

SITE LOGIC Report

QuantArray[®]-Chlor Study

Contact: Bernard Bockisch

Phone: (505) 234-1105

Address: EA Engineering
320 Gold Ave SW
Suite 1300
Albuquerque, NM 87102

Email: bbockisch@eaest.com

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Project: KAFB-Bulk Fuels Fac Vadose Zone, 62735DM02

Comments: Please note the total bacterial biomass on samples BMS9-181018-477, BM-S5-231018-491, BM-S5-231018-506, BM-S4-051118-494, BM-S2-161118-499, BM-S3-201118-489, and BM-S3-211118-494 was low, and the samples may have PCR inhibition.

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The QuantArray®-Chlor Approach

Quantification of *Dehalococcoides*, the only known bacterial group capable of complete reductive dechlorination of PCE and TCE to ethene, has become an indispensable component of assessment, remedy selection, and performance monitoring at sites impacted by chlorinated solvents. While undeniably a key group of halorespiring bacteria, *Dehalococcoides* are not the only bacteria of interest in the subsurface because reductive dechlorination is not the only potential biodegradation pathway operative at contaminated sites, and chlorinated ethenes are not always the primary contaminants of concern. The QuantArray®-Chlor not only includes a variety of halorespiring bacteria (*Dehalococcoides*, *Dehalobacter*, *Dehalogenimonas*, etc.) to assess the potential for reductive dechlorination of chloroethenes, chloroethanes, chlorobenzenes, chlorophenols, and chloroform, but also provides quantification of functional genes involved in aerobic (co)metabolic pathways for biodegradation of chlorinated solvents and even competing biological processes. Thus, the QuantArray®-Chlor will give site managers the ability to simultaneously yet economically evaluate the potential for biodegradation of a spectrum of common chlorinated contaminants through a multitude of anaerobic and aerobic (co) metabolic pathways to give a much more clear and comprehensive view of contaminant biodegradation.

The QuantArray®-Chlor is used to quantify specific microorganisms and functional genes to evaluate the following:

Anaerobic Reductive Dechlorination

Quantification of important halorespiring bacteria (e.g. *Dehalococcoides*, *Dehalobacter*, *Dehalogenimonas*, *Desulfotobacterium* spp.) and key functional genes (e.g. vinyl chloride reductases, TCE reductase, chloroform reductase) responsible for reductive dechlorination of a broad spectrum of chlorinated solvents.

Aerobic Cometabolism

Several different types of bacteria including methanotrophs and some toluene/phenol utilizing bacteria can co-oxidize TCE, DCE, and vinyl chloride. The QuantArray®-Chlor quantifies functional genes like soluble methane monooxygenase encoding enzymes capable of co-oxidation of chlorinated ethenes.

Aerobic (Co)metabolism of Vinyl Chloride

Ethene oxidizing bacteria are capable of cometabolism of vinyl chloride. In some cases, ethenotrophs can also utilize vinyl chloride as a growth supporting substrate. The QuantArray®-Chlor targets key functional genes in ethene metabolism.

How do QuantArrays® work?

The QuantArray®-Chlor in many respects is a hybrid technology combining the highly parallel detection of microarrays with the accurate and precise quantification provided by qPCR into a single platform. The key to highly parallel qPCR reactions is the nanoliter fluidics platform for low volume, solution phase qPCR reactions.

How are QuantArray® results reported?

One of the primary advantages of the QuantArray®-Chlor is the simultaneous quantification of a broad spectrum of different microorganisms and key functional genes involved in a variety of pathways for chlorinated hydrocarbon biodegradation. However, highly parallel quantification combined with the various metabolic and cometabolic capabilities of different target organisms can complicate data presentation. Therefore, in addition to Summary Tables, QuantArray® results will be presented as Microbial Population Summary and Comparison Figures to aid in data interpretation and subsequent evaluation of site management activities.

Types of Tables and Figures:

Microbial Population Summary

Figure presenting the concentrations of QuantArray®-Chlor target populations (e.g. *Dehalococcoides*) and functional genes (e.g. vinyl chloride reductase) relative to typically observed values.

Summary Tables

Tables of target population concentrations grouped by biodegradation pathway and contaminant type.

Comparison Figures

Depending on the project, sample results can be presented to compare changes over time or examine differences in microbial populations along a transect of the dissolved plume.

Results

Table 1: Summary of the QuantArray®-Chlor results obtained for samples BM-S9-171018-475, BM-S9-181018-477, BM-S5-231018-491, BM-S5-231018-506, and BM-S9-301018-496.

Sample Name	BM-S9-171018-475	BM-S9-181018-477	BM-S5-231018-491	BM-S5-231018-506	BM-S9-301018-496
Sample Date	10/17/2018	10/18/2018	10/23/2018	10/23/2018	10/30/2018
<i>Reductive Dechlorination</i>	cells/g	cells/g	cells/g	cells/g	cells/g
<i>Dehalococcoides</i> (DHC)	<1.00E+03	<2.00E+03	<1.00E+03	<1.00E+03	<1.00E+03
tceA Reductase (TCE)	<1.00E+03	<2.00E+03	<1.00E+03	<1.00E+03	<1.00E+04
BAV1 Vinyl Chloride Reductase (BVC)	<1.00E+03	<2.00E+03	<1.00E+03	<1.00E+03	<1.00E+04
Vinyl Chloride Reductase (VCR)	<1.00E+03	<2.00E+03	<1.00E+03	<1.00E+03	<1.00E+04
<i>Dehalobacter</i> spp. (DHBt)	2.12E+06	<2.00E+04	<1.00E+04	<1.00E+04	3.83E+05
<i>Dehalobacter</i> DCM (DCM)	<1.00E+04	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalogenimonas</i> spp. (DHG)	<1.00E+04	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
cerA Reductase (CER)	<1.00E+04	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
trans-1,2-DCE Reductase (TDR)	<1.00E+04	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Desulfitobacterium</i> spp. (DSB)	8.01E+03 (J)	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalobium chlorocoercia</i> (DECO)	3.02E+05	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Desulfuromonas</i> spp. (DSM)	<1.00E+04	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
PCE Reductase (PCE-1)	<1.00E+04	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
PCE Reductase (PCE-2)	<1.00E+04	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Chloroform Reductase (CFR)	<1.00E+04	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
1,1 DCA Reductase (DCA)	<1.00E+04	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
1,2 DCA Reductase (DCAR)	<1.00E+04	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Aerobic (Co)Metabolic</i>					
Soluble Methane Monooxygenase (SMMO)	<1.00E+04	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Dioxygenase (TOD)	<1.00E+04	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Phenol Hydroxylase (PHE)	1.06E+06	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Trichlorobenzene Dioxygenase (TCBO)	<1.00E+04	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Monooxygenase 2 (RDEG)	5.30E+05	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Monooxygenase (RMO)	6.73E+05	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Ethene Monooxygenase (EtnC)	<1.00E+04	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Epoxyalkane Transferase (EtnE)	<1.00E+04	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Dichloromethane Dehalogenase (DCMA)	<1.00E+04	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Other</i>					
Total Eubacteria (EBAC)	2.69E+07	3.22E+04 (I)	2.86E+05 (I)	3.39E+04 (I)	1.02E+06
Sulfate Reducing Bacteria (APS)	1.08E+07	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Methanogens (MGN)	<1.00E+04	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>CENSUS Targets</i>					
Benzoyl Coenzyme A Reductase (BCR)	3.80E+04				1.94E+03 (J)
Benzene Carboxylase (abcA)	<1.00E+04				9.67E+02 (J)

Legend:

NA = Not Analyzed
I = Inhibited

NS = Not Sampled
< = Result Not Detected

J = Estimated Gene Copies Below PQL but Above LQL

Table 2: Summary of the QuantArray®-Chlor results obtained for samples BM-S4-041118-480, BM-S4-051118-494, BM-S2-161118-474, and BM-S2-161118-499.

Sample Name	BM-S4-041118-480	BM-S4-051118-494	BM-S2-161118-474	BM-S2-161118-499
Sample Date	11/04/2018	11/05/2018	11/16/2018	11/16/2018
<i>Reductive Dechlorination</i>	cells/g	cells/g	cells/g	cells/g
<i>Dehalococcoides</i> (DHC)	<1.00E+03	<1.00E+03	<1.00E+03	<1.00E+03
tceA Reductase (TCE)	<1.00E+04	<1.00E+04	<1.00E+03	<1.00E+03
BAV1 Vinyl Chloride Reductase (BVC)	<1.00E+04	<1.00E+04	<1.00E+03	<1.00E+03
Vinyl Chloride Reductase (VCR)	<1.00E+04	<1.00E+04	<1.00E+03	<1.00E+03
<i>Dehalobacter</i> spp. (DHBt)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalobacter</i> DCM (DCM)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalogenimonas</i> spp. (DHG)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
cerA Reductase (CER)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
trans-1,2-DCE Reductase (TDR)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Desulfitobacterium</i> spp. (DSB)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalobium chlorocoercia</i> (DECO)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Desulfuromonas</i> spp. (DSM)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
PCE Reductase (PCE-1)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
PCE Reductase (PCE-2)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Chloroform Reductase (CFR)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
1,1 DCA Reductase (DCA)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
1,2 DCA Reductase (DCAR)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Aerobic (Co)Metabolic</i>				
Soluble Methane Monooxygenase (SMMO)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Dioxygenase (TOD)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Phenol Hydroxylase (PHE)	2.42E+03 (J)	<1.00E+04	1.17E+06	<1.00E+04
Trichlorobenzene Dioxygenase (TCBO)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Monooxygenase 2 (RDEG)	1.53E+05	<1.00E+04	1.32E+06	<1.00E+04
Toluene Monooxygenase (RMO)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Ethene Monooxygenase (EtnC)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Epoxyalkane Transferase (EtnE)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Dichloromethane Dehalogenase (DCMA)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Other</i>				
Total Eubacteria (EBAC)	1.16E+06	5.57E+05 (I)	2.43E+06	3.59E+05 (I)
Sulfate Reducing Bacteria (APS)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Methanogens (MGN)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>CENSUS Targets</i>				
Benzoyl Coenzyme A Reductase (BCR)	5.26E+04		1.57E+03 (J)	
Benzene Carboxylase (abcA)	<1.00E+04		<1.00E+04	

Legend:

NA = Not Analyzed

I = Inhibited

NS = Not Sampled

< = Result Not Detected

J = Estimated Gene Copies Below PQL but Above LQL

Table 3: Summary of the QuantArray®-Chlor results obtained for samples BM-S3-201118-477, BM-S3-201118-489, and BM-S3-211118-494.

Sample Name Sample Date	BM-S3-201118-477 11/20/2018	BM-S3-201118-489 11/20/2018	BM-S3-211118-494 11/21/2018
<i>Reductive Dechlorination</i>	cells/g	cells/g	cells/g
<i>Dehalococcoides</i> (DHC)	<1.00E+03	<1.00E+03	<1.00E+03
tceA Reductase (TCE)	<1.00E+03	<1.00E+03	<1.00E+03
BAV1 Vinyl Chloride Reductase (BVC)	<1.00E+03	<1.00E+03	<1.00E+03
Vinyl Chloride Reductase (VCR)	<1.00E+03	<1.00E+03	<1.00E+03
<i>Dehalobacter</i> spp. (DHBt)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalobacter</i> DCM (DCM)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalogenimonas</i> spp. (DHG)	<1.00E+04	<1.00E+04	<1.00E+04
cerA Reductase (CER)	<1.00E+04	<1.00E+04	<1.00E+04
trans-1,2-DCE Reductase (TDR)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Desulfitobacterium</i> spp. (DSB)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalobium chlorocoercia</i> (DECO)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Desulfuromonas</i> spp. (DSM)	<1.00E+04	<1.00E+04	<1.00E+04
PCE Reductase (PCE-1)	<1.00E+04	<1.00E+04	<1.00E+04
PCE Reductase (PCE-2)	<1.00E+04	<1.00E+04	<1.00E+04
Chloroform Reductase (CFR)	<1.00E+04	<1.00E+04	<1.00E+04
1,1 DCA Reductase (DCA)	<1.00E+04	<1.00E+04	<1.00E+04
1,2 DCA Reductase (DCAR)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Aerobic (Co)Metabolic</i>			
Soluble Methane Monooxygenase (SMMO)	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Dioxygenase (TOD)	<1.00E+04	<1.00E+04	<1.00E+04
Phenol Hydroxylase (PHE)	<1.00E+04	<1.00E+04	<1.00E+04
Trichlorobenzene Dioxygenase (TCBO)	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Monooxygenase 2 (RDEG)	9.93E+04	<1.00E+04	<1.00E+04
Toluene Monooxygenase (RMO)	<1.00E+04	<1.00E+04	<1.00E+04
Ethene Monooxygenase (EtnC)	<1.00E+04	<1.00E+04	<1.00E+04
Epoxyalkane Transferase (EtnE)	<1.00E+04	<1.00E+04	<1.00E+04
Dichloromethane Dehalogenase (DCMA)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Other</i>			
Total Eubacteria (EBAC)	5.73E+05	9.78E+03 (I)	6.99E+03 (I)
Sulfate Reducing Bacteria (APS)	<1.00E+04	<1.00E+04	<1.00E+04
Methanogens (MGN)	<1.00E+04	<1.00E+04	<1.00E+04
<i>CENSUS Targets</i>			
Benzoyl Coenzyme A Reductase (BCR)	1.42E+04		
Benzene Carboxylase (abcA)	<1.00E+04		

Legend:

NA = Not Analyzed
I = Inhibited

NS = Not Sampled
< = Result Not Detected

J = Estimated Gene Copies Below PQL but Above LQL

Microbial Populations BM-S9-171018-475

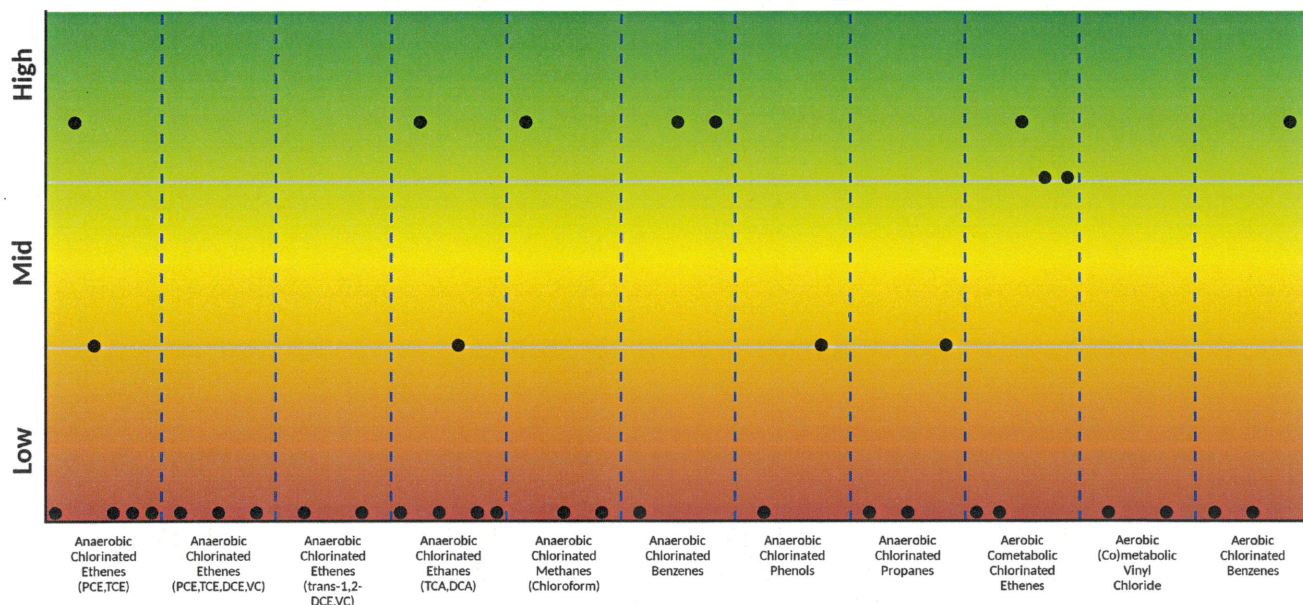


Figure 1: Microbial population summary to aid in evaluating potential pathways and biodegradation of specific contaminants.

Anaerobic - Reductive Dechlorination or Dichloroelimination		Aerobic - (Co)metabolism	
Chlorinated Ethenes (PCE, TCE)	DHC, DHBt, DSB, DSM, PCE-1, PCE-2	Chlorinated Ethenes (TCE,DCE,VC)	sMMO, TOD, PHE, RDEG, RMO
Chlorinated Ethenes (PCE, TCE, DCE, VC)	DHC, BVC, VCR	(Co)metabolic Vinyl Chloride	etnC, etnE
Chlorinated Ethenes (trans-1,2-DCE, VC)	TDR, CER	Chlorinated Benzenes	TOD, TCBO, PHE
Chlorinated Ethanes (TCA and 1,2-DCA)	DHC, DHBt, DHG, DSB ¹ , DCA, DCAR		
Chlorinated Methanes (Chloroform)	DHBt, DCM, CFR		
Chlorinated Benzenes	DHC, DHBt ² , DECO		
Chlorinated Phenols	DHC, DSB		
Chlorinated Propanes	DHC, DHG, DSB ¹		

¹Desulfotobacterium dichloroeliminans DCA1. ²Implicated in reductive dechlorination of dichlorobenzene and potentially chlorobenzene.

Microbial Populations BM-S9-181018-477

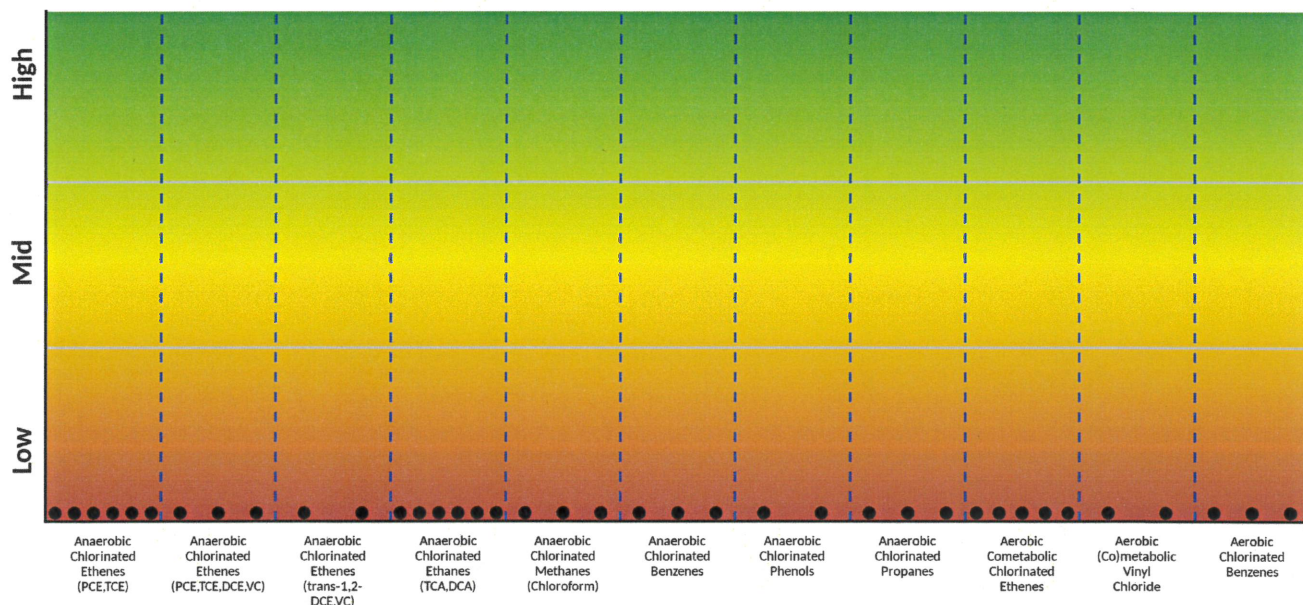


Figure 2: Microbial population summary to aid in evaluating potential pathways and biodegradation of specific contaminants.

