u> </u>	-20101208 21-3	2
[] Domestic Well Data ∭Monitoring Well Data [] Other Well Type: [] QA Sample Type:					Low				
SAMPLING DATA:									4
Date: 12-8-10	Color	рН	S.C.	Temp.	Turbidity	DO	Salinity	Other	-
Time: 1425	(Visual)	(S.U.)	(mS/cm)	(⁰ C)	(NTU)	(mg/l)	(%)	ORP	
Method: submersible Dump	clear	4.11	0.096	10.81	0.0	4.52	0.0	305	1
PURGE DATA:	(Gellons)								
Date: 12-8-10	Volume	рН	S.C.	Temp.	Turbidity	DO	Salinity	Other	Time
Method: submersible pump	1.0	4.14	0.083	11.11	1.6	5.61	0.0	303	1305
Monitor Reading (ppm): 0.0	45	4.11	0.095	10.89	0,0	4.77	0.0	316	1325
Well Casing Diameter & Material	90	4.11	0.096		0.0	4.55	0.0	309	1345
Type: Zinch PVC	135	4.12	0.096		0.0	4.48	0.0	304	1405
Total Well Depth (TD): 412	180	4.11	0,096		0.0	4.52		1	
Static Water Level (WL): 29.99	100	0.000	0.076	10.01		1.22	0.0	305	1425
One Casing Volume (gal/L): 62.3									-
									-
Start Purge (hrs): 1305		<u> </u>							-
End Purge (hrs): 1425			<u> </u>						_
Total Purge Time (min): 80									_
Total Vol. Purged (gal/L): 195 SAMPLE COLLECTION INFORMA			L		L				
	TION:	1 -							4
Analysis Vocs		Prese		384		equirements	****	Collected	-
<u> </u>		17	<u> </u>	384	fo mi v	rels		6	
									-
••••••••••••••••••••••••••••••••••••••									
									1
									-
									1
OBSERVATIONS / NOTES:	. Martin ad The State of States of States of		e antique an an an an an an an an						1
-punp set in we	ll at ~	395	BGS			water le	vel		
- Du o Mar a la	. 7 . 0		-			Sector Se	30 00	1	
- Sample split 10	ite AR	CHUE	5		Ì	3.30 ->	30,22	0	
- Sample split to - No stains, adars, or	elevate	d PID	readil	ngs obse					
Circle if Applicable:					Signature(s	$\frac{1+24}{2}$	30.23		-
MS/MSD Duplicate ID No.	***	an a			Signature(S	1 . 1	1		
- BPOW-		20101	208		6	1AF	1		

ATTACHMENT 4 SAMPLE CHAIN OF CUSTODY FORMS





ATTACHMENT 5 DATA VALIDATION PACKAGES



Tetra Tech NUS

INTERNAL CORRESPONDENCE

TO: D. BRAYACK

DATE: NOVEMBER 30, 2010

FROM: L. GANSER

COPIES: DV FILE

SUBJECT: ORGANIC DATA VALIDATION – VOC NWIRP BETHPAGE CTO 066 SAMPLE DELIVERY GROUP (SDG) 1010101

SAMPLES: 3/Aqueous/VOC

BP-OW-1-3 BP-OWTB-101810

<u>Overview</u>

The sample set for NWIRP Bethpage, CTO 066, SDG 1010101 consists of one (1) environmental aqueous sample and one (1) trip blank. The samples were analyzed for volatile organic compounds (VOCs). No field duplicates were included within this SDG.

The samples were collected on October 18, 2010 and analyzed by CompuChem, a division of Liberty Analytical Corporation. VOC analyses were conducted in accordance with EPA Method SW-846 8260B. The data contained in this SDG were validated with regard to the following parameters:

- * Data completeness
- * Holding times
- * GC/MŠ Tune
 - Initial/continuing calibrations
 - Laboratory Method Blank Results
- * Surrogate Recoveries
- Laboratory Control Sample/Laboratory Control Sample Duplicate Recoveries
- Internal Standard Recoveries
- * Compound Quantitation
- Compound Identification
- * Detection Limits

The symbol (*) indicates that all quality control criteria were met for this parameter. Qualified analytical results are presented in Appendix A, results as reported by the laboratory are presented in Appendix B, Region II data validation forms are presented in Appendix C, and documentation supporting these findings is presented in Appendix D.

Volatile Organic Compounds

Initial and continuing calibration relative response factor (RRF) was <0.05 for acetone. Positive results for acetone were qualified as estimated, "J".