Anaerobic - Reductive Dechlorination or Dichloroelimination			Aerobic - (Co)metabolism	
Chlorinated Ethenes (PCE, TCE)	DHC, DHBt, DSB, DSM, PCE-1, PCE-2	Chlorinated Ethenes (TCE,DCE,VC)	sMMO, TOD, PHE, RDEG, RMO	
Chlorinated Ethenes (PCE, TCE, DCE, VC)	DHC, BVC, VCR	(Co)metabolic Vinyl Chloride	etnC, etnE	
Chlorinated Ethenes (trans-1,2-DCE, VC)	TDR, CER	Chlorinated Benzenes	TOD, TCBO, PHE	
Chlorinated Ethanes (TCA and 1,2-DCA)	DHC, DHBt, DHG, DSB ¹ , DCA, DCAR			
Chlorinated Methanes (Chloroform)	DHBt, DCM, CFR			
Chlorinated Benzenes	DHC, DHBt ² , DECO			
Chlorinated Phenols	DHC, DSB			
Chlorinated Propanes	DHC, DHG, DSB ¹			

¹ *Desulfotobacterium dichloroeliminans* DCA1. ² Implicated in reductive dechlorination of dichlorobenzene and potentially chlorobenzene.

Microbial Populations BM-S5-231018-491

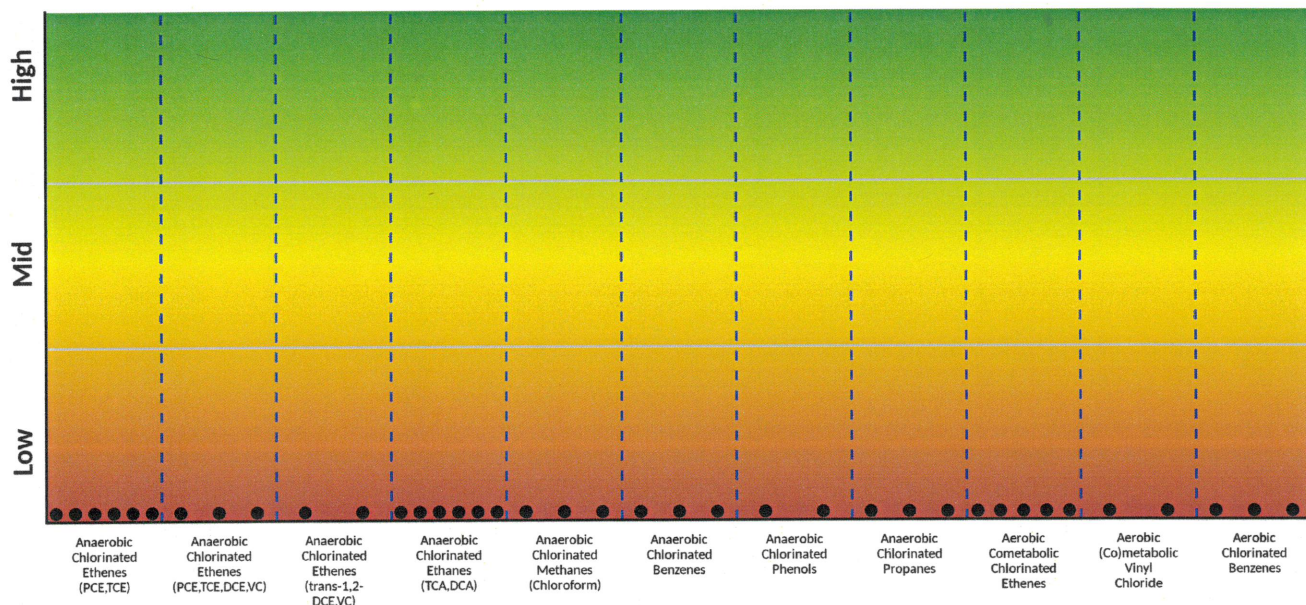


Figure 3: Microbial population summary to aid in evaluating potential pathways and biodegradation of specific contaminants.

Anaerobic - Reductive Dechlorination or Dichloroelimination

Chlorinated Ethenes (PCE, TCE)	DHC, DHBt, DSB, DSM, PCE-1, PCE-2
Chlorinated Ethenes (PCE, TCE, DCE, VC)	DHC, BVC, VCR
Chlorinated Ethenes (trans-1,2-DCE, VC)	TDR, CER
Chlorinated Ethanes (TCA and 1,2-DCA)	DHC, DHBt, DHG, DSB ¹ , DCA, DCAR
Chlorinated Methanes (Chloroform)	DHBt, DCM, CFR
Chlorinated Benzenes	DHC, DHBt ² , DECO
Chlorinated Phenols	DHC, DSB
Chlorinated Propanes	DHC, DHG, DSB ¹

Aerobic - (Co)metabolism

Chlorinated Ethenes (TCE,DCE,VC)	sMMO, TOD, PHE, RDEG, RMO
(Co)metabolic Vinyl Chloride	etnC, etnE
Chlorinated Benzenes	TOD, TCBO, PHE

¹ *Desulfotobacterium dichloroeliminans* DCA1. ² Implicated in reductive dechlorination of dichlorobenzene and potentially chlorobenzene.

Microbial Populations BM-S5-231018-506

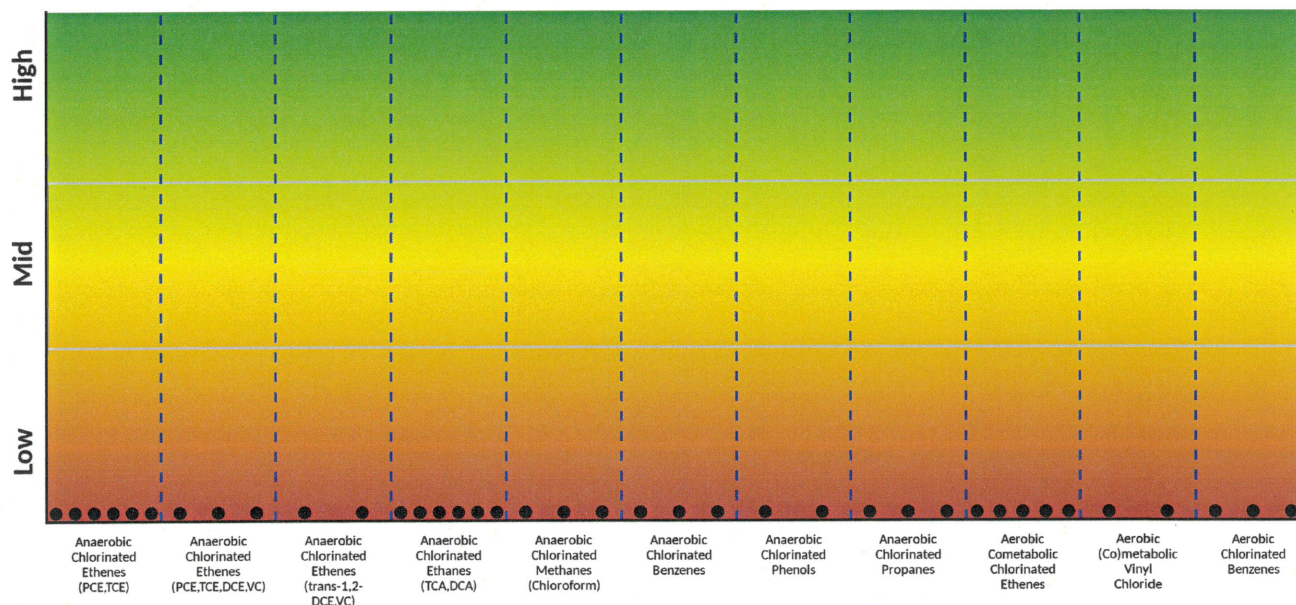


Figure 4: Microbial population summary to aid in evaluating potential pathways and biodegradation of specific contaminants.

Anaerobic - Reductive Dechlorination or Dichloroelimination		Aerobic - (Co)metabolism	
Chlorinated Ethenes (PCE, TCE)	DHC, DHBt, DSB, DSM, PCE-1, PCE-2	Chlorinated Ethenes (TCE,DCE,VC)	sMMO, TOD, PHE, RDEG, RMO
Chlorinated Ethenes (PCE, TCE, DCE, VC)	DHC, BVC, VCR	(Co)metabolic Vinyl Chloride	etnC, etnE
Chlorinated Ethenes (trans-1,2-DCE, VC)	TDR, CER	Chlorinated Benzenes	TOD, TCBO, PHE
Chlorinated Ethanes (TCA and 1,2-DCA)	DHC, DHBt, DHG, DSB ¹ , DCA, DCAR		
Chlorinated Methanes (Chloroform)	DHBt, DCM, CFR		
Chlorinated Benzenes	DHC, DHBt ² , DECO		
Chlorinated Phenols	DHC, DSB		
Chlorinated Propanes	DHC, DHG, DSB ¹		

¹ *Desulfotobacterium dichloroeliminans* DCA1. ² Implicated in reductive dechlorination of dichlorobenzene and potentially chlorobenzene.

Microbial Populations BM-S9-301018-496

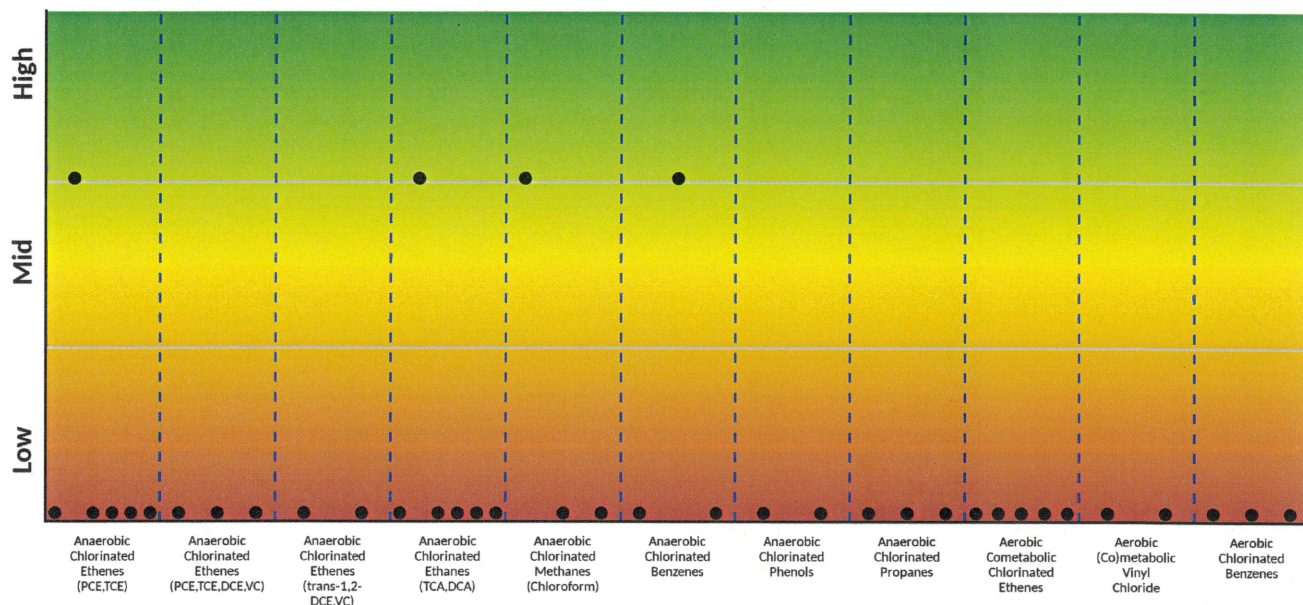


Figure 5: Microbial population summary to aid in evaluating potential pathways and biodegradation of specific contaminants.

Anaerobic - Reductive Dechlorination or Dichloroelimination			Aerobic - (Co)metabolism	
Chlorinated Ethenes (PCE, TCE)	DHC, DHBt, DSB, DSM, PCE-1, PCE-2	Chlorinated Ethenes (TCE,DCE,VC)	sMMO, TOD, PHE, RDEG, RMO	
Chlorinated Ethenes (PCE, TCE, DCE, VC)	DHC, BVC, VCR	(Co)metabolic Vinyl Chloride	etnC, etnE	
Chlorinated Ethenes (trans-1,2-DCE, VC)	TDR, CER	Chlorinated Benzenes	TOD, TCBO, PHE	
Chlorinated Ethanes (TCA and 1,2-DCA)	DHC, DHBt, DHG, DSB ¹ , DCA, DCAR			
Chlorinated Methanes (Chloroform)	DHBt, DCM, CFR			
Chlorinated Benzenes	DHC, DHBt ² , DECO			
Chlorinated Phenols	DHC, DSB			
Chlorinated Propanes	DHC, DHG, DSB ¹			

¹ *Desulfotobacterium dichloroeliminans* DCA1. ² Implicated in reductive dechlorination of dichlorobenzene and potentially chlorobenzene.

Microbial Populations BM-S4-041118-480

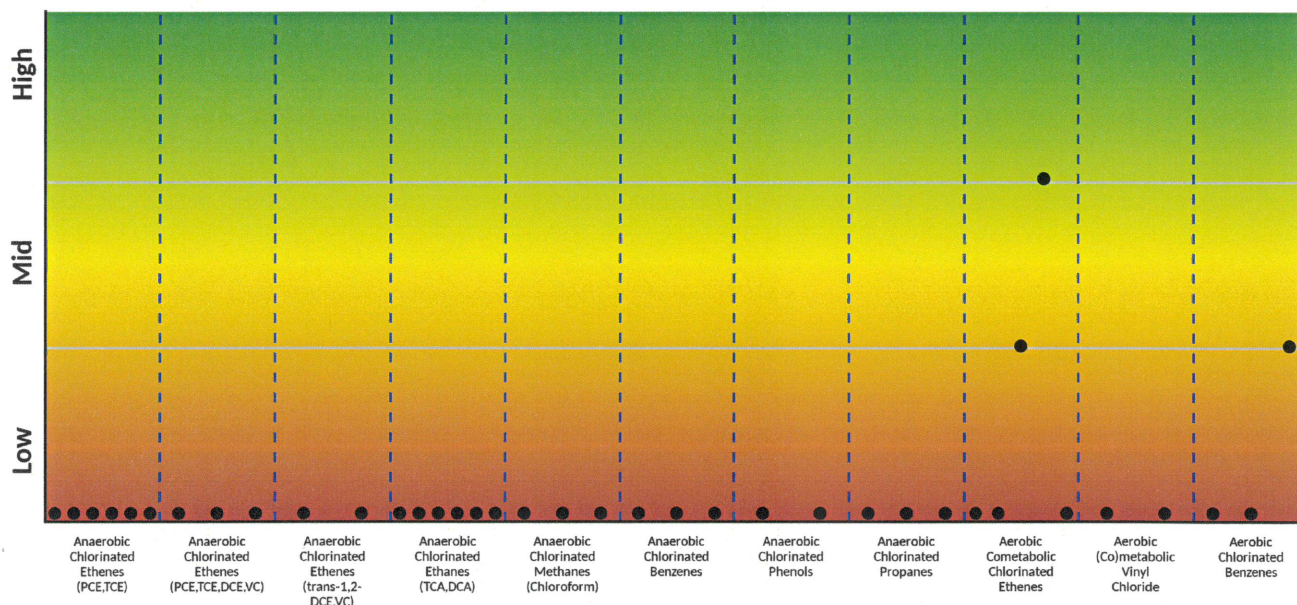


Figure 6: Microbial population summary to aid in evaluating potential pathways and biodegradation of specific contaminants.

Anaerobic - Reductive Dechlorination or Dichloroelimination

Chlorinated Ethenes (PCE, TCE)	DHC, DHBt, DSB, DSM, PCE-1, PCE-2
Chlorinated Ethenes (PCE, TCE, DCE, VC)	DHC, BVC, VCR
Chlorinated Ethenes (trans-1,2-DCE, VC)	TDR, CER
Chlorinated Ethanes (TCA and 1,2-DCA)	DHC, DHBt, DHG, DSB ¹ , DCA, DCAR
Chlorinated Methanes (Chloroform)	DHBt, DCM, CFR
Chlorinated Benzenes	DHC, DHBt ² , DECO
Chlorinated Phenols	DHC, DSB
Chlorinated Propanes	DHC, DHG, DSB ¹

Aerobic - (Co)metabolism

Chlorinated Ethenes (TCE,DCE,VC) (Co)metabolic Vinyl Chloride	sMMO, TOD, PHE, RDEG, RMO etnC, etnE
Chlorinated Benzenes	TOD, TCBO, PHE

¹ *Desulfotobacterium dichloroeliminans* DCA1. ² Implicated in reductive dechlorination of dichlorobenzene and potentially chlorobenzene.

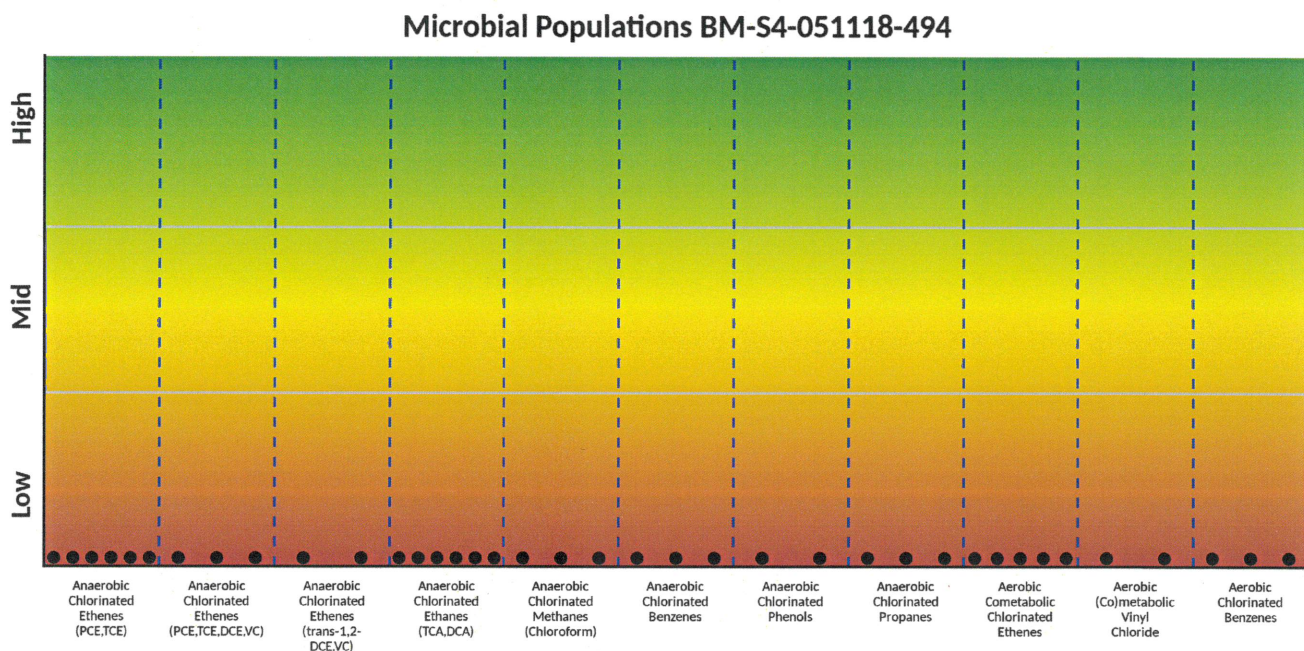


Figure 7: Microbial population summary to aid in evaluating potential pathways and biodegradation of specific contaminants.