Initial calibration percent relative standard deviation was >15% quality control limit (<90%) for bromomethane. Nondetected results for bromomethane were qualified as estimated, "UJ".

Contaminants were detected in laboratory method blank VBLKBO at the following maximum concentrations.

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Contaminant	Maximum Concentration (ug/L)	Action Level (ug/L)
1,2,4-Trichlorobenzene	1.6	8.0
Naphthalene	4.1	20.5
Toluene	0.92	4.6

An action level of 5X the maximum contaminant concentration was established to evaluate the samples for laboratory method blank contamination. Sample aliquot and dilution factors were taken into consideration during application of the blank action level. Positive results less than the action level were qualified as nondetected, "U", due to blank contamination. The trip blank was not qualified for laboratory blank contamination.

Additional Comments

Nondetected results are reported at the limit of detection (LOD).

Positive results below the limit of quantitation (LOQ) and above the detection limit were qualified as estimated, "J", due to uncertainty near the detection limit.

No matrix spike/matrix spike duplicate samples were requested.

The trip blank was reanalyzed by the laboratory to verify results. The results of the original analysis are presented in the EDD.

EXECUTIVE SUMMARY

Laboratory Performance Issues: 1,2,4-Trichlorobenzene, naphthalene, and toluene were detected in a laboratory method blank. Initial and continuing calibration RRF was <0.05 for acetone. Initial calibration percent relative standard deviation was greater than the quality control limit for bromomethane.

Other Factors Affecting Data Quality: None.

The data for these analyses were reviewed with reference to the EPA National Functional Guidelines for Organic Data Validation (10/99), USEPA Region II Standard Operating Procedures for Validating Volatile Organic Compounds by SW-846 Method 8260B HW-24 Revision 2 (August 2008) and the Department of Defense (DoD) document entitled "Quality Systems Manual (QSM) for Environmental Laboratories" (January 2006).

The text of this report has been formulated to address only those problem areas affecting data quality.

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Tetra Tech NUS Leanne Ganser Data Validator

TetraTech NUS Joseph A. Samchuck Data Validation Quality Assurance Officer

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Attachments:

- 1.
- Appendix A Qualified Analytical Results Appendix B Results as Reported by the Laboratory 2.

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- Appendix C Region II Data Validation Forms Appendix D Support Documentation З.
- 4.

Appendix A

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Qualified Analytical Results

Data Validation Qualifier Codes:

- A = Lab Blank Contamination
- B = Field Blank Contamination
- C = Calibration Noncompliance (e.g. % RSDs, %Ds, ICVs, CCVs, RRFs, etc.)
- C01 = GC/MS Tuning Noncompliance
- D = MS/MSD Recovery Noncompliance
- E = LCS/LCSD Recovery Noncompliance
- F = Lab Duplicate Imprecision
- G = Field Duplicate Imprecision
- H = Holding Time Exceedance
- I = ICP Serial Dilution Noncompliance
- J = GFAA PDS-GFAA MSA's r < 0.995 / ICP PDS Recovery Noncompliance
- K = ICP Interference includes ICS % R Noncompliance
- L = Instrument Calibration Range Exceedance
- M = Sample Preservation Noncompliance
- N = Internal Standard Noncompliance
- N01 = Internal Standard Recovery Noncompliance Dioxins
- N02 = Recovery Standard Noncompliance Dioxins
- N03 = Clean-up Standard Noncompliance Dioxins
- O = Poor Instrument Performance (e.g. base-line drifting)
- P = Uncertainty near detection limit (< 2 x IDL for inorganics and <CRQL for organics)
- Q = Other problems (can encompass a number of issues; e.g. chromatography, interferences, etc.)
- R = Surrogates Recovery Noncompliance
- S = Pesticide/PCB Resolution
- T = % Breakdown Noncompliance for DDT and Endrin
- U = % Difference between columns/detectors >25% for positive results determined via GC/HPLC
- V = Non-linear calibrations; correlation coefficient r < 0.995
- W = EMPC result
- X = Signal to noise response drop
- Y = Percent solids <30%
- Z = Uncertainty at 2 sigma deviation is greater than sample activity

					0.0	
	NSAIVIPLE	BP-OW-1-3		BP-OWIB-101810	1810	
				10-1010101		
FRACTION: OV	SAMP_DATE	10/18/2010		10/18/2010		
MEDIA: WATER	QC_TYPE	NM		TB		
	UNITS	NG/L		NG/L		
	PCT_SOLIDS	0.0		0.0		
	DUP_OF					
PARAMETER			QLCD	RESULT	VaL	QLCD
1,1,1-TRICHLOROETHANE		0.5 U		0.5	n	
1,1,2,2-TETRACHLOROETHANE	HANE	2 U		2	D	
1,1,2-TRICHLOROETHANE		2 U		2	2 U	
1,1,2-TRICHLOROTRIFLUOROETHANE	ROETHANE	2 U		2	D	
1,1-DICHLOROETHANE		0.5 U		0.5 U	5	
1,1-DICHLOROETHENE		2 N		2	5	
1,1-DICHLOROPROPENE		0.5 U		0.5	D	
1,2,4-TRICHLOROBENZEN	Е	0.5 U		4.1	-	٩
1,2-DIBROMO-3-CHLOROPROPANE	ROPANE	2 U		2	∍	
1,2-DIBROMOETHANE		0.5 U		0.5 U	∍	
1,2-DICHLOROBENZENE		0.5 U		0.5		
1,2-DICHLOROETHANE		0.5 U		0.5	 	
1,2-DICHLOROPROPANE		2 U		2		
1,3-DICHLOROBENZENE		0.5 U	-	0.5 U	D	
1,4-DICHLOROBENZENE		0.5 U .		0.5 U	Э	
2-BUTANONE		5 U		5		
2-HEXANONE		1.3 U		1.3 U	_	
4-METHYL-2-PENTANONE		1.3 U		1.3 U		
ACETONE		6.1 J	СР	34	ر	U
BENZENE		0.5 U		0.5 U	5	
BROMODICHLOROMETHANE	NE	0.5 U		0.5	۰.	
BROMOFORM		2 U		2	n	
BROMOMETHANE		2 UJ	U	2	2 UJ	U
CARBON DISULFIDE		0.5 U		0.5 U	D	
CARBON TETRACHLORIDE		0.5 U		0.5 U	N	
CHLOROBENZENE		0.5 U		0.5	D	
CHLORODIBROMOMETHANE	Ш	0.5 U		0.5	N	
CHLOROETHANE		2 U	•	9.3	,	
CHLOROFORM		0.5 U		0.5	5	
CHLOROMETHANE		0.5 U		0.5	5	
CIS-1,2-DICHLOROETHENE	Int	2 U		2	5	
CIS-1,3-DICHLOROPROPENE	NE	2 U		7	n	
CYCLOHEXANE		0.5 U		0.5 U	5	
DICHLORODIFLUOROMETHANE	HANE	0.5 U		0.5		
ETHYLBENZENE		0.5 U		0.5		
ISOPROPYLBENZENE		0.5 U		0.5	 	
1 of 2						

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PRUJ_NU: UU622	INDAINIPLE	BP-UW-1-3			BP-UW1B-101810	1810		
SDG: 1010101	LAB_ID	1010101-02			1010101-01			
FRACTION: OV	SAMP_DATE	10/18/2010			10/18/2010			1
MEDIA: WATER	QC_TYPE	NM			TB			r
	UNITS	NG/L			NG/L			
	PCT_SOLIDS	0.0			0.0			-
	DUP_OF							-
PARAMETER		RESULT	VQL	arcd	RESULT	VaL	QLCD	1
M+P-XYLENES		-	1 U			1 U		1
METHYL ACETATE		2	2 U		5	2 U		1
METHYL CYCLOHEXANE		0.5 U	Ъ		0.5 U	<u>р</u>		T
METHYL TERT-BUTYL ETHER	HER	0.5	5		0.5 U	5		1
METHYLENE CHLORIDE		0.5	_		0.5 U	D		1
NAPHTHALENE		2	⊃		3.1	٦ ۲	٩.	
O-XYLENE		0.5 U	D		0.5 U	5		1
STYRENE		0.5 U	∍		0.5 U	n		r
TETRACHLOROETHENE		0.5 U	D		0.5 U	<u> </u>		γ <u> </u>
TOLUENE		0.86 U	n	A	-	1 J	٩	r
TOTAL 1,2-DICHLOROETHENE	IENE	0.5	D		0.5 U	Л		r
TOTAL XYLENES		0.5	n		0.5 U	5		,
TRANS-1,2-DICHLOROETHENE	HENE	0.5 U	n		0.5 U	5		r
TRANS-1,3-DICHLOROPROPENE	OPENE	0.5 U	5		0.5 U	Л		r
TRICHLOROETHENE		0.5 U	Ъ		0.5 U	<u> </u>		η
TRICHLOROFLUOROMETHANE	HANE	0.5 U	N		0.5 U	U		
VINYL CHLORIDE		0.5 U	D		0.5	5		r1
	·							