Anaerobic - Reductive Dechlorination or Dichloroelimination

Chlorinated Ethenes (PCE, TCE)	DHC, DHBt, DSB, DSM, PCE-1, PCE-2
Chlorinated Ethenes (PCE, TCE, DCE, VC)	DHC, BVC, VCR
Chlorinated Ethenes (trans-1,2-DCE, VC)	TDR, CER
Chlorinated Ethanes (TCA and 1,2-DCA)	DHC, DHBt, DHG, DSB ¹ , DCA, DCAR
Chlorinated Methanes (Chloroform)	DHBt, DCM, CFR
Chlorinated Benzenes	DHC, DHBt ² , DECO
Chlorinated Phenols	DHC, DSB
Chlorinated Propanes	DHC, DHG, DSB ¹

Aerobic - (Co)metabolism

Chlorinated Ethenes (TCE,DCE,VC)	sMMO, TOD, PHE, RDEG, RMO
(Co)metabolic Vinyl Chloride	etnC, etnE
Chlorinated Benzenes	TOD, TCBO, PHE

¹*Desulfotobacterium dichloroelimians* DCA1. ²Implicated in reductive dechlorination of dichlorobenzene and potentially chlorobenzene.

Microbial Populations BM-S2-161118-474

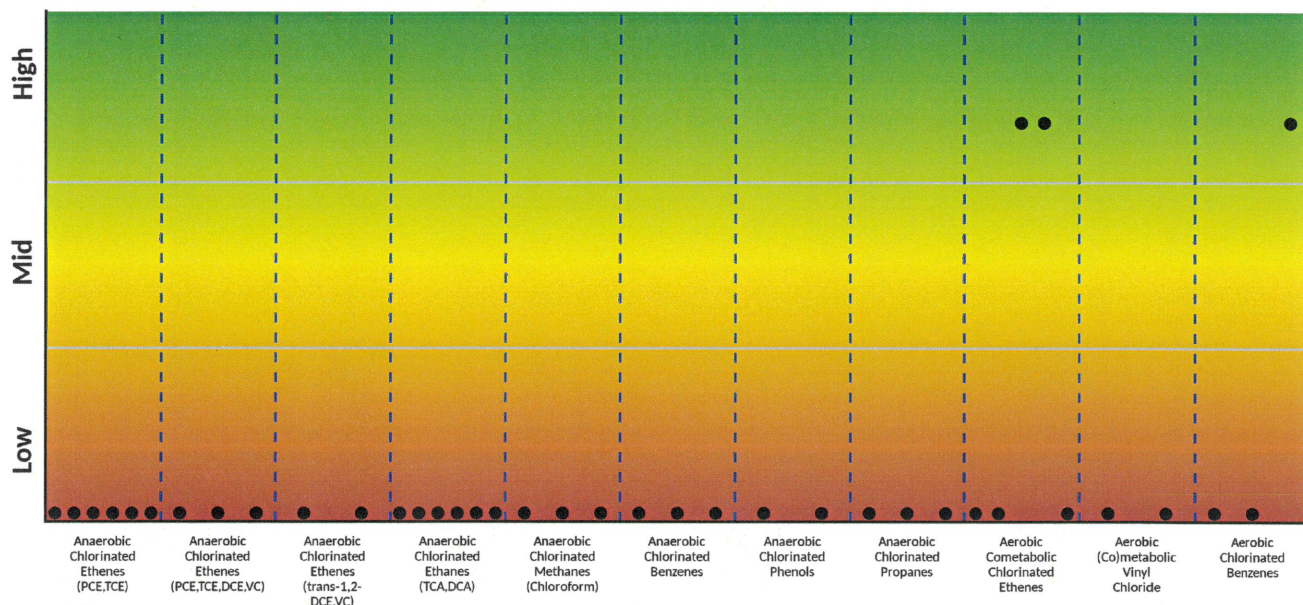


Figure 8: Microbial population summary to aid in evaluating potential pathways and biodegradation of specific contaminants.

Anaerobic - Reductive Dechlorination or Dichloroelimination		Aerobic - (Co)metabolism	
Chlorinated Ethenes (PCE, TCE)	DHC, DHBt, DSB, DSM, PCE-1, PCE-2	Chlorinated Ethenes (TCE,DCE,VC)	sMMO, TOD, PHE, RDEG, RMO
Chlorinated Ethenes (PCE, TCE, DCE, VC)	DHC, BVC, VCR	(Co)metabolic Vinyl Chloride	etnC, etnE
Chlorinated Ethenes (trans-1,2-DCE, VC)	TDR, CER	Chlorinated Benzenes	TOD, TCBO, PHE
Chlorinated Ethanes (TCA and 1,2-DCA)	DHC, DHBt, DHG, DSB ¹ , DCA, DCAR		
Chlorinated Methanes (Chloroform)	DHBt, DCM, CFR		
Chlorinated Benzenes	DHC, DHBt ² , DECO		
Chlorinated Phenols	DHC, DSB		
Chlorinated Propanes	DHC, DHG, DSB ¹		

¹ *Desulfotobacterium dichloroelimians* DCA1. ² Implicated in reductive dechlorination of dichlorobenzene and potentially chlorobenzene.

Microbial Populations BM-S2-161118-499

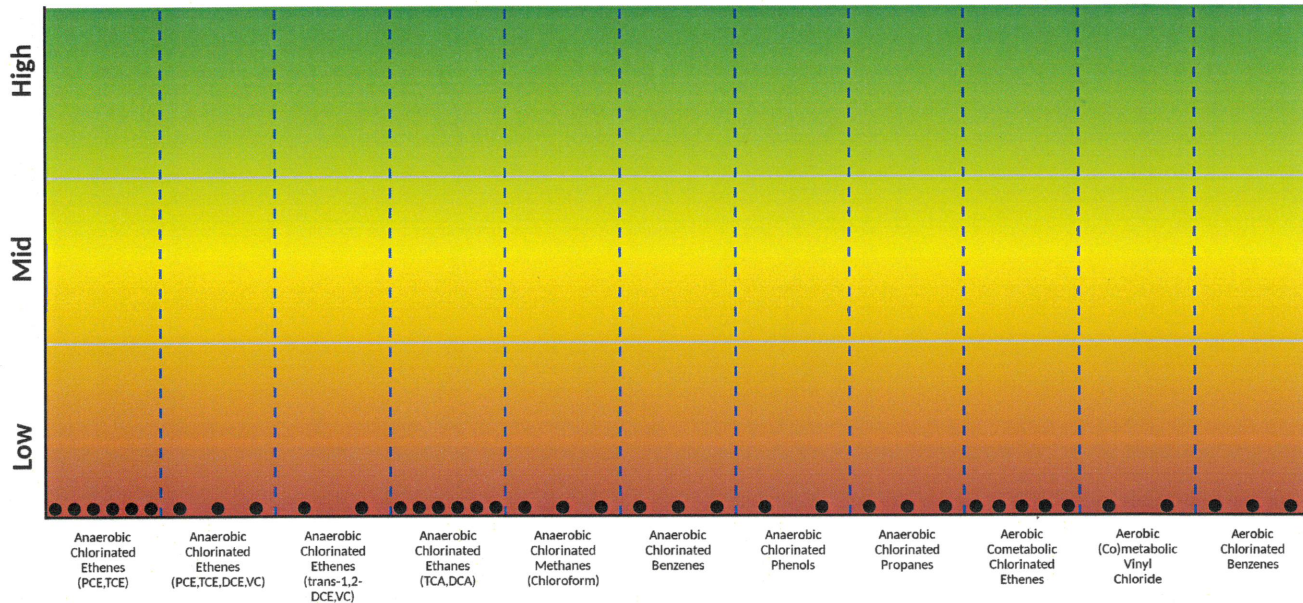


Figure 9: Microbial population summary to aid in evaluating potential pathways and biodegradation of specific contaminants.

Anaerobic - Reductive Dechlorination or Dichloroelimination			Aerobic - (Co)metabolism	
Chlorinated Ethenes (PCE, TCE)	DHC, DHBt, DSB, DSM, PCE-1, PCE-2	Chlorinated Ethenes (TCE,DCE,VC)	sMMO, TOD, PHE, RDEG, RMO	
Chlorinated Ethenes (PCE, TCE, DCE, VC)	DHC, BVC, VCR	(Co)metabolic Vinyl Chloride	etnC, etnE	
Chlorinated Ethenes (trans-1,2-DCE, VC)	TDR, CER	Chlorinated Benzenes	TOD, TCBO, PHE	
Chlorinated Ethanes (TCA and 1,2-DCA)	DHC, DHBt, DHG, DSB ¹ , DCA, DCAR			
Chlorinated Methanes (Chloroform)	DHBt, DCM, CFR			
Chlorinated Benzenes	DHC, DHBt ² , DECO			
Chlorinated Phenols	DHC, DSB			
Chlorinated Propanes	DHC, DHG, DSB ¹			

¹ *Desulfotobacterium dichloroeliminans* DCA1. ² Implicated in reductive dechlorination of dichlorobenzene and potentially chlorobenzene.

Microbial Populations BM-S3-201118-477

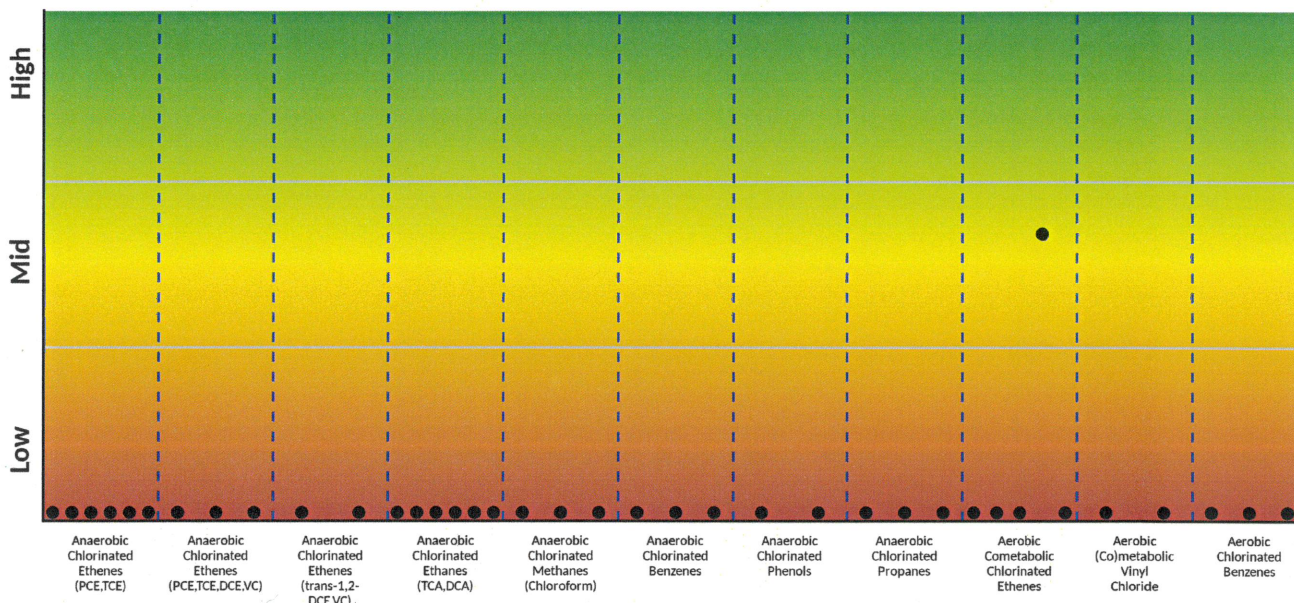


Figure 10: Microbial population summary to aid in evaluating potential pathways and biodegradation of specific contaminants.

Anaerobic - Reductive Dechlorination or Dichloroelimination

Chlorinated Ethenes (PCE, TCE)	DHC, DHBt, DSB, DSM, PCE-1, PCE-2
Chlorinated Ethenes (PCE, TCE, DCE, VC)	DHC, BVC, VCR
Chlorinated Ethenes (trans-1,2-DCE, VC)	TDR, CER
Chlorinated Ethanes (TCA and 1,2-DCA)	DHC, DHBt, DHG, DSB ¹ , DCA, DCAR
Chlorinated Methanes (Chloroform)	DHBt, DCM, CFR
Chlorinated Benzenes	DHC, DHBt ² , DECO
Chlorinated Phenols	DHC, DSB
Chlorinated Propanes	DHC, DHG, DSB ¹

Aerobic - (Co)metabolism

Chlorinated Ethenes (TCE,DCE,VC) (Co)metabolic Vinyl Chloride	sMMO, TOD, PHE, RDEG, RMO etnC, etnE
Chlorinated Benzenes	TOD, TCBO, PHE

¹*Desulfitobacterium dichloroeliminans* DCA1. ²Implicated in reductive dechlorination of dichlorobenzene and potentially chlorobenzene.

Microbial Populations BM-S3-201118-489

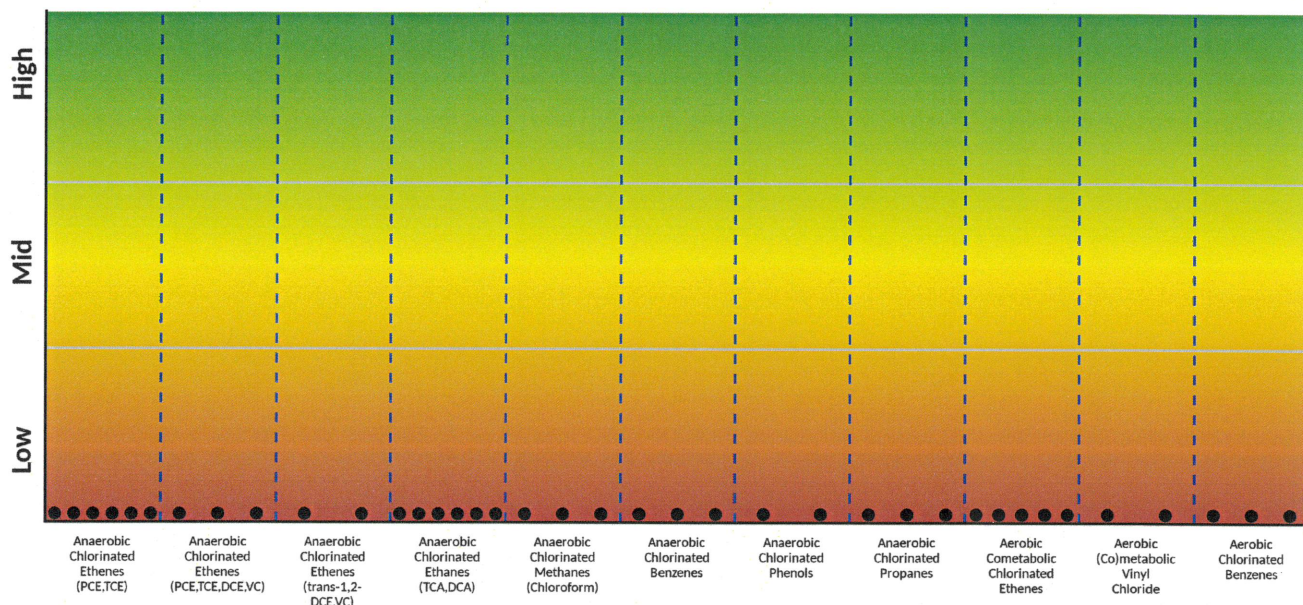


Figure 11: Microbial population summary to aid in evaluating potential pathways and biodegradation of specific contaminants.

Anaerobic - Reductive Dechlorination or Dichloroelimination

Chlorinated Ethenes (PCE, TCE)	DHC, DHBt, DSB, DSM, PCE-1, PCE-2
Chlorinated Ethenes (PCE, TCE, DCE, VC)	DHC, BVC, VCR
Chlorinated Ethenes (trans-1,2-DCE, VC)	TDR, CER
Chlorinated Ethanes (TCA and 1,2-DCA)	DHC, DHBt, DHG, DSB ¹ , DCA, DCAR
Chlorinated Methanes (Chloroform)	DHBt, DCM, CFR
Chlorinated Benzenes	DHC, DHBt ² , DECO
Chlorinated Phenols	DHC, DSB
Chlorinated Propanes	DHC, DHG, DSB ¹

Aerobic - (Co)metabolism

Chlorinated Ethenes (TCE,DCE,VC)	sMMO, TOD, PHE, RDEG, RMO
(Co)metabolic Vinyl Chloride	etnC, etnE
Chlorinated Benzenes	TOD, TCBO, PHE

¹ *Desulfotobacterium dichloroeliminans* DCA1. ² Implicated in reductive dechlorination of dichlorobenzene and potentially chlorobenzene.

Microbial Populations BM-S3-211118-494

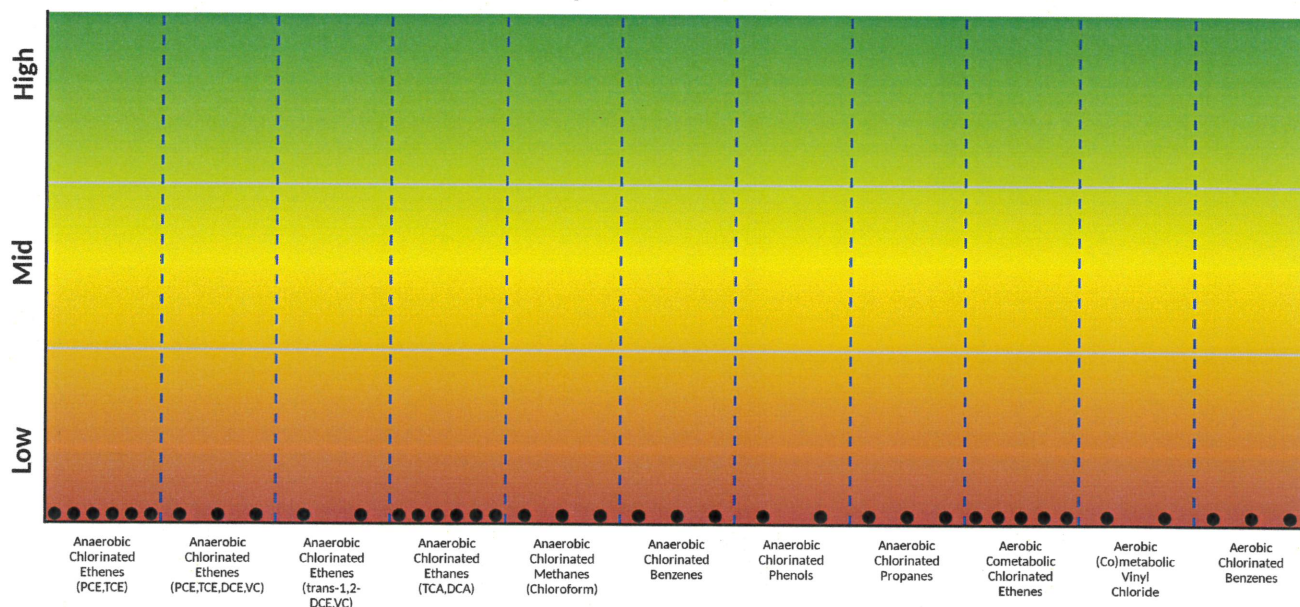


Figure 12: Microbial population summary to aid in evaluating potential pathways and biodegradation of specific contaminants.

Anaerobic - Reductive Dechlorination or Dichloroelimination			Aerobic - (Co)metabolism	
Chlorinated Ethenes (PCE, TCE)	DHC, DHBt, DSB, DSM, PCE-1, PCE-2	Chlorinated Ethenes (TCE,DCE,VC)	sMMO, TOD, PHE, RDEG, RMO	
Chlorinated Ethenes (PCE, TCE, DCE, VC)	DHC, BVC, VCR	(Co)metabolic Vinyl Chloride	etnC, etnE	
Chlorinated Ethenes (trans-1,2-DCE, VC)	TDR, CER	Chlorinated Benzenes	TOD, TCBO, PHE	
Chlorinated Ethanes (TCA and 1,2-DCA)	DHC, DHBt, DHG, DSB ¹ , DCA, DCAR			
Chlorinated Methanes (Chloroform)	DHBt, DCM, CFR			
Chlorinated Benzenes	DHC, DHBt ² , DECO			
Chlorinated Phenols	DHC, DSB			
Chlorinated Propanes	DHC, DHG, DSB ¹			

¹ *Desulfotobacterium dichloroeliminans* DCA1. ² Implicated in reductive dechlorination of dichlorobenzene and potentially chlorobenzene.

Table 4: Summary of the QuantArray®-Chlor results for microorganisms responsible for reductive dechlorination for samples BM-S9-171018-475, BM-S9-181018-477, BM-S5-231018-491, BM-S5-231018-506, and BM-S9-301018-496.

Sample Name	BM-S9-171018-475	BM-S9-181018-477	BM-S5-231018-491	BM-S5-231018-506	BM-S9-301018-496
Sample Date	10/17/2018	10/18/2018	10/23/2018	10/23/2018	10/30/2018
Reductive Dechlorination	cells/g	cells/g	cells/g	cells/g	cells/g
<i>Dehalococcoides</i> (DHC)	<1.00E+03	<2.00E+03	<1.00E+03	<1.00E+03	<1.00E+03
tceA Reductase (TCE)	<1.00E+03	<2.00E+03	<1.00E+03	<1.00E+03	<1.00E+04
BAV1 Vinyl Chloride Reductase (BVC)	<1.00E+03	<2.00E+03	<1.00E+03	<1.00E+03	<1.00E+04
Vinyl Chloride Reductase (VCR)	<1.00E+03	<2.00E+03	<1.00E+03	<1.00E+03	<1.00E+04
<i>Dehalobacter</i> spp. (DHBt)	2.12E+06	<2.00E+04	<1.00E+04	<1.00E+04	3.83E+05
<i>Dehalobacter</i> DCM (DCM)	<1.00E+04	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalogenimonas</i> spp. (DHG)	<1.00E+04	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Desulfotobacterium</i> spp. (DSB)	8.01E+03 (J)	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalobium chlorocoercia</i> (DECO)	3.02E+05	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Desulfuromonas</i> spp. (DSM)	<1.00E+04	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04

Microbial Populations - Reductive Dechlorination

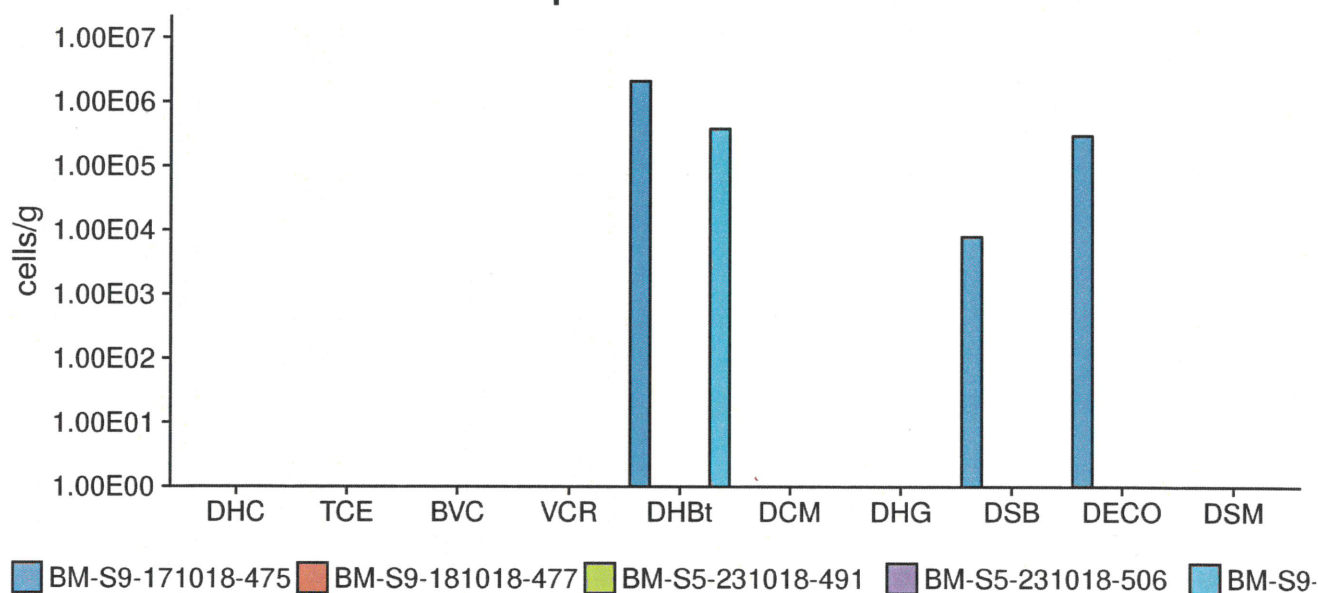


Figure 13: Comparison - microbial populations involved in reductive dechlorination.

Table 5: Summary of the QuantArray®-Chlor results for microorganisms responsible for reductive dechlorination for samples BM-S9-171018-475, BM-S9-181018-477, BM-S5-231018-491 , BM-S5-231018-506 , and BM-S9-301018-496.

Sample Name	BM-S9-171018-475	BM-S9-181018-477	BM-S5-231018-491	BM-S5-231018-506	BM-S9-301018-496
Sample Date	10/17/2018	10/18/2018	10/23/2018	10/23/2018	10/30/2018
<i>Reductive Dechlorination</i>	cells/g	cells/g	cells/g	cells/g	cells/g
Chloroform Reductase (CFR)	<1.00E+04	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
1,1 DCA Reductase (DCA)	<1.00E+04	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
1,2 DCA Reductase (DCAR)	<1.00E+04	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
PCE Reductase (PCE-1)	<1.00E+04	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
PCE Reductase (PCE-2)	<1.00E+04	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalogenimonas trans</i> -1,2-DCE Reductase (TDR)	<1.00E+04	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalogenimonas cerA</i> Reductase (CER)	<1.00E+04	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04

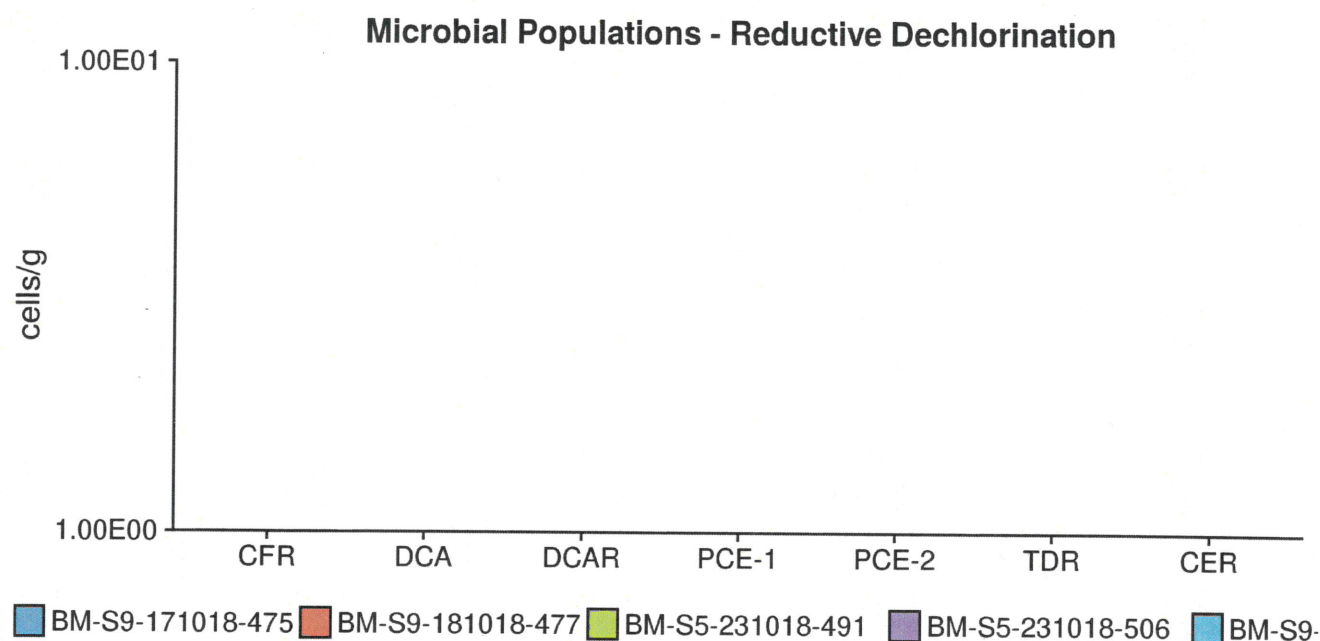


Figure 14: Comparison - microbial populations involved in reductive dechlorination.

Table 6: Summary of the QuantArray®-Chlor results for microorganisms responsible for reductive dechlorination for samples BM-S4-041118-480, BM-S4-051118-494, BM-S2-161118-474, and BM-S2-161118-499.

Sample Name	BM-S4-041118-480	BM-S4-051118-494	BM-S2-161118-474	BM-S2-161118-499
Sample Date	11/04/2018	11/05/2018	11/16/2018	11/16/2018
Reductive Dechlorination	cells/g	cells/g	cells/g	cells/g
<i>Dehalococcoides</i> (DHC)	<1.00E+03	<1.00E+03	<1.00E+03	<1.00E+03
tceA Reductase (TCE)	<1.00E+04	<1.00E+04	<1.00E+03	<1.00E+03
BAV1 Vinyl Chloride Reductase (BVC)	<1.00E+04	<1.00E+04	<1.00E+03	<1.00E+03
Vinyl Chloride Reductase (VCR)	<1.00E+04	<1.00E+04	<1.00E+03	<1.00E+03
<i>Dehalobacter</i> spp. (DHBt)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalobacter</i> DCM (DCM)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalogenimonas</i> spp. (DHG)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Desulfitobacterium</i> spp. (DSB)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalobium chlorocoercia</i> (DECO)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Desulfuromonas</i> spp. (DSM)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04

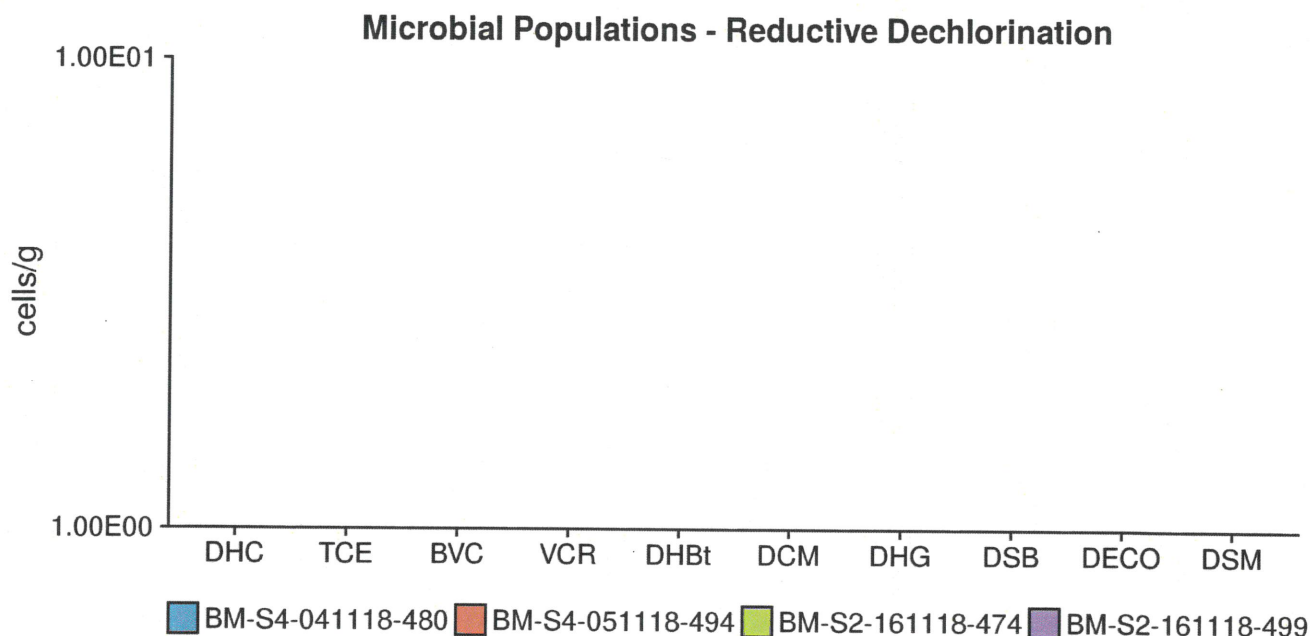


Figure 15: Comparison - microbial populations involved in reductive dechlorination.

Table 7: Summary of the QuantArray®-Chlor results for microorganisms responsible for reductive dechlorination for samples BM-S4-041118-480, BM-S4-051118-494, BM-S2-161118-474, and BM-S2-161118-499.

Sample Name	BM-S4-041118-480	BM-S4-051118-494	BM-S2-161118-474	BM-S2-161118-499
Sample Date	11/04/2018	11/05/2018	11/16/2018	11/16/2018
<i>Reductive Dechlorination</i>	cells/g	cells/g	cells/g	cells/g
Chloroform Reductase (CFR)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
1,1 DCA Reductase (DCA)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
1,2 DCA Reductase (DCAR)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
PCE Reductase (PCE-1)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
PCE Reductase (PCE-2)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalogenimonas trans-1,2-DCE</i> Reductase (TDR)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalogenimonas cerA</i> Reductase (CER)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04

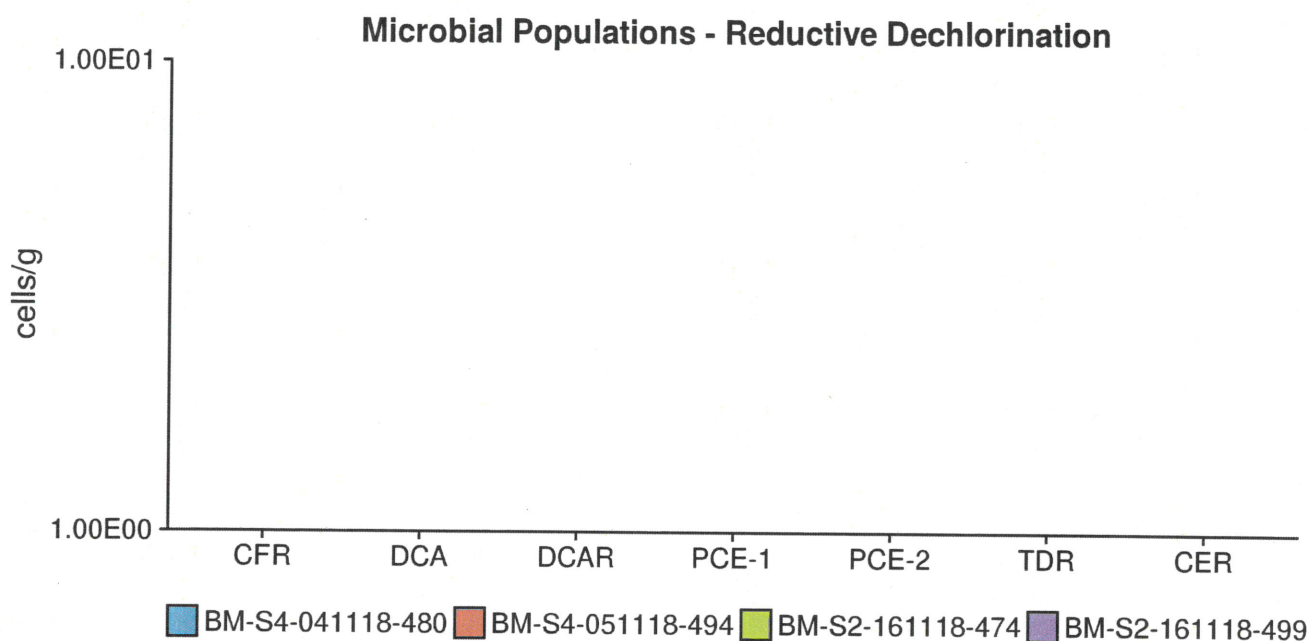


Figure 16: Comparison - microbial populations involved in reductive dechlorination.

Table 8: Summary of the QuantArray®-Chlor results for microorganisms responsible for reductive dechlorination for samples BM-S3-201118-477, BM-S3-201118-489, and BM-S3-211118-494.