12/1/2010



Tetra Tech NUS

D. BRAYACK

TO:

DATE: JANUARY 14, 2011

FROM: MICHELLE L. ALLEN COPIES: DV FILE

SUBJECT: ORGANIC DATA VALIDATION – VOC, SVOC, PEST, and PCB NAVAL WEAPONS INDUSTRIAL RESERVE PLANT (NWIRP), BETHPAGE CTO 066 SAMPLE DELIVERY GROUP (SDG) B4488

SAMPLES: 5/Aqueous/VOC

BP-TB01-20101208 BPOW2-1-20101208 BPOW-DUP01-20101208 BPOW2-2-20101208 BPOW1-3-20101208

1/IDW/VOC/SVOC/PEST/PCB

BP-FRACIDW-20101209

<u>Overview</u>

The sample sets for NWIRP Bethpage, SDG B4488 consisted of four (4) aqueous environmental samples, one (1) aqueous waste sample, and one (1) aqueous trip blank. All six (6) aqueous samples were analyzed for volatile organic compounds (VOC). The one (1) aqueous waste sample was analyzed for semi-volatile organic compounds (SVOC), pesticides (PEST), and polychlorinated biphenyls (PCB). One field duplicate sample pair was associated with this sample data group (SDG); BPOW-DUP01-20101208/BPOW1-3-20101208.

The samples were collected by Tetra Tech on December 8 and 9, 2010 and analyzed by Chemtech. All analyses were conducted in accordance with EPA Methods SW-846 8260B, 8270C, 8081, 8082 and EPA Method 624 analytical and reporting protocols. The data contained in this SDG was validated with regard to the following parameters:

- Data completeness
- Hold times
 - GC/MS System Tuning and Performance
 - Initial/continuing calibrations
 - Laboratory Method Blank Results
 - Surrogate Spike Recoveries
 - Internal Standard Recoveries
 - Laboratory Control Sample/Laboratory Control Sample Duplicate Recoveries
 - Matrix Spike/Matrix Spike Duplicate Results
 - Field Duplicate Precision Results
- Compound Identification
- Compound Quantitation
- Detection Limits

The symbol (*) indicates that all quality control criteria were met for this parameter. Qualified analytical results are presented in Appendix A, results as reported by the laboratory are presented in Appendix B, Region II data validation forms are presented in Appendix C, and documentation supporting these findings is presented in Appendix D.

TO: D. BRAYACK SDG: B4488

Volatile (VOC)

The Percent Differences (%Ds) for 2-hexanone and bromoform exceeded the 20% quality control limit for the continuing calibration performed on instrument MSVOAD on 12/15/10 @ 11:33. Sample BPOW-DUP01-20101208 was affected. Only non-detected results were reported for these compounds in the affected sample and these non-detects were qualified as estimated, (UJ).

The continuing calibration %Ds for acetone and methyl acetate were greater than 20% quality control criteria on instrument MSVOAG on 12/13/10 @ 10:41 affecting samples TB01-20101208, BPOW1-3-20101208, BPOW2-1-20101208, and BPOW2-2-20101208. The non-detected results reported for these compounds were qualified as estimated, (UJ).

The Relative Percent Difference (RPD) for acetone, methyl acetate, 2-butanone, 1,1,2,2tetrachloroethene, and 1,2-dibromo-3-chloropropane exceeded the 20% quality control limit in the Matrix Spike/Matrix Spike Duplicate (MS/MSD) analyses of sample BPOW2-1-20101208. No action was taken for the non-detected results reported for these compounds in the environmental sample since the Percent Recoveries (%Rs) were acceptable in the MS and MSD samples.

The Laboratory Control Sample (LCS), BSG1213W1, had %Rs for acetone and methyl acetate above the upper quality control limits. No action was taken in the affected samples since no positive results were reported for these compounds.

The LCS/Laboratory Control Sample Duplicate (LCSD) analyses, samples BSG1209W3/BSG1209W4, had RPDs for dichlorofluoromethane, chloromethane, vinyl chloride, bromomethane, chloroethane, trichlorofluoromethane, 1,1-dichloroethene, acetone, carbon disulfide, methyl acetate, and 2-butanone that exceeded 20%. In addition, the %R for acetone was greater that the upper quality control limit. No action was taken in the affected waste water sample since only non-detects were reported for the noncompliant compounds.