Sample Name	BM-S3-201118-477	BM-S3-201118-489	BM-S3-211118-494
Sample Date	11/20/2018	11/20/2018	11/21/2018
Reductive Dechlorination	cells/g	cells/g	cells/g
<i>Dehalococcoides</i> (DHC)	<1.00E+03	<1.00E+03	<1.00E+03
tceA Reductase (TCE)	<1.00E+03	<1.00E+03	<1.00E+03
BAV1 Vinyl Chloride Reductase (BVC)	<1.00E+03	<1.00E+03	<1.00E+03
Vinyl Chloride Reductase (VCR)	<1.00E+03	<1.00E+03	<1.00E+03
<i>Dehalobacter</i> spp. (DHBt)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalobacter</i> DCM (DCM)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalogenimonas</i> spp. (DHG)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Desulfotobacterium</i> spp. (DSB)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalobium chlorocoercia</i> (DECO)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Desulfuromonas</i> spp. (DSM)	<1.00E+04	<1.00E+04	<1.00E+04

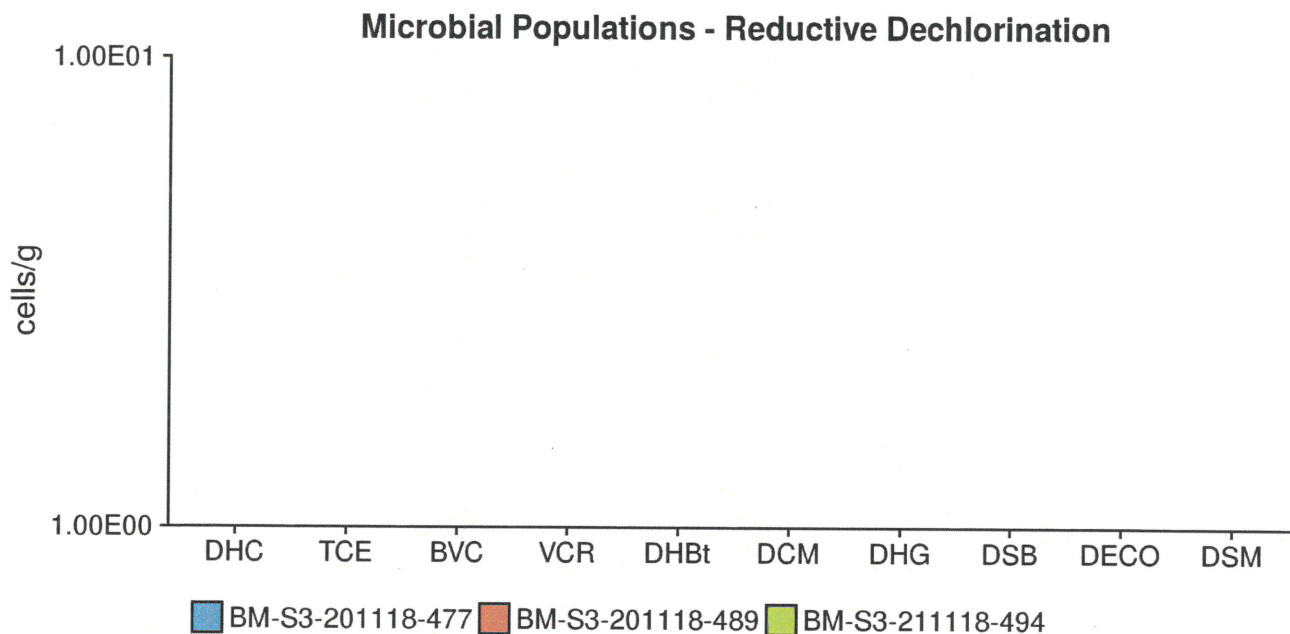


Figure 17: Comparison - microbial populations involved in reductive dechlorination.

Table 9: Summary of the QuantArray®-Chlor results for microorganisms responsible for reductive dechlorination for samples BM-S3-201118-477, BM-S3-201118-489, and BM-S3-211118-494.

Sample Name	BM-S3-201118-477	BM-S3-201118-489	BM-S3-211118-494
Sample Date	11/20/2018	11/20/2018	11/21/2018
Reductive Dechlorination	cells/g	cells/g	cells/g
Chloroform Reductase (CFR)	<1.00E+04	<1.00E+04	<1.00E+04
1,1 DCA Reductase (DCA)	<1.00E+04	<1.00E+04	<1.00E+04
1,2 DCA Reductase (DCAR)	<1.00E+04	<1.00E+04	<1.00E+04
PCE Reductase (PCE-1)	<1.00E+04	<1.00E+04	<1.00E+04
PCE Reductase (PCE-2)	<1.00E+04	<1.00E+04	<1.00E+04
Dehalogenimonas trans-1,2-DCE Reductase (TDR)	<1.00E+04	<1.00E+04	<1.00E+04
Dehalogenimonas cerA Reductase (CER)	<1.00E+04	<1.00E+04	<1.00E+04

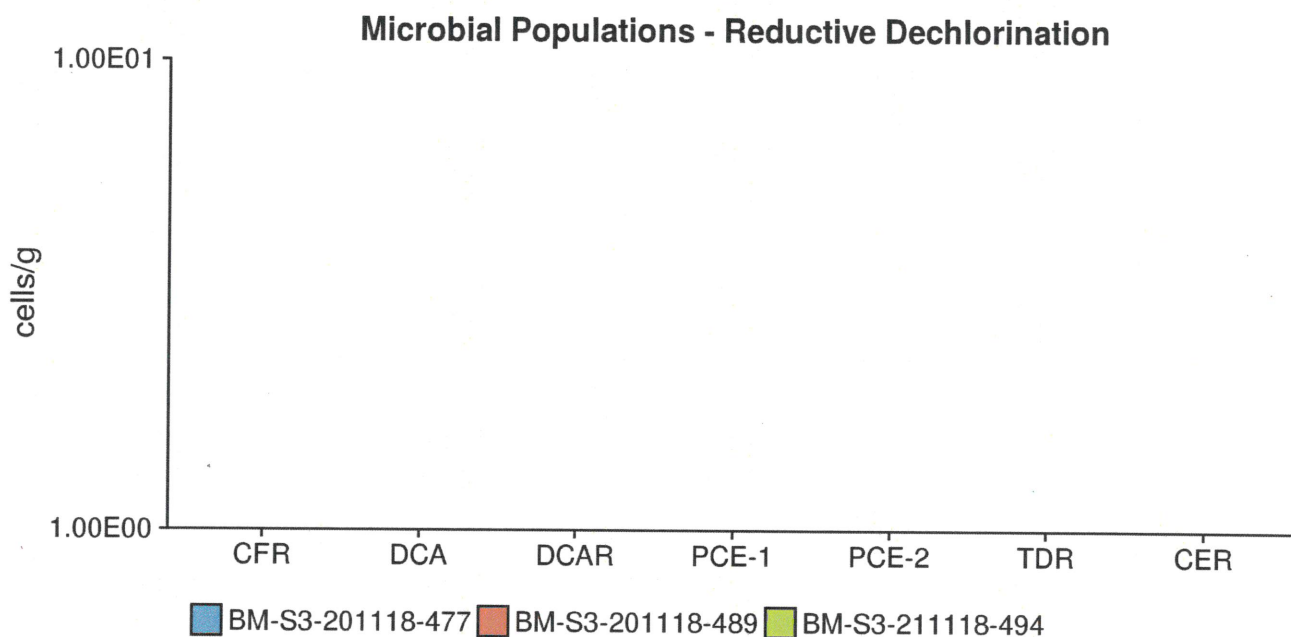


Figure 18: Comparison - microbial populations involved in reductive dechlorination.

Table 10: Summary of the QuantArray®-Chlor results for microorganisms responsible for aerobic (co)metabolism for samples BM-S9-171018-475, BM-S9-181018-477, BM-S5-231018-491, BM-S5-231018-506, and BM-S9-301018-496.

Sample Name	BM-S9-171018-475	BM-S9-181018-477	BM-S5-231018-491	BM-S5-231018-506	BM-S9-301018-496
Sample Date	10/17/2018	10/18/2018	10/23/2018	10/23/2018	10/30/2018
<i>Aerobic (Co)Metabolic</i>	cells/g	cells/g	cells/g	cells/g	cells/g
Soluble Methane Monooxygenase (SMMO)	<1.00E+04	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Dioxygenase (TOD)	<1.00E+04	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Phenol Hydroxylase (PHE)	1.06E+06	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Trichlorobenzene Dioxygenase (TCBO)	<1.00E+04	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Monooxygenase 2 (RDEG)	5.30E+05	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Monooxygenase (RMO)	6.73E+05	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Ethene Monooxygenase (EtnC)	<1.00E+04	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Epoxyalkane Transferase (EtnE)	<1.00E+04	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Dichloromethane Dehalogenase (DCMA)	<1.00E+04	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04

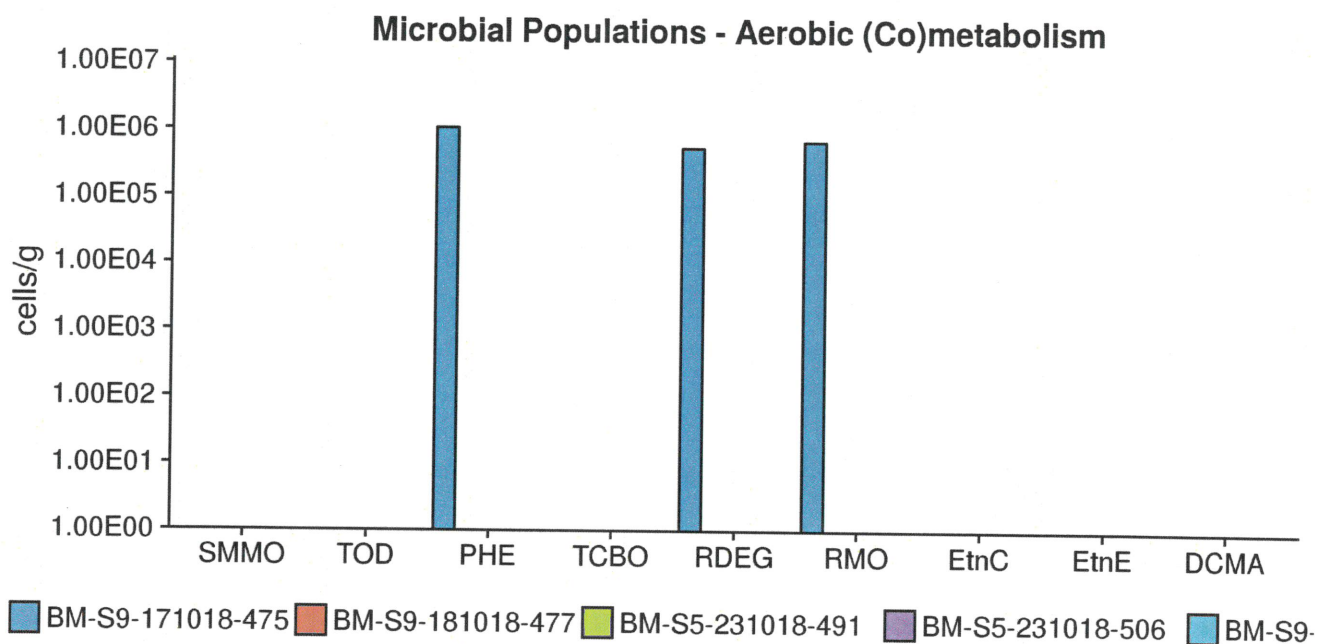


Figure 19: Comparison - microbial populations involved in aerobic (co)metabolism.

Table 11: Summary of the QuantArray®-Chlor results for microorganisms responsible for aerobic (co)metabolism for samples BM-S4-041118-480, BM-S4-051118-494, BM-S2-161118-474, and BM-S2-161118-499.

Sample Name	BM-S4-041118-480	BM-S4-051118-494	BM-S2-161118-474	BM-S2-161118-499
Sample Date	11/04/2018	11/05/2018	11/16/2018	11/16/2018
<i>Aerobic (Co)Metabolic</i>	cells/g	cells/g	cells/g	cells/g
Soluble Methane Monooxygenase (SMMO)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Dioxygenase (TOD)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Phenol Hydroxylase (PHE)	2.42E+03 (J)	<1.00E+04	1.17E+06	<1.00E+04
Trichlorobenzene Dioxygenase (TCBO)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Monooxygenase 2 (RDEG)	1.53E+05	<1.00E+04	1.32E+06	<1.00E+04
Toluene Monooxygenase (RMO)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Ethene Monooxygenase (EtnC)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Epoxyalkane Transferase (EtnE)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Dichloromethane Dehalogenase (DCMA)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04

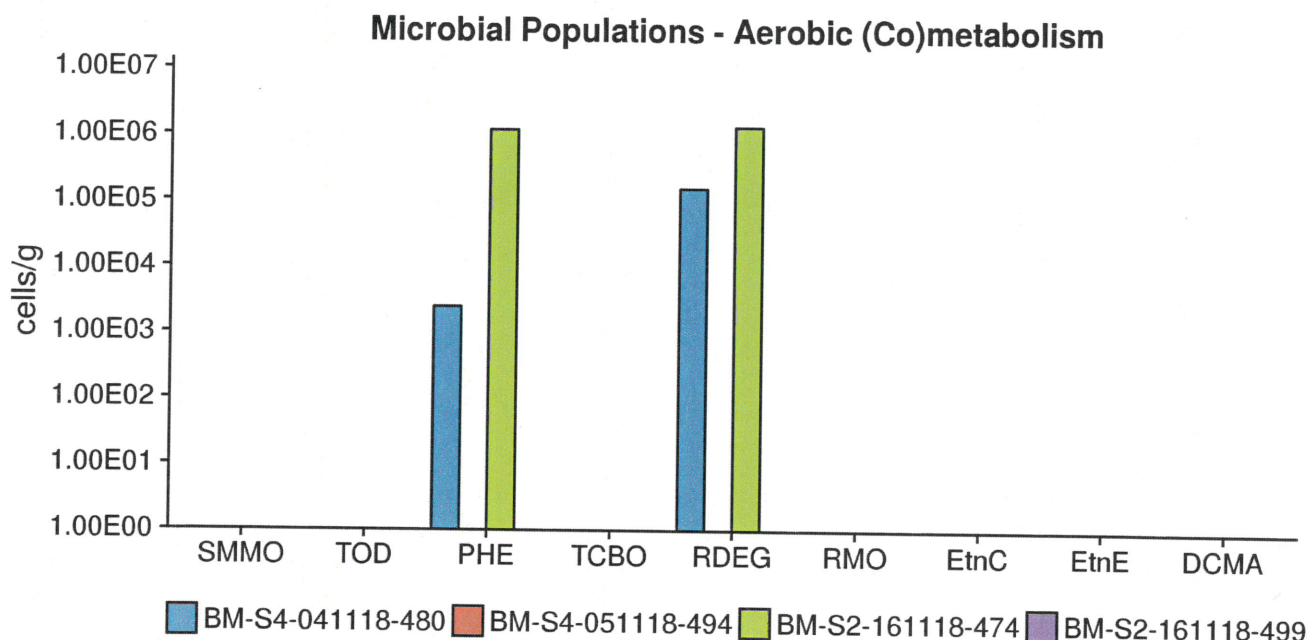


Figure 20: Comparison - microbial populations involved in aerobic (co)metabolism.

Table 12: Summary of the QuantArray®-Chlor results for microorganisms responsible for aerobic (co)metabolism for samples BM-S3-201118-477, BM-S3-201118-489, and BM-S3-211118-494.

Sample Name	BM-S3-201118-477	BM-S3-201118-489	BM-S3-211118-494
Sample Date	11/20/2018	11/20/2018	11/21/2018
Aerobic (Co)Metabolic	cells/g	cells/g	cells/g
Soluble Methane Monooxygenase (SMMO)	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Dioxygenase (TOD)	<1.00E+04	<1.00E+04	<1.00E+04
Phenol Hydroxylase (PHE)	<1.00E+04	<1.00E+04	<1.00E+04
Trichlorobenzene Dioxygenase (TCBO)	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Monooxygenase 2 (RDEG)	9.93E+04	<1.00E+04	<1.00E+04
Toluene Monooxygenase (RMO)	<1.00E+04	<1.00E+04	<1.00E+04
Ethene Monooxygenase (EtnC)	<1.00E+04	<1.00E+04	<1.00E+04
Epoxyalkane Transferase (EtnE)	<1.00E+04	<1.00E+04	<1.00E+04
Dichloromethane Dehalogenase (DCMA)	<1.00E+04	<1.00E+04	<1.00E+04

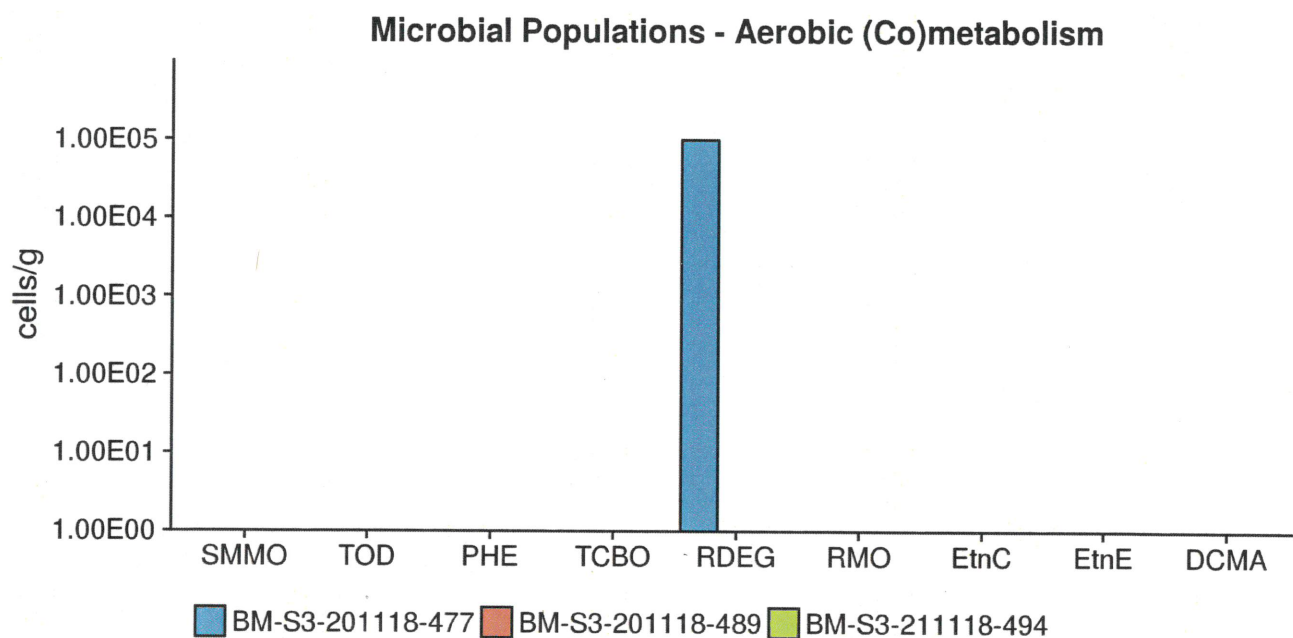


Figure 21: Comparison - microbial populations involved in aerobic (co)metabolism.

Table 13: Summary of the QuantArray® results for total bacteria and other populations for samples BM-S9-171018-475, BM-S9-181018-477, BM-S5-231018-491, BM-S5-231018-506, and BM-S9-301018-496.

Sample Name	BM-S9-171018-475	BM-S9-181018-477	BM-S5-231018-491	BM-S5-231018-506	BM-S9-301018-496
Sample Date	10/17/2018	10/18/2018	10/23/2018	10/23/2018	10/30/2018
Other	cells/g	cells/g	cells/g	cells/g	cells/g
Total Eubacteria (EBAC)	2.69E+07	3.22E+04 (I)	2.86E+05 (I)	3.39E+04 (I)	1.02E+06
Sulfate Reducing Bacteria (APS)	1.08E+07	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Methanogens (MGN)	<1.00E+04	<2.00E+04	<1.00E+04	<1.00E+04	<1.00E+04

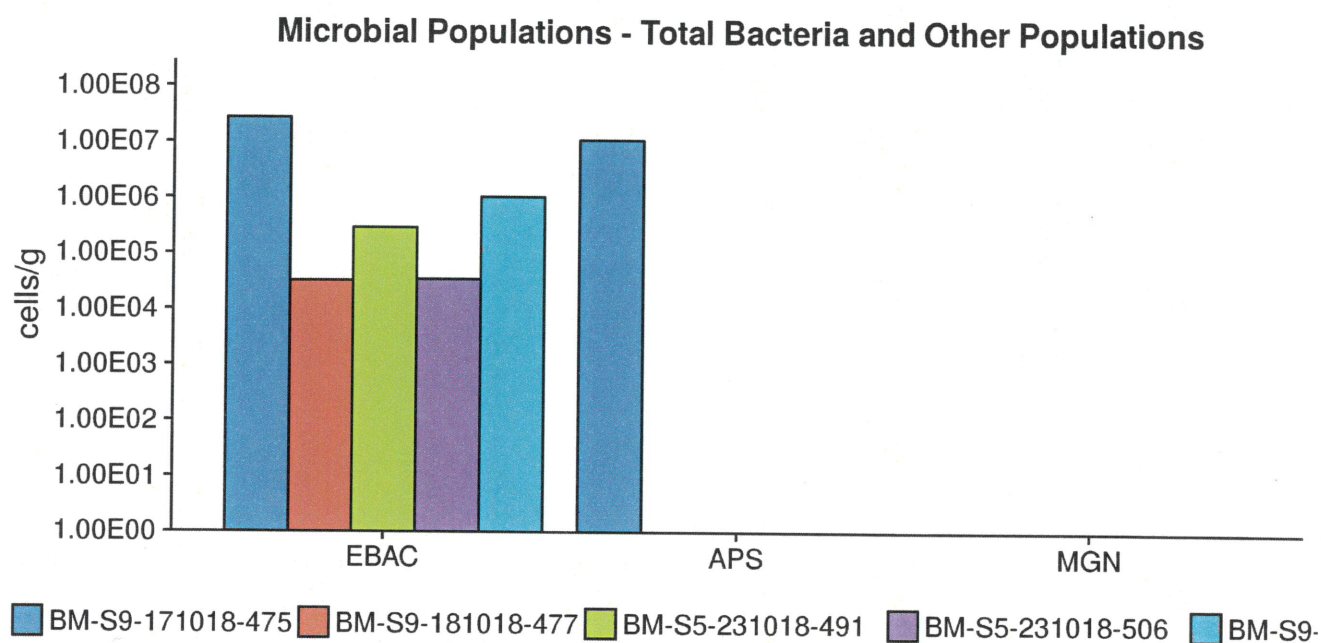


Figure 22: Comparison - microbial populations.

Table 14: Summary of the QuantArray® results for total bacteria and other populations for samples BM-S4-041118-480, BM-S4-051118-494, BM-S2-161118-474, and BM-S2-161118-499.

Sample Name	BM-S4-041118-480	BM-S4-051118-494	BM-S2-161118-474	BM-S2-161118-499
Sample Date	11/04/2018	11/05/2018	11/16/2018	11/16/2018
Other	cells/g	cells/g	cells/g	cells/g
Total Eubacteria (EBAC)	1.16E+06	5.57E+05 (I)	2.43E+06	3.59E+05
Sulfate Reducing Bacteria (APS)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Methanogens (MGN)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04

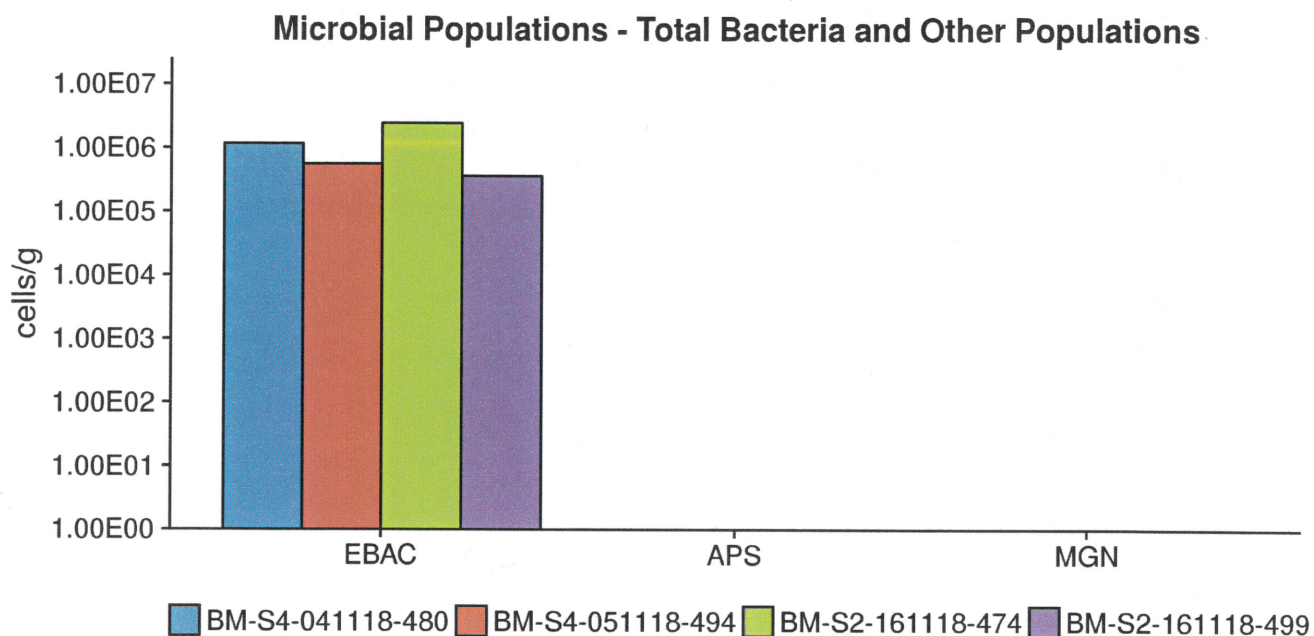


Figure 23: Comparison - microbial populations.

Table 15: Summary of the QuantArray® results for total bacteria and other populations for samples BM-S3-201118-477, BM-S3-201118-489, and BM-S3-211118-494.

Sample Name	BM-S3-201118-477	BM-S3-201118-489	BM-S3-211118-494
Sample Date	11/20/2018	11/20/2018	11/21/2018
Other	cells/g	cells/g	cells/g
Total Eubacteria (EBAC)	5.73E+05	9.78E+03 (I)	6.99E+03 (I)
Sulfate Reducing Bacteria (APS)	<1.00E+04	<1.00E+04	<1.00E+04
Methanogens (MGN)	<1.00E+04	<1.00E+04	<1.00E+04

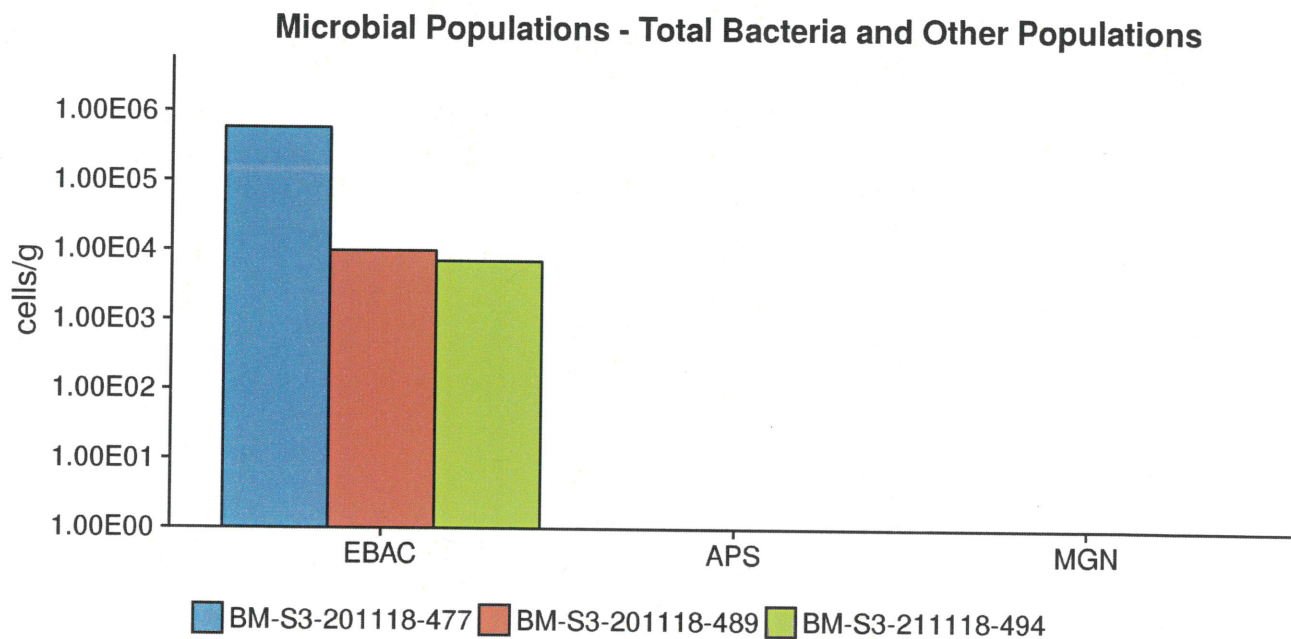


Figure 24: Comparison - microbial populations.

Interpretation

The overall purpose of the QuantArray®-Chlor is to give site managers the ability to simultaneously yet economically evaluate the potential for biodegradation of a spectrum of common chlorinated contaminants through a multitude of anaerobic and aerobic (co)metabolic pathways in order to provide a clearer and more comprehensive view of contaminant biodegradation. The following discussion describes the interpretation of results in general terms and is meant to serve as a guide.

Reductive Dechlorination - Chlorinated Ethenes: While a number of bacterial cultures including *Dehalococcoides*, *Dehalobacter*, *Desulfotobacterium*, and *Desulfuromonas* spp. capable of utilizing PCE and TCE as growth-supporting electron acceptors have been isolated [1–5], *Dehalococcoides* may be the most important because they are the only bacterial group that has been isolated to date which is capable of complete reductive dechlorination of PCE to ethene [6]. In fact, the presence of *Dehalococcoides* has been associated with complete reductive dechlorination to ethene at sites across North America and Europe [7], and Lu et al. [8] have proposed using a *Dehalococcoides* concentration of 1×10^4 cells/mL as a screening criterion to identify sites where biological reductive dechlorination is predicted to proceed at “generally useful” rates.

At chlorinated ethene sites, any “stall” leading to the accumulation of daughter products, especially vinyl chloride, would be a substantial concern. While *Dehalococcoides* concentrations greater than 1×10^4 cells/mL correspond to ethene production and useful rates of dechlorination, the range of chlorinated ethenes degraded varies by strain within the *Dehalococcoides* genus [6, 9], and the presence of co-contaminants and competitors can have complex impacts on the halo-respiring microbial community [10–15]. Therefore, QuantArray®-Chlor also provides quantification of a suite of reductive dehalogenase genes (PCE, TCE, BVC, VCR, CER, and TDR) to more definitively confirm the potential for reductive dechlorination of all chlorinated ethene compounds including vinyl chloride.

Perhaps most importantly, QuantArray®-Chlor quantifies TCE reductase (TCE) and both known vinyl chloride reductase genes (BVC, VCR) from *Dehalococcoides* to conclusively evaluate the potential for complete reductive dechlorination of chlorinated ethenes to non-toxic ethene [16–18]. In addition, the analysis also includes quantification of reductive dehalogenase genes from *Dehalogenimonas* spp. capable of reductive dechlorination of chlorinated ethenes. More specifically, these are the trans-1,2-DCE dehalogenase gene (TDR) from strain WBC-2 [19] and the vinyl chloride reductase gene (CER) from GP, the only known organisms other than *Dehalococcoides* capable of vinyl chloride reduction [20]. Finally, PCE reductase genes responsible for sequential reductive dechlorination of PCE to cis-DCE by *Sulfurospirillum* and *Geobacter* spp. are also quantified. In mixed cultures, evidence increasingly suggests that partial dechlorinators like *Sulfurospirillum* and *Geobacter* may be responsible for the majority of reductive dechlorination of PCE to TCE and cis-DCE while *Dehalococcoides* functions more as cis-DCE and vinyl chloride reducing specialists [10, 21].

Reductive Dechlorination - Chlorinated Ethanes: Under anaerobic conditions, chlorinated ethanes are susceptible to reductive dechlorination by several groups of halo-respiring bacteria including *Dehalobacter*, *Dehalogenimonas*, and *Dehalococcoides*. While the reported range of chlorinated ethanes utilized varies by genus, species, and sometimes at the strain level, several general observations can be made regarding biodegradation pathways and daughter product formation. *Dehalobacter* spp. have been isolated that are capable of sequential reductive dechlorination of 1,1,1-TCA through 1,1-DCA to chloroethane [13]. Biodegradation of 1,1,2-TCA by several halo-respiring bacteria including *Dehalobacter* and *Dehalogenimonas* spp. proceeds via dichloroelimination producing vinyl chloride [22–24]. Similarly, 1,2-DCA biodegradation by *Dehalobacter*, *Dehalogenimonas*, and *Dehalococcoides* occurs via dichloroelimination producing ethene. While not utilized by many *Desulfotobacterium* isolates, at least one strain, *Desulfotobacterium dichloroeliminans* strain DCA1, is also capable of dichloroelimination of 1,2-DCA [25]. The 1,2-dichloroethane reductive dehalogenase gene (DCAR) from members of *Desulfotobacterium* and *Dehalobacter* is known to dechlorinate 1,2-DCA to ethene, while the 1,1-dichloroethane reductive dehalogenase (DCA) targets the gene responsible for 1,1-DCA dechlorination in some strains of *Dehalobacter*. In addition to chloroform, chloroform reductase (CFR) has also been shown to be responsible for reductive dechlorination of 1,1,1-TCA [26].

Reductive Dechlorination - Chlorinated Methanes: Chloroform is a common co-contaminant at chlorinated solvent sites and can inhibit reductive dechlorination of chlorinated ethenes. Grostern et al. demonstrated that a *Dehalobacter* population was capable of reductive dechlorination of chloroform to produce dichloromethane [27]. The *cfrA* gene encodes the reductase which catalyzes this initial step in chloroform biodegradation [26]. Justicia-Leon et al. have since shown that dichloromethane can support growth of a distinct group of *Dehalobacter* strains via fermentation [28]. The *Dehalobacter* DCM assay targets the 16S rRNA gene of these strains.

Reductive Dechlorination - Chlorinated Benzenes: Chlorinated benzenes are an important class of industrial solvents and chemical intermediates in the production of drugs, dyes, herbicides, and insecticides. The physical-chemical properties of chlorinated benzenes as well as susceptibility to biodegradation are functions of their degree of chlorination and the positions of chlorine substituents. Under anaerobic conditions, reductive dechlorination of higher chlorinated benzenes including hexachlorobenzene (HCB),

pentachlorobenzene (PeCB), tetrachlorobenzene (TeCB) isomers, and trichlorobenzene (TCB) isomers has been well documented [29], although biodegradation of individual compounds and isomers varies between isolates. For example, *Dehalococcoides* strain CBDB1 reductively dechlorinates HCB, PeCB, all three TeCB isomers, 1,2,3-TCB, and 1,2,4-TCB [9, 30]. *Dehalobium chlorocoercia* DF-1 has been shown to be capable of reductive dechlorination of HCB, PeCB, and 1,2,3,5-TeCB [31]. The dichlorobenzene (DCB) isomers and chlorobenzene (CB) were considered relatively recalcitrant under anaerobic conditions. However, new evidence has demonstrated reductive dechlorination of DCBs to CB and CB to benzene [32] with corresponding increases in concentrations of *Dehalobacter* spp. [33].

Reductive Dechlorination - Chlorinated Phenols: Pentachlorophenol (PCP) was one of the most widely used biocides in the U.S. and despite residential use restrictions, is still extensively used industrially as a wood preservative. Along with PCP, the tetrachlorophenol and trichlorophenol isomers were also used as fungicides in wood preserving formulations. 2,4-Dichlorophenol and 2,4,5-TCP were used as chemical intermediates in herbicide production (e.g. 2,4-D) and chlorophenols are known byproducts of chlorine bleaching in the pulp and paper industry. While the range of compounds utilized varies by strain, some *Dehalococcoides* isolates are capable of reductive dechlorination of PCP and other chlorinated phenols. For example, *Dehalococcoides* strain CBDB1 is capable of utilizing PCP, all three tetrachlorophenol (TeCP) congeners, all six trichlorophenol (TCP) congeners, and 2,3-dichlorophenol (2,3-DCP). PCP dechlorination by strain CBDB1 produces a mixture of 3,5-DCP, 3,4-DCP, 2,4-DCP, 3-CP, and 4-CP [34]. In the same study, however, *Dehalococcoides* strain 195 dechlorinated a more narrow spectrum of chlorophenols which included 2,3-DCP, 2,3,4-TCP, and 2,3,6-TCP, but no other TCPs or PCP. Similar to *Dehalococcoides*, some species and strains of *Desulfotobacterium* are capable of utilizing PCP and other chlorinated phenols. *Desulfotobacterium hafniense* PCP-1 is capable of reductive dechlorination of PCP to 3-CP [35]. However, the ability to biodegrade PCP is not universal among *Desulfotobacterium* isolates. *Desulfotobacterium* sp. strain PCE1 and *D. chlororespirans* strain Co23, for example, can utilize some TCP and DCP isomers, but not PCP for growth [2, 36].

Reductive Dechlorination - Chlorinated Propanes: *Dehalogenimonas* is a recently described bacterial genus of the phylum Chloroflexi which also includes the well-known chloroethene-respiring *Dehalococcoides* [23]. The *Dehalogenimonas* isolates characterized to date are also halorespiring bacteria, but utilize a rather unique range of chlorinated compounds as electron acceptors including chlorinated propanes (1,2,3-TCP and 1,2-DCP) and a variety of other vicinally chlorinated alkanes including 1,1,2,2-tetrachloroethane, 1,1,2-trichloroethane, and 1,2-dichloroethane [23].

Aerobic - Chlorinated Ethene Cometabolism: Under aerobic conditions, several different types of bacteria including methane-oxidizing bacteria (methanotrophs), and many benzene, toluene, ethylbenzene, xylene, and (BTEX)-utilizing bacteria can cometabolize or co-oxidize TCE, DCE, and vinyl chloride [37]. In general, cometabolism of chlorinated ethenes is mediated by monooxygenase enzymes with "relaxed" specificity that oxidize a primary (growth supporting) substrate (e.g. methane) and co-oxidize the chlorinated compound (e.g. TCE). QuantArray[®]-Chlor provides quantification of a suite of genes encoding oxygenase enzymes capable of co-oxidation of chlorinated ethenes including soluble methane monooxygenase (sMMO). Soluble methane monooxygenases co-oxidize a broad range of chlorinated compounds [38–41] including TCE, *cis*-DCE, and vinyl chloride. Furthermore, soluble methane monooxygenases are generally believed to support greater rates of aerobic cometabolism [40]. QuantArray[®]-Chlor also quantifies aromatic oxygenase genes encoding ring hydroxylating toluene monooxygenase genes (RMO, RDEG), toluene dioxygenase (TOD) and phenol hydroxylases (PHE) capable of TCE co-oxidation [42–46]. TCE or a degradation product has been shown to induce expression of toluene monooxygenases in some laboratory studies [43, 47] raising the possibility of TCE cometabolism with an alternative (non-aromatic) growth substrate. Moreover, while a number of additional factors must be considered, recent research under ESTCP Project 201584 has shown positive correlations between concentrations of monooxygenase genes (soluble methane monooxygenase, ring hydroxylating monooxygenases, and phenol hydroxylase) and the rate of TCE degradation [48].

Aerobic - Chlorinated Ethane Cometabolism: While less widely studied than cometabolism of chlorinated ethenes, some chlorinated ethanes are also susceptible to co-oxidation. As mentioned previously, soluble methane monooxygenases (sMMO) exhibit very relaxed specificity. In laboratory studies, sMMO has been shown to co-oxidize a number of chlorinated ethanes including 1,1,1-TCA and 1,2-DCA [38, 40].