The positive result for 1,1-dichloroethane in sample BPOW2-2-20101208 reported below the Limit of Quantitation (LOQ) but above the Method Detection Limit (MDL) was qualified as estimated, (J). Non-detected results are reported to the Limit of Detection (LOD).

Semi-Volatile Organic Compounds (SVOC)

The internal standard, perylene-d12, was below the lower quality control limit in sample BP-FRACIDW-20101209. The sample was reanalyzed yielding similar results. The initial analysis of this sample was used in the data validation. The non-detected results reported the compounds associated with this internal standard were qualified as estimated, (UJ).

Pesticides (PEST)

No problems were noted.

Polychlorinated Biphenyls (PCB)

The surrogate spike compound, decachlorobiphenyl, had %Rs below the lower quality control limit in sample BP-FRACIDW-20101209 and its reanalysis. The initial analysis was used in the validation of the data. The non-detected results reported for the PCBs in this fraction were qualified as estimated, (UJ).

Additional Comments

The VOC analysis of the waste sample, BP-FRACIDW-20101209, was analyzed via EPA Method 624 and evaluated accordingly.

EXECUTIVE SUMMARY

Laboratory Performance Issues: Some compounds were estimated due to continuing calibration %Ds greater than their respective quality control limit. The VOC LCS/LSD had %Rs and RPDs outside the quality control limits. Noncompliant surrogate %Rs resulted in the qualification the waste sample in the PCB fraction. One internal standard was below the lower quality control limit in the SVOC analysis of the waste sample. Affected compounds were estimated.

Other Factors Affecting Data Quality: The MS/MSD sample had noncompliant %Rs and RPDs. Nondetected results were not qualified. A positive result reported below the LOQ but above the MDL was qualified as estimated, (J). Non-detected results are reported to the LOD.

The data for these analyses were reviewed with reference to the following: SOP #HW-24 Revision #2, August 2008, USEPA Region II Hazardous Waste Support Branch Validating Volatile Organic Compounds by Gas Chromatography/Mass Spectrometry, SOP #HW-22 Revision #4, August 2008, USEPA Region II Hazardous Waste Support Branch Validating Semivolatile Organic Compounds by Gas Chromatography/Mass Spectrometry, SOP #HW-44 Revision #1, October 2006, USEPA Region II Hazardous Waste Support Branch Validating Pesticides by Gas Chromatography, SOP #HW-45 Revision #1, October 2006, USEPA Region II Hazardous Waste Support Branch Validating Pesticides by Gas Chromatography, SOP #HW-45 Revision #1, October 2006, USEPA Region II Hazardous Waste Support Branch Validating Pesticides by Gas Chromatography, SOP #HW-45 Revision #1, October 2006, USEPA Region II Hazardous Waste Support Branch Validating Pesticides by Gas Chromatography, SOP #HW-45 Revision #1, October 2006, USEPA Region II Hazardous Waste Support Branch Validating Pesticides by Gas Chromatography, SOP #HW-45 Revision #1, October 2006, USEPA Region II Hazardous Waste Support Branch Validating Polychlorinated Biphenyls by Gas Chromatography by SW-846 Methods 8260B, 8270C, 8081, and 8082, EPA Method 624, and the Department of Defense (DoD) document entitled "Quality Systems Manual (QSM) for Environmental Laboratories" (January 2006).

VetraTech NUS Michelle L. Allen Chemist/Data Validator

TetraTech NUS Joseph A. Samchuck Data Validation Quality Assurance Officer

Attachments:

- 1. Appendix A Qualified Analytical Results
- 2. Appendix B Results as Reported by the Laboratory
- 3. Appendix C Region II Data Validation Forms
- 4. Appendix D Support Documentation

Appendix A

Qualified Analytical Results

Data Validation Qualifier Codes:

- A = Lab Blank Contamination
- B = Field Blank Contamination
- C = Calibration Noncompliance (e.g. % RSDs, %Ds, ICVs, CCVs, RRFs, etc.)
- C01 = GC/MS Tuning Noncompliance
- D = MS/MSD Recovery Noncompliance
- E = LCS/LCSD Recovery Noncompliance
- F = Lab Duplicate Imprecision
- G = Field Duplicate Imprecision
- H = Holding Time Exceedance
- I = ICP Serial Dilution Noncompliance
- J = GFAA PDS GFAA MSA's r < 0.995 / ICP PDS Recovery Noncompliance
- K = ICP Interference includes ICS % R Noncompliance
- L = Instrument Calibration Range Exceedance
- M = Sample Preservation Noncompliance
- N = Internal Standard Noncompliance
- N01 = Internal Standard Recovery Noncompliance Dioxins
- N02 = Recovery Standard Noncompliance Dioxins
- N03 = Clean-up Standard Noncompliance Dioxins
- O = Poor Instrument Performance (e.g. base-line drifting)
- P = Uncertainty near detection limit (< 2 x IDL for inorganics and <CRQL for organics)
- Q = Other problems (can encompass a number of issues; e.g. chromatography, interferences, etc.)
- R = Surrogates Recovery Noncompliance
- S = Pesticide/PCB Resolution
- T = % Breakdown Noncompliance for DDT and Endrin
- U = % Difference between columns/detectors >25% for positive results determined via GC/HPLC
- V = Non-linear calibrations; correlation coefficient r < 0.995
- W = EMPC result
- X = Signal to noise response drop
- Y = Percent solids <30%
- Z = Uncertainty at 2 sigma deviation is greater than sample activity

SDG: B4488 LAB_ID FRACTION: OV SAMP_DA MEDIA: WATER QC_TYPE MEDIA: WATER UNITS PCT_SOLI DUP_OF PARAMETER DUP_OF 1,1,2-TRICHLOROETHANE 1,1,2-TRICHLOROETHANE 1,1,2-TRICHLOROETHANE 1,1,2-TRICHLOROETHANE	LAB_ID B4488-06 SAMP_DATE 12/8/2010 QC_TYPE NM UNITS UGA1			B4488-02		B4488-05			B4488-07	
FRACTION: OV SAMP MEDIA: WATER QC_T MEDIA: WATER UNITS PCT_5 DUP_1 1,1,1-TRICHLOROETHANE 1,1,2,2-TETRACHLOROETHANE 1,1,2-TRICHLOROETHANE 1,1,2-TRICHLOROETHANE	DATE PE									
MEDIA: WATER QC_T UNITS PCT_G PCT_G PCT_G PCT_G PCT_G PCT_G PCT_G PCT_G PCT_G PCT_G PCT_G 1,1,2-TRICHLOROETHANE 1,1,2-TRICHLOROETHANE	ЪЕ			12/8/2010		12/8/2010			12/8/2010	
UNITS				MN		WN			MN	
PCT_G PARAMETER 1,1,1-TRICHLOROETHANE 1,1,2,2-TETRACHLOROETHANE 1,1,2,2-TRICHLOROETHANE				NG/L		NG/L			NG/L	
DUP_0 PARAMETER 1,1,1-TRICHLOROETHANE 1,1,2,2-TETRACHLOROETHANE 1,1,2-TRICHLOROETHANE	PCT_SOLIDS 0.0			0.0		0.0			0.0	
PARAMETER 1,1,1-TRICHLOROETHANE 1,1,2,2-TETRACHLOROETHANE 1,1,2-TRICHLOROETHANE			, .						BP0W1-3-20101208	8
1,1,1-TRICHLOROETHANE 1,1,2,2-TETRACHLOROETHANE 1,1,2-TRICHLOROETHANE	RESULT	Кa	QLCD	RESULT VQL	L QLCD	RESULT	VaL	alcd	RESULT VQL	L QLCD
1,1,2,2-TETRACHLOROETHANE 1,1,2-TRICHLOROETHANE	0	0.5 U		0.5 U		o	0.5 U		0.5 U	
1,1,2-TRICHLOROETHANE	0	0.5 U		0.5 U		o	0.5 U		0.5 U	
		0.5 U		0.5 U		0	0.5 U		0.5 U	
1,1,2-TRICHLOROTRIFLUOROETHANE		0.5 U		0.5 U		Ö	0.5 U		0.5 U	
1,1-DICHLOROETHANE	0	0.5 U		0.5 U		0.74	74 J	٩	0.5 U	
1,1-DICHLOROETHENE	0	0.5 U		0.5 U		0	0.5 U		0.5 U	
1,2,4-TRICHLOROBENZENE		0.5 U		0.5 U		0	0.5 U		0.5 U	
1,2-DIBROMO-3-CHLOROPROPANE		0.5 U		0.5 U		0	0.5 U		0.5 U	
1,2-DIBROMOETHANE	0	0.5 U		0.5 U		° O	0.5 U		0.5 U	
1,2-DICHLOROBENZENE	0	0.5 U		0.5 U		Ö	0.5 U		0.5 U	
1,2-DICHLOROETHANE	0	0.5 U		0.5 U		0	0.5 U		0.5 U	
1,2-DICHLOROPROPANE	0	0.5 U		0.5 U		ō	0.5 U		0.5 U	
1,3-DICHLOROBENZENE	0	0.5 U		0.5 U		O	0.5 U		0.5 U	
1,4-DICHLOROBENZENE	0	0.5 U		0.5 U		0	0.5 U		0.5 U	
2-BUTANONE	2	2.5 U		2.5 U		2	2.5 U		2.5 U	
2-HEXANONE	2	2.5 U		2.5 U		~	2.5 U		2.5 UJ	0
4-METHYL-2-PENTANONE	2	2.5 U		2.5 U		2.	2.5 U		2.5 U	
ACETONE	0	2.5 UJ	U	2.5 UJ	С	2	2.5 UJ	U	2.5 U	
BENZENE	0	0.5 U		0.5 U		O	0.5 U		0.5 U	
BROMODICHLOROMETHANE	0	0.5 U		0.5 U		0	0.5 U		0.5 U	
BROMOFORM	Ő	0.5 U		0.5 U		o 	0.5 U		0.5 UJ	U
BROMOMETHANE	Ö.	0.5 U		0.5 U		Ö	.5 U		0.5 U	
CARBON DISULFIDE	Ő	0.5 U		0.5 U		Ö	0.5 U		0.5 U	
CARBON TETRACHLORIDE	Ő	0.5 U		0.5 U		Ö	0.5 U		0.5 U	
CHLOROBENZENE	Ő	0.5 U		0.5 U		0	0.5 U		0.5 U	
CHLORODIBROMOMETHANE	O	0.5 U		0.5 U		0	0.5 U		0.5 U	
CHLOROETHANE	0	0.5 U		0.5 U		Ö	0.5 U		0.5 U	
CHLOROFORM	Ö	0.5 U		0.5 U		ö	0.5 U		0.5 U	
CHLOROMETHANE	Ő	0.5 U		0.5 U		.0	0.5 U		0.5 U	
CIS-1,2-DICHLOROETHENE	Ó	0.5 U		0.5 U		0	0.5 U		0.5 U	
CIS-1,3-DICHLOROPROPENE	0	0.5 U		0.5 U		o	0.5 U		0.5 U	
CYCLOHEXANE	0	0.5 U		0.5 U		Ö	0.5 U		0.5 U	
DICHLORODIFLUOROMETHANE	Ó	0.5 U		0.5 U		0	0.5 U		0.5 U	
ETHYLBENZENE	Ó	0.5 U		0.5 U		0	0.5 U		0.5 U	
ISOPROPYLBENZENE	Ö	0.5 U		0.5 U		0	0.5 U		0.5 U	
M+P-XYLENES		1 U		1			1 U		10	