Aerobic - Vinyl Chloride Cometabolism: Beginning in the early 1990s, numerous microcosm studies demonstrated aerobic oxidation of vinyl chloride under MNA conditions without the addition of exogenous primary substrates. Since then, strains of

Mycobacterium, *Nocardioideis*, *Pseudomonas*, *Ochrobactrum*, and *Ralstonia* species have been isolated which are capable of aerobic growth on both ethene and vinyl chloride (see Mattes et al. [49] for a review). The initial steps in the pathway are the monooxygenase (*etnABCD*) catalyzed conversion of ethene and vinyl chloride to their respective epoxyalkanes (epoxyethane and chlorooxirane), followed by epoxyalkane:CoM transferase (*etnE*) mediated conjugation and breaking of the epoxide [50].

Aerobic - Chlorinated Benzenes: In general, chlorobenzenes with four or less chlorine groups are susceptible to aerobic biodegradation and can serve as growth-supporting substrates. Toluene dioxygenase (TOD) has a relatively relaxed substrate specificity and mediates the incorporation of both atoms of oxygen into the aromatic ring of benzene and substituted benzenes (toluene and chlorobenzene). Comparison of TOD levels in background and source zone samples from a CB-impacted site suggested that CBs promoted growth of TOD-containing bacteria [51]. In addition, aerobic biodegradation of some trichlorobenzene and even tetrachlorobenzene isomers is initiated by a group of related trichlorobenzene dioxygenase genes (TCBO). Finally, phenol hydroxylases catalyze the continued oxidation and in some cases, the initial oxidation of a variety of monoaromatic compounds. In an independent study, significant increases in numbers of bacteria containing PHE genes corresponded to increases in biodegradation of DCB isomers [51].

Aerobic - Chlorinated Methanes: Many aerobic methylotrophic bacteria, belonging to diverse genera (*Hyphomicrobium*, *Methylobacterium*, *Methylophilus*, *Pseudomonas*, *Paracoccus*, and *Alibacter*) have been isolated which are capable of utilizing dichloromethane (DCM) as a growth substrate. The DCM metabolic pathway in methylotrophic bacteria is initiated by a dichloromethane dehalogenase (DCMA) gene. DCMA is responsible for aerobic biodegradation of dichloromethane by methylotrophs by first producing formaldehyde which is then further oxidized [52]. As discussed in previous sections, soluble methane monooxygenase (sMMO) exhibits relaxed specificity and co-oxidizes a broad spectrum of chlorinated hydrocarbons. In addition to chlorinated ethenes, sMMO has been shown to co-oxidize chloroform in laboratory studies [38, 41].

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SITE LOGIC Report

Abiotic Potential

Contact: Bernard Bockisch

Phone: 505-234-1105

Address: EA Engineering - Albuquerque, NM
320 Gold Ave SW, Suite 1300
Albuquerque, NM 87102

Email: bbockisch@eaest.com

MI Identifier: 072PJ

Report Date: 11-30-2018

Project Name: KAFB-Bulk Fuels Fac Vadose Zone

Comments:

Overview

Although not always fully considered, abiotic degradation can be a substantial or even the primary process for chlorinated hydrocarbon destruction at sites undergoing or transitioning to monitored natural attenuation (MNA). A variety of iron-bearing minerals including iron sulfides (mackinawite and pyrite), iron oxides (magnetite), green rust, and iron-bearing clays are capable of complete or nearly complete degradation of PCE, TCE, and carbon tetrachloride (He et al. 2009). Some iron-bearing minerals also catalyze the degradation of chlorinated ethanes and the lesser chlorinated ethenes cis-dichloroethene (DCE) and vinyl chloride. While the quantities and types will vary, these reactive iron minerals are frequently identified in subsurface environments under iron reducing and sulfate reducing conditions.

Brown et al. (2007) recommend four avenues for evaluating the role of abiotic processes in contaminant attenuation. First, examining contaminant concentrations along the flow path - decreasing parent compound concentrations with no evidence of accumulation of chlorinated transformation products like cis-DCE and vinyl chloride suggest abiotic degradation. Performing compound specific isotope analysis (CSIA) or monitoring for products unique to abiotic reactions such as acetylene can also provide a strong line of evidence. Microcosm studies with native sediment and killed controls can also be performed. Finally, Brown et al. (2007) suggest performing mineralogical analyses on aquifer sediment to characterize reactive minerals such as magnetite or iron monosulfides

Magnetic Susceptibility - Magnetite

Magnetite (Fe_3O_4) is a mixed valence iron mineral shown to react with PCE, TCE, and carbon tetrachloride. Furthermore, Ferrey et al. (2004) conclusively linked the observed degradation of cis-DCE at a former ammunition plant to magnetite in the subsurface. No direct chemical test is available for quantification of magnetite. However, magnetite is the most abundant mineral in natural sediments that exhibits magnetic behavior. Therefore, magnetic susceptibility provides an inexpensive and valuable estimate of the quantity of magnetite in environmental samples.

X-ray Diffraction (XRD) - Mackinawite, Pyrite, Magnetite and Green Rust

XRD is one of the primary techniques used to identify unknown crystalline materials. Most minerals are crystalline and will scatter X-rays in a regular, characteristic manner dependent on their crystal structure.

- Mackinawite is the most reactive of the iron-bearing minerals and a crystalline form (tetragonal FeS) can be detected by XRD. Mackinawite will transform PCE and TCE primarily by elimination to acetylene. Carbon tetrachloride is transformed mainly to chloroform but carbon dioxide, formate, and carbon disulfide have also been detected. Finally, the more heavily chlorinated ethanes including hexachloroethane, pentachloroethane, and tetrachloroethanes react to form chlorinated ethenes which can be further degraded.
- Pyrite (FeS_2) catalyzes beta elimination transforming PCE, TCE, and cis-DCE to acetylene and ethene. Vinyl chloride is transformed to ethene and ethane. Pyrite is also capable of degradation of carbon tetrachloride potentially forming a number of products including chloroform, carbon dioxide, carbon disulfide, and formate depending on reaction conditions.
- While not quantitative like the magnetic susceptibility test, XRD can also detect magnetite when present at between 2% and 5% on a weight basis.
- Green rusts have been reported to transform a number of common chlorinated contaminants including cis-DCE, vinyl chloride, trichloroethanes, and tetrachloroethanes. While special sample care to prevent oxidation would be needed, XRD can be used to detect green rust.

Percent Clay

Clays have large surface areas, balanced by exchangeable cations, which can bind a large number of both organic and inorganic molecules impacting their availability and reactivity in the subsurface. While less well studied than the other iron-bearing minerals, various phyllosilicate clays have been shown to be capable of degradation of PCE, TCE, cis-DCE, vinyl chloride, and carbon tetrachloride

Sample Location	BM-S9-171018-475	BM-S9-181018-477	BM-S5-231018-491
Date Sampled	10/17/2018	10/18/2018	10/23/2018

Magnetic Susceptibility Analysis

Magnetic Susceptibility (m3/kg)	$3.2\text{E-}6 \pm 6.4\text{E-}8$	$1.3\text{E-}6 \pm 2.2\text{E-}8$	$1.9\text{E-}6 \pm 4.0\text{E-}8$
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Sample Location	BM-S5-231018-506	BM-S9-181018-483	BM-S9-301018-496
Date Sampled	10/23/2018	10/18/2018	10/30/2018

Magnetic Susceptibility Analysis

Magnetic Susceptibility (m3/kg)	$2.4\text{E-}6 \pm 1.1\text{E-}7$	$1.7\text{E-}6 \pm 4.2\text{E-}8$	$2.2\text{E-}6 \pm 8.4\text{E-}9$
---------------------------------	-----------------------------------	-----------------------------------	-----------------------------------

Sample Location	BM-S4-041118-480	BM-S4-051118-494	BM-S2-161118-474
Date Sampled	11/04/2018	11/05/2018	11/16/2018

Magnetic Susceptibility Analysis

Magnetic Susceptibility (m3/kg)	$1.9\text{E-}6 \pm 8.4\text{E-}8$	$2.0\text{E-}6 \pm 2.6\text{E-}8$	$6.0\text{E-}6 \pm 9.7\text{E-}8$
---------------------------------	-----------------------------------	-----------------------------------	-----------------------------------

Sample Location	BM-S2-161118-499	BM-S3-201118-477	BM-S3-201118-489
Date Sampled	11/16/2018	11/20/2018	11/20/2018

Magnetic Susceptibility Analysis

Magnetic Susceptibility (m3/kg)	$2.6\text{E-}6 \pm 9.3\text{E-}8$	$3.1\text{E-}6 \pm 3.3\text{E-}8$	$2.0\text{E-}6 \pm 2.4\text{E-}7$
---------------------------------	-----------------------------------	-----------------------------------	-----------------------------------

* Analysis performed in triplicate and results reported as the mean followed by +/- standard deviation.

Sample Location

BM-S3-211118-494

Date Sampled

11/21/2018

Magnetic Susceptibility Analysis

Magnetic Susceptibility (m3/kg)

$2.5\text{E-}6 \pm 1.9\text{E-}7$

* Analysis performed in triplicate and results reported as the mean followed by +/- standard deviation.

References

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Abiotic Reactions of Chlorinated Compounds with Iron Bearing Minerals and Zero Valent Iron (ZVI). Summaries for iron bearing minerals are based on He et al. (2009) and references therein. He et al. available at <http://nepis.epa.gov/>. Summary of ZVI based on Liu et al. (2005) and Song et al. (2005).

Contaminant	Mineral	Degradation	Reported Degradation Intermediates and Products ¹
PCE	FeS	Yes	Acetylene, TCE, cis-DCE, 1,1-DCE, ethene
	Pyrite	Yes	TCE, acetylene, ethene
	Magnetite	Yes	Unknown ²
	³ GR(SO ₄) phyllosilicate clays	Reports Differ	
TCE	ZVI	Yes	TCE, 1,1-DCE, vinyl chloride, 1,1,2-TCA, 1,1-DCA, chloroacetylene, acetylene, ethene, ethane
	FeS	Yes	Ethene and ethane
	Pyrite	Yes	Acetylene, cis-DCE, vinyl chloride, 1,1-DCE
	Magnetite	Yes	Acetylene, ethene, cis-DCE, (organic acids with DO present)
cis-DCE	GR(SO ₄), GR(CO ₃)	No	Unknown ¹
	phyllosilicate clays	Yes	Only observed degradation when Cu(II) added
	ZVI	Yes	cDCE, vinyl chloride, acetylene, ethene, ethane
	FeS	Yes	Ethane, ethene, acetylene with minor amounts of DCE, VC depending on conditions
	Pyrite	No	None detected
	Magnetite	Yes	Acetylene, ethene
Vinyl chloride	GR(SO ₄)	Yes	Unknown ²
	phyllosilicate clays	Yes	
	ZVI	Yes	Primarily acetylene and ethene but also much lesser amounts of ethane and VC and traces of methane, propane, propene, butane and butene
	FeS	Unknown	
	Pyrite	Yes	Ethene, ethane
	Magnetite	Yes	Unknown ²
Vinyl chloride	GR(SO ₄)	Yes	
	phyllosilicate clays	Yes	
	ZVI	Yes	Ethene, ethane, (no evidence of acetylene)

Contaminant	Mineral	Degradation	Reported Degradation Intermediates and Products ¹
1,1-DCA	FeS	Not Significant	None detected
1,1-DCA	GR(SO ₄)	Low conversion	Ethene and ethane (w/ Cu or Ag)
1,1-DCA	ZVI	Yes (low)	Ethane
1,2-DCA	FeS	Not Significant	None detected
1,2-DCA	FeS (Biogenic)	Yes	Not monitored
1,2-DCA	GR(SO ₄)	No	
1,2-DCA	ZVI	No	
1,1,1-TCA	FeS	Yes	1,1-DCA, ethene, 2-butyne
1,1,1-TCA	GR(SO ₄)	Yes	1,1-DCA, CA, ethene ethane
1,1,1-TCA	ZVI	Yes	1,1-DCA, ethane
1,1,2-TCA	FeS	Rate not significant	Small amounts of 1,1-DCE and vinyl chloride but rate not significant
1,1,2-TCA	GR(SO ₄)	Yes	Vinyl chloride, 1,1-DCE, ethene, ethane
1,1,2-TCA	ZVI	Yes	Ethane
1,1,1,2-TeCA	FeS	Yes	1,1-DCE
1,1,1,2-TeCA	GR(SO ₄)	Yes	1,1-DCE and minor (<1%) vinyl chloride, ethene, ethane
1,1,1,2-TeCA	phyllosilicate clays	Yes	1,1-DCE
1,1,1,2-TeCA	ZVI	Yes	TCE, 1,1-DCE
1,1,2,2-TeCA	FeS	Yes	TCE, cis-DCE, trans-DCE, acetylene
1,1,2,2-TeCA	GR(SO ₄)	Yes	TCE (major), cis-DCE, trans-DCE
1,1,2,2-TeCA	phyllosilicate clays	Yes	TCE
1,1,2,2-TeCA	ZVI	Yes	TCE, trans-DCE, cis-DCE
Carbon Tetrachloride	FeS	Yes	Chloroform, carbon disulfide, possibly methane, ethene, ethane
CT	Pyrite	Yes	Chloroform, CO ₂ , carbon disulfide, formate (highly dependent on conditions)
CT	Magnetite	Yes	Chloroform, carbon monoxide, methane, formate (highly dependent on conditions)
CT	GR(SO ₄)	Yes	Chloroform and hexachloroethane; Chloroform, DCM, methane, ethene
CT	phyllosilicate clays	Yes	Chloroform
CT	ZVI	Yes	Chloroform, dichloromethane, methane (depending on conditions)

Notes: GR(SO₄) sulfate green rust. GR(CO₃) carbonate green rust. ZVI zero valent iron

¹Compilation of reported degradation products. Mass recovery of products typically low --- additional undetected and unreported products are likely. Reported reaction products or proportions of reaction products were often a function of environmental conditions.

²No published studies that identify the transformation products of PCE, TCE, cis-DCE or vinyl chloride with magnetite. Ferrey et al (2004) analyzed for products of cis-DCE dechlorination including vinyl chloride, ethene, and ethane and did not find them. If Fe²⁺ sorbed to magnetite stabilizes carbene ions, the ultimate degradation product of cis-DCE on magnetite would be CO₂.

SITE LOGIC Report

Abiotic Potential

Contact: Bernard Bockisch
Address: EA Engineering
320 Gold Ave SW
Suite 1300
Albuquerque, NM 87102

Phone: 505-234-1105

Email: bbockisch@eaest.com

MI
Identifier: 072PJ

Report Date: January 11, 2019

Comments: KAFB-Bulk Fuels Fac Vadose Zone, 62735DM02

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Results

Table 1. Summary of X-Ray Diffraction and X-ray fluorescence results.

Sample Location	BM-S9-171018-475	BM-S9-181018-477	BM-S5-231018-491	BM-S5-231018-506	BM-S9-181018-483
Sample Type	Solid	Solid	Solid	Solid	Solid
Sample Date	10/17/2018	10/18/2018	10/23/2018	10/23/2018	10/18/2018
MI ID	072PJ1	072PJ2	072PJ3	072PJ4	072PJ5
X-Ray Diffraction Results					
Mineral Constituent	Relative Abundance (%)				
Quartz	44	27	28	40	31
Cristobalite	ND	ND	ND	ND	ND
Plagioclase Feldspar	39	23	39	40	16
Microcline	6	4	8	7	4
Calcite	ND	5	1		44
Magnetite	5	6	5	5	3.5
Hematite	<0.5	ND	ND	ND	ND
Hornblende	<0.5	ND	ND	ND	ND
Clinoptilolite	ND	0.5	ND	<0.5	ND
Goethite	ND	ND	1	ND	ND
Kaolinite	0.5	1	<0.5	0.5	<0.5
Chlorite	<0.5	0.5	ND	0.5	<0.5
Illite/Mica	0.5	1	0.5	1	<0.5
Montmorillonite	5	32	17.5	6	1.5
Total	100	100	100	100	100
X-Ray Fluorescence Results					
Compound	Results (mass %)				
MgO	0.4032	0.7938	0.5018	0.2765	0.2179
Al ₂ O ₃	8.4816	10.4114	8.7385	8.3363	5.1776
SiO ₂	74.8852	67.2286	74.2316	75.9823	50.5207
P ₂ O ₅	1.0978	0.9791	0.9459	0.9571	0.7796
SO ₃	ND	0.1093	ND	0.0392	ND
Cl	ND	ND	0.0288	ND	ND
K ₂ O	4.5492	5.1831	5.1344	5.2099	3.6163
CaO	3.7306	6.7662	3.4171	3.1390	34.966
TiO ₂	0.5864	0.9050	0.5983	0.5586	0.3818
MnO	0.0863	0.1278	0.1063	0.0832	0.0795
Fe ₂ O ₃	6.0153	6.8537	5.9017	5.2802	4.1215
CuO	ND	ND	ND	ND	ND
ZnO	ND	0.0288	ND	ND	ND
Rb ₂ O	0.0183	0.0326	0.0264	0.0243	0.0184
SrO	0.1089	0.1206	0.0974	0.0831	0.0859
Y ₂ O ₃	0.0034	0.0002	ND	ND	ND
ZrO ₂	0.0339	0.0793	0.0380	0.0303	0.0347
BaO	ND	0.3805	0.2338	ND	ND

Legend: J = Estimated result below PQL but above LQL ND = Result not detected < = Below detection limit

Table 2. Summary of X-Ray Diffraction and X-ray fluorescence results continued.

Sample Location	BM-S9-301018-496	BM-S4-041118-480	BM-S4-051118-494	BM-S2-161118-474
Sample Type	Solid	Solid	Solid	Solid
Sample Date	10/30/2018	11/04/2018	11/05/2018	11/16/2018
MI ID	072PJ6	072PJ7	072PJ8	072PJ9
X-Ray Diffraction Results				
Mineral Constituent	Relative Abundance (%)			
Quartz	34	36	44	35
Cristobalite	ND	ND	ND	ND
Plagioclase Feldspar	48	21	32	33
Microcline	9	5	8	6
Calcite	ND	ND	ND	ND
Magnetite	4	5	4	7
Hematite	ND	ND	ND	1
Hornblende	ND	ND	ND	ND
Clinoptilolite	ND	ND	ND	ND
Goethite	ND	ND	ND	ND
Kaolinite	0.5	1	1	1
Chlorite	ND	ND	ND	ND
Illite/Mica	0.5	1	1	1
Montmorillonite	4	31	10	16
TOTAL	100	100	100	100
X-Ray Fluorescence Results				
Compound	Results (mass %)			
MgO	0.2769	0.6241	0.3689	0.2856
Al ₂ O ₃	8.1006	10.2034	8.8896	8.9350
SiO ₂	72.7651	72.4060	75.3079	70.905
P ₂ O ₅	1.0564	1.0487	1.0403	0.9873
SO ₃	ND	ND	0.0371	0.0485
Cl	ND	ND	ND	0.0154
K ₂ O	5.5193	4.7968	5.6238	4.7379
CaO	6.1292	2.9698	3.2911	3.5409
TiO ₂	0.6463	0.7515	0.5222	1.3590
MnO	ND	0.1901	0.0834	0.1178
Fe ₂ O ₃	5.1149	6.8266	4.5973	8.8067
CuO	ND	ND	ND	ND
ZnO	ND	0.0202	ND	0.0267
Rb ₂ O	0.0243	0.0273	0.0234	0.0291
SrO	0.1142	0.0701	0.0975	0.1062
Y ₂ O ₃	ND	ND	0.0564	ND
ZrO ₂	0.0222	0.0653	0.0612	0.0989
BaO	0.2305	ND	ND	ND

Legend: J = Estimated result below PQL but above LQL ND = Result not detected < = Below detection limit

Table 3. Summary of X-Ray Diffraction and X-ray fluorescence results continued.