PROJ_NO: 00622	NSAMPLE	BP-TB01-20101208	01208	
SDG: B4488	LAB_ID	B4488-01		
FRACTION: OV	SAMP_DATE	12/8/2010		
MEDIA: WATER	QC_TYPE	MΝ		
	UNITS	NG/L		
	PCT_SOLIDS	0.0		
	DUP_OF			
PARAMETER		RESULT	VQL	arcd
1,1,1-TRICHLOROETHANE		0.5	n	
1,1,2,2-TETRACHLOROETHANE	HANE	0.5	D	
1,1,2-TRICHLOROETHANE		0.5	<u>р</u>	
1,1,2-TRICHLOROTRIFLUOROETHANE	ROETHANE	0.5	D	
1,1-DICHLOROETHANE		0.5	D	
1,1-DICHLOROETHENE		0.5	D	
1,2,4-TRICHLOROBENZENE	Ш	0.5 U	D	
1,2-DIBROMO-3-CHLOROPROPANE	ROPANE	0.5 U	5	
1,2-DIBROMOETHANE		0.5 U	n	
1,2-DICHLOROBENZENE		0.5	n	
1,2-DICHLOROETHANE		0.5	D	
1,2-DICHLOROPROPANE		0.5	D	
1,3-DICHLOROBENZENE		0.5	D	
1,4-DICHLOROBENZENE		0.5	D	
2-BUTANONE		2.5	D	
2-HEXANONE		2.5	n	
4-METHYL-2-PENTANONE		2.5	n	
ACETONE		2.5	n	c
BENZENE		0.5	D	
BROMODICHLOROMETHANE	NE	0.5	D	
BROMOFORM		0.5	5	
BROMOMETHANE		0.5	n	
CARBON DISULFIDE		0.5	n	
CARBON TETRACHLORIDE	111	0.5	D	
CHLOROBENZENE		0.5	D	
CHLORODIBROMOMETHANE	NE	0.5	D	
CHLOROETHANE		0.5	D	
CHLOROFORM		0.5	n	
CHLOROMETHANE		9.0	n	
CIS-1,2-DICHLOROETHENE	1111	0.5	D	
CIS-1,3-DICHLOROPROPENE	ШN	0.5 U	D	
CYCLOHEXANE		0.5	5	
DICHLORODIFLUOROMETHANE	HANE	0.5	5	
ETHYLBENZENE		0.5	D	
ISOPROPYLBENZENE		0.5	n	
M+P-XYLENES		ł	D	
2 of 4				

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PROJ_NO: 00622	NSAMPLE	BP0W1-3-20101208	208	BP0W2-1-20101208	01208		BP0W2-2-20101208	8	BP0W-DUP01-20101208	1-201012	8
SDG: B4488	LAB_ID	B4488-06		B4488-02			B4488-05		B4488-07		
FRACTION: OV	SAMP_DATE	12/8/2010		12/8/2010			12/8/2010		12/8/2010		
MEDIA: WATER	QC_TYPE	NM		MN			MN		MN		
	UNITS	NG/L		NG/L			NG/L		NG/L		
	PCT_SOLIDS	0.0		0.0			0.0		0.0		
	DUP_OF								BP0W1-3-20101208	101208	
PARAMETER		RESULT V	VQL QLCD	RESULT	VaL	QLCD	RESULT VQL	L QLCD	RESULT	Val	OLCD -
METHYL ACETATE		0.5 UJ) C	0.5	0.5 UJ	с	0.5 UJ	U	0	0.5 U	
METHYL CYCLOHEXANE		0.5 U		0.5	D		0.5 U		ō	0.5 U	
METHYL TERT-BUTYL ETHER	HER	0.5 U		0.5	D		0.5 U		0.5	5 U	
METHYLENE CHLORIDE		0.5 U		0.5	n		0.5 U		0.5	5 U	
O-XYLENE		0.5 U		0.5	D		0.5 U	-	0.5	5 U	
STYRENE		0.5 U		0.5	D		0.5 U		0.5	5 U	
TETRACHLOROETHENE		0.5 U		0.5 U	D		0.5 U		0.5	5 U	
TOLUENE		0.5 U		0.5	n		0.5 U		0.5	5 U	
TRANS-1,2-DICHLOROETHENE	IENE	0.5 U		0.5	n		0.5 U		0.5	5 U	
TRANS-1,3-DICHLOROPROPENE	OPENE	0.5 U		0.5	5		0.5 U		0.5	5 U	
TRICHLOROETHENE		0.5 U		0.5	5		0.5 U		°0	0.5 U	
TRICHLOROFLUOROMETHANE	HANE	0.5 U		0.5 U	D		0.5 U		0.5	5 U	
VINYL CHLORIDE		0.5 U		0.5	∍		0.5 U		0.5	5 U	

1/14/2011

PROJ_NO: 00622	NSAMPLE	BP-TB01-20101208	01208	
SDG: B4488		B4488-01		
FRACTION: OV	SAMP_DATE	12/8/2010		
MEDIA: WATER	QC_TYPE	MN		
	UNITS	NG/L		
	PCT_SOLIDS	0.0		
	DUP_OF			
PARAMETER		RESULT	VQL	alcd
METHYL ACETATE		0.5	0.5 UJ	
METHYL CYCLOHEXANE		0.5 U	5	
METHYL TERT-BUTYL ETHER	IER	0.5	_	
METHYLENE CHLORIDE		0.5	D	
O-XYLENE		0.5 U	Ъ	
STYRENE		0.5	5	
TETRACHLOROETHENE		0.5	D	
TOLUENE		0.5 U	n	
TRANS-1,2-DICHLOROETHENE	IENE	0.5 U	Ъ	
TRANS-1,3-DICHLOROPROPENE	DENE	0.5 U	5	
TRICHLOROETHENE		0.5 U	Л	
TRICHLOROFLUOROMETHANE	IANE	0.5 U	n	
VINYL CHLORIDE		0.5 U	n	