Sample Location	BM-S2-161118-499	BM-S3-201118-477	BM-S3-201118-489	BM-S3-211118-494
Sample Type	Solid	Solid	Solid	Solid
Sample Date	11/16/2018	11/20/2018	11/20/2018	11/21/2018
MI ID	072PJ10	072PJ11	072PJ12	072PJ13
X-Ray Diffraction Results				
Mineral Constituent	Relative Abundance (%)			
Quartz	46	34	32	29
Cristobalite	ND	2	3	ND
Plagioclase Feldspar	32	33	41	43
Microcline	8	7	11	10
Calcite	ND	ND	ND	1
Magnetite	3.5	4	5	5
Hematite	0.5	1	ND	ND
Hornblende	ND	ND	ND	ND
Clinoptilolite	<0.5	ND	ND	0.5
Goethite	ND	ND	ND	ND
Kaolinite	0.5	0.5	0.5	0.5
Chlorite	ND	<0.5	ND	ND
Illite/Mica	1	1.5	0.5	1
Montmorillonite	8.5	17	7	10
TOTAL	100	100	100	100
X-Ray Fluorescence Results				
Compound	Results (mass %)			
MgO	ND	0.2307	0.3615	0.1955
Al ₂ O ₃	7.941	8.8761	9.4752	9.9017
SiO ₂	77.2565	75.9737	73.5919	72.5321
P ₂ O ₅	1.005	0.9390	0.9608	1.0373
SO ₃	ND	0.0339	0.0590	ND
Cl	0.0159	0.0412	0.0108	ND
K ₂ O	5.6548	4.9802	5.4124	5.6524
CaO	2.8252	3.1696	3.6045	3.6527
TiO ₂	0.5312	0.569	0.6124	0.6062
MnO	0.1002	0.066	0.1083	0.1282
Fe ₂ O ₃	4.5199	4.952	5.6333	5.8372
CuO	ND	ND	ND	0.0253
ZnO	ND	ND	ND	0.0220
Rb ₂ O	0.0219	0.0205	0.0210	0.0264
SrO	0.0961	0.0900	0.0948	0.0927
Y ₂ O ₃	ND	ND	0.0027	ND
ZrO ₂	0.0323	0.0581	0.0512	0.0315
BaO	ND	ND	ND	0.2587

Legend: J = Estimated result below PQL but above LQL ND = Result not detected < = Below detection limit

References

- Brown, R. A., J. T. Wilson and M. Ferrey (2007). "Monitored natural attenuation forum: The case for abiotic MNA." Remediation Journal **17**(2): 127-137.
- Ferrey, M. L., R. T. Wilkin, R. G. Ford and J. T. Wilson (2004). "Nonbiological Removal of cis-Dichloroethylene and 1,1-Dichloroethylene in Aquifer Sediment Containing Magnetite." Environmental Science & Technology **38**(6): 1746-1752.
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- Liu, Y., S. A. Majetich, R. D. Tilton, D. S. Sholl and G. V. Lowry (2005). "TCE Dechlorination Rates, Pathways, and Efficiency of Nanoscale Iron Particles with Different Properties." Environmental Science & Technology **39**(5): 1338-1345.
- Song, H. and E. R. Carraway (2005). "Reduction of Chlorinated Ethanes by Nanosized Zero-Valent Iron: Kinetics, Pathways, and Effects of Reaction Conditions." Environmental Science & Technology **39**(16): 6237-6245.

Abiotic Reactions of Chlorinated Compounds with Iron Bearing Minerals and Zero Valent Iron (ZVI). Summaries for iron bearing minerals are based on He et al. (2009) and references therein. He et al. available at <http://nepis.epa.gov/>. Summary of ZVI based on Liu et al. (2005) and Song et al. (2005).

Contaminant	Mineral	Degradation	Reported Degradation Intermediates and Products ¹
PCE	FeS	Yes	Acetylene, TCE, cis-DCE, 1,1-DCE, ethene
	Pyrite	Yes	TCE, acetylene, ethene
	Magnetite	Yes	Unknown ²
	³ GR(SO ₄)	Reports Differ	
	phyllosilicate clays	Yes	TCE, 1,1-DCE, vinyl chloride, 1,1,2-TCA, 1,1-DCA, chloroacetylene, acetylene, ethene, ethane
	ZVI	Yes	Ethene and ethane
TCE	FeS	Yes	Acetylene, cis-DCE, vinyl chloride, 1,1-DCE
	Pyrite	Yes	Acetylene, ethene, cis-DCE, (organic acids with DO present)
	Magnetite	Yes	Unknown ¹
	GR(SO ₄), GR(CO ₃)	No	Only observed degradation when Cu(II) added
	phyllosilicate clays	Yes	cDCE, vinyl chloride, acetylene, ethene, ethane
	ZVI	Yes	Ethane, ethene, acetylene with minor amounts of DCE, VC depending on conditions
cis-DCE	FeS	No	None detected
	Pyrite	Yes	Acetylene, ethene
	Magnetite	Yes	Unknown ²
	GR(SO ₄)	Yes	
	phyllosilicate clays	Yes	
	ZVI	Yes	Primarily acetylene and ethene but also much lesser amounts of ethane and VC and traces of methane, propane, propene, butane and butene
Vinyl chloride	FeS	Unknown	
	Pyrite	Yes	Ethene, ethane
	Magnetite	Yes	Unknown ²
	GR(SO ₄)	Yes	
	phyllosilicate clays	Yes	
	ZVI	Yes	Ethene, ethane, (no evidence of acetylene)

Contaminant	Mineral	Degradation	Reported Degradation Intermediates and Products ¹
1,1-DCA	FeS	Not Significant	None detected
1,1-DCA	GR(SO ₄)	Low conversion	Ethene and ethane (w/ Cu or Ag)
1,1-DCA	ZVI	Yes (low)	Ethane
1,2-DCA	FeS	Not Significant	None detected
1,2-DCA	FeS (Biogenic)	Yes	Not monitored
1,2-DCA	GR(SO ₄)	No	
1,2-DCA	ZVI	No	
1,1,1-TCA	FeS	Yes	1,1-DCA, ethene, 2-butyne
1,1,1-TCA	GR(SO ₄)	Yes	1,1-DCA, CA, ethene ethane
1,1,1-TCA	ZVI	Yes	1,1-DCA, ethane
1,1,2-TCA	FeS	Rate not significant	Small amounts of 1,1-DCE and vinyl chloride but rate not significant
1,1,2-TCA	GR(SO ₄)	Yes	Vinyl chloride, 1,1-DCE, ethene, ethane
1,1,2-TCA	ZVI	Yes	Ethane
1,1,1,2-TeCA	FeS	Yes	1,1-DCE
1,1,1,2-TeCA	GR(SO ₄)	Yes	1,1-DCE and minor (<1%) vinyl chloride, ethene, ethane
1,1,1,2-TeCA	phyllosilicate clays	Yes	1,1-DCE
1,1,1,2-TeCA	ZVI	Yes	TCE, 1,1-DCE
1,1,2,2-TeCA	FeS	Yes	TCE, cis-DCE, trans-DCE, acetylene
1,1,2,2-TeCA	GR(SO ₄)	Yes	TCE (major), cis-DCE, trans-DCE
1,1,2,2-TeCA	phyllosilicate clays	Yes	TCE
1,1,2,2-TeCA	ZVI	Yes	TCE, trans-DCE, cis-DCE
Carbon Tetrachloride	FeS	Yes	Chloroform, carbon disulfide, possibly methane, ethene, ethane
CT	Pyrite	Yes	Chloroform, CO ₂ , carbon disulfide, formate (highly dependent on conditions)
CT	Magnetite	Yes	Chloroform, carbon monoxide, methane, formate (highly dependent on conditions)
CT	GR(SO ₄)	Yes	Chloroform and hexachloroethane; Chloroform, DCM, methane, ethene
CT	phyllosilicate clays	Yes	Chloroform
CT	ZVI	Yes	Chloroform, dichloromethane, methane (depending on conditions)

Notes: GR(SO₄) sulfate green rust. GR(CO₃) carbonate green rust. ZVI zero valent iron

¹Compilation of reported degradation products. Mass recovery of products typically low - additional undetected and unreported products are likely. Reported reaction products or proportions of reaction products were often a function of environmental conditions.

²No published studies that identify the transformation products of PCE, TCE, cis-DCE or vinyl chloride with magnetite. Ferrey et al (2004) analyzed for products of cis-DCE dechlorination including vinyl chloride, ethene, and ethane and did not find them. If Fe²⁺ sorbed to magnetite stabilizes carbene ions, the ultimate degradation product of cis-DCE on magnetite would be CO₂.

REPORT TO:

Name: P. Moss
Company: EA Engineering, Science & Technology, Inc.
Address: 7995 E Prentice Ave.
Greenwood Village CO 80111
Email: p.moss@eaeast.com
Phone: 303-590-9143
Fax:

Project Manager: Bernie Bockisch
Project Name: Kirtland Vadose Zone Coring
Project No.: 62735DM02.1017

Report Type: ☒ Standard (default) ☐ Microbial Insights Level III raw data(15% surcharge) ☐ Microbial Insights Level IV (25% surcharge) ☐ Comprehensive Interpretive(15%) ☐ Historical Interpretive (35%)

EDD type: ☒ Microbial Insights Standard (default) ☐ All other available EDDs (5% surcharge) Specify EDD Type:

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 Phone: _____
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Purchase Order No. 18091
Subcontract No.
RFI Quote No. Q2018628.0001

Please Check One:

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Sample Information				Analyses		CENSUS: Please select the target organism/gene								
MI ID (Laboratory Use Only)	Sample Name	Date Sampled	Time Sampled	Matrix	Total Number of Containers	PLFA	NGS	QuantArray Chior	QuantArray Petro	DHC (Dehalococcoides)	Magnetic suscept/LXR/RD/ERD	Other	Other	Other
072015	3 BM-231018-489	10/23/18	1331	S	1	1	1	X	1	1	X			
	4 BM-231018-506	10/23/18	1645	S	1	1	1	X	1	1	X			
CONTACT ABOVE per above														

Relinquished by: Kevin M. Kopp

Date: 10/23/18 1800

Received by: [Signature]

Date: 10/24/18 10:57

It is vital that chain of custody is filled out correctly & that all relative information is provided. Failure to provide sufficient and/or correct information regarding reporting, invoicing & analyses requested information may result in delays for which MI will not be liable.

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Knoxville, TN 37932
865-573-8188
www.microbe.com

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REPORT TO:

Name: P. Moss
Company: EA Engineering, Science & Technology, Inc.
Address: 7995 E Prentic Ave.
Greenwood Village CO 80111

Email: p.moss@eaest.com
Phone: 303-590-9143
Fax:

Project Manager:	Bernie Bockisch
Project Name:	Kirtland Vadose Zone Coring
Project No.:	62735DM02.1017

Report Type: ☒ Standard (default) ☐ Microbial Insights Level III raw data (15% surcharge) ☐ Microbial Insights Level IV (25% surcharge) ☐ Comprehensive Interpretive (15%) ☐ Historical Interpretive (35%)

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COC1-S2-MJ-20181116

REPORT TO:

Name: P. Moss
Company: EA Engineering, Science & Technology, Inc.
Address: 7995 E Prentic Ave.
Greenwood Village CO 80111
E-mail: p.moss@eaest.com
Phone: 303-590-9143
Fax:

Project Manager:	Bernie Bockisch
Project Name:	Kirtland Vadose Zone Coring
Project No.:	62735DM02.1017

Report Type:

EDD type: ☒ Microbial Insights Standard (default) ☐ All o

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INVOICE TO: (For Invoices paid by a third party it is imperative that all information be provided)

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 Company: _____
 Address: _____

Dallas Accounts Payable via Email below
EA Engineering, Science & Technology, Inc.

E-mail: dallasap@east.com
Phone: _____
Fax: _____

Purchase Order No.	18091
Subcontract No.	
MI Quote No.	Q2018628.0001

Please Check One:

☐ More samples to follow

☐ No Additional Samples

Report Type: ☒ Standard (default) ☐ Microbial Insights Level III raw data (15% surcharge) ☐ Microbial Insights Level IV (25% surcharge) ☐ Comprehensive Interpretive (15%) ☐ Historical Interpretive (35%)

EDD type: ☒ Microbial Insights Standard (default) ☐ All other available EDDs (5% surcharge) Specify EDD Type:

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02118102-Inv-35-1707

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Name: _____
 Company: Dallas Accounts Payable via Email below
 Address: EA Engineering, Science & Technology, Inc.

 Email: dallasap@eaeast.com
 Phone: _____
 Fax: _____

Purchase Order No.	18091
Subcontract No.	
MI Quote No.	Q2018628.0001

Please Check One:

☐ More samples to follow

☐ No Additional Samples

Report Type: ☒ Standard (default) ☐ Microbial Insights Level III raw data (15% surcharge) ☐ Microbial Insights Level IV (25% surcharge) ☐ Comprehensive Interpretive (15%) ☐ Historical Interpretive (35%)

EDD type: ☒ Microbial Insights Standard (default) ☐ All other available EDDs (5% surcharge) Specify EDD Type:

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SITE LOGIC Report

QuantArray®-Chlor Study

Contact: Pamela J. Moss

Phone: (303) 590-9143

Address: EA Engineering- Greenwood
Village, CO
7995 E. Prentice Ave
Suite 206E
Greenwood Village, CO 80111

Email: pmoss@eaest.com

MI Identifier: 053QA

Report Date: 02/21/2019

Project: KAFB-Bulk Fuels Fac Vadose Zone, 627350DM02
Comments: Please note the total bacterial biomass on all samples was low, and the samples may have PCR inhibition.

NOTICE: This report is intended only for the addressee shown above and may contain confidential or privileged information. If the recipient of this material is not the intended recipient or if you have received this in error, please notify Microbial Insights, Inc. immediately. The data and other information in this report represent only the sample(s) analyzed and are rendered upon condition that it is not to be reproduced without approval from Microbial Insights, Inc. Thank you for your cooperation.

The QuantArray®-Chlor Approach

Quantification of *Dehalococcoides*, the only known bacterial group capable of complete reductive dechlorination of PCE and TCE to ethene, has become an indispensable component of assessment, remedy selection, and performance monitoring at sites impacted by chlorinated solvents. While undeniably a key group of halo-respiring bacteria, *Dehalococcoides* are not the only bacteria of interest in the subsurface because reductive dechlorination is not the only potential biodegradation pathway operative at contaminated sites, and chlorinated ethenes are not always the primary contaminants of concern. The QuantArray®-Chlor not only includes a variety of halo-respiring bacteria (*Dehalococcoides*, *Dehalobacter*, *Dehalogenimonas*, etc.) to assess the potential for reductive dechlorination of chloroethenes, chloroethanes, chlorobenzenes, chlorophenols, and chloroform, but also provides quantification of functional genes involved in aerobic (co)metabolic pathways for biodegradation of chlorinated solvents and even competing biological processes. Thus, the QuantArray®-Chlor will give site managers the ability to simultaneously yet economically evaluate the potential for biodegradation of a spectrum of common chlorinated contaminants through a multitude of anaerobic and aerobic (co) metabolic pathways to give a much more clear and comprehensive view of contaminant biodegradation.

The QuantArray®-Chlor is used to quantify specific microorganisms and functional genes to evaluate the following:

Anaerobic Reductive Dechlorination

Quantification of important halo-respiring bacteria (e.g. *Dehalococcoides*, *Dehalobacter*, *Dehalogenimonas*, *Desulfotobacterium* spp.) and key functional genes (e.g. vinyl chloride reductases, TCE reductase, chloroform reductase) responsible for reductive dechlorination of a broad spectrum of chlorinated solvents.

Aerobic Cometabolism

Several different types of bacteria including methanotrophs and some toluene/phenol utilizing bacteria can co-oxidize TCE, DCE, and vinyl chloride. The QuantArray®-Chlor quantifies functional genes like soluble methane monooxygenase encoding enzymes capable of co-oxidation of chlorinated ethenes.

Aerobic (Co)metabolism of Vinyl Chloride

Ethene oxidizing bacteria are capable of cometabolism of vinyl chloride. In some cases, ethenotrophs can also utilize vinyl chloride as a growth supporting substrate. The QuantArray®-Chlor targets key functional genes in ethene metabolism.

How do QuantArrays® work?

The QuantArray®-Chlor in many respects is a hybrid technology combining the highly parallel detection of microarrays with the accurate and precise quantification provided by qPCR into a single platform. The key to highly parallel qPCR reactions is the nanoliter fluidics platform for low volume, solution phase qPCR reactions.

How are QuantArray® results reported?

One of the primary advantages of the QuantArray®-Chlor is the simultaneous quantification of a broad spectrum of different microorganisms and key functional genes involved in a variety of pathways for chlorinated hydrocarbon biodegradation. However, highly parallel quantification combined with the various metabolic and cometabolic capabilities of different target organisms can complicate data presentation. Therefore, in addition to Summary Tables, QuantArray® results will be presented as Microbial Population Summary and Comparison Figures to aid in data interpretation and subsequent evaluation of site management activities.

Types of Tables and Figures:

Microbial Population Summary

Figure presenting the concentrations of QuantArray®-Chlor target populations (e.g. *Dehalococcoides*) and functional genes (e.g. vinyl chloride reductase) relative to typically observed values.

Summary Tables

Tables of target population concentrations grouped by biodegradation pathway and contaminant type.

Comparison Figures

Depending on the project, sample results can be presented to compare changes over time or examine differences in microbial populations along a transect of the dissolved plume.

Results

Table 1: Summary of the QuantArray®-Chlor results obtained for samples BM-S8-180119-475, BM-S8-180119-499, BM-S7-220119-469, BM-S7-220119-485, and BM-S7-220119-495.

Sample Name	BM-S8-180119-475	BM-S8-180119-499	BM-S7-220119-469	BM-S7-220119-485	BM-S7-220119-495
Sample Date	01/18/2019	01/18/2019	01/22/2019	01/22/2019	01/22/2019
<i>Reductive Dechlorination</i>	cells/g	cells/g	cells/g	cells/g	cells/g
<i>Dehalococcoides</i> (DHC)	<1.00E+03	<1.00E+03	<1.00E+03	<1.00E+03	<1.00E+03
tceA Reductase (TCE)	<1.00E+03	<1.00E+03	<1.00E+03	<1.00E+03	<1.00E+03
BAV1 Vinyl Chloride Reductase (BVC)	<1.00E+03	<1.00E+03	<1.00E+03	<1.00E+03	<1.00E+03
Vinyl Chloride Reductase (VCR)	<1.00E+03	<1.00E+03	<1.00E+03	<1.00E+03	<1.00E+03
<i>Dehalobacter</i> spp. (DHBt)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalobacter</i> DCM (DCM)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalogenimonas</i> spp. (DHG)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
cerA Reductase (CER)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
trans-1,2-DCE Reductase (TDR)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Desulfitobacterium</i> spp. (DSB)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalobium chlorocoercia</i> (DECO)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Desulfuromonas</i> spp. (DSM)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
PCE Reductase (PCE-1)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
PCE Reductase (PCE-2)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Chloroform Reductase (CFR)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
1,1 DCA Reductase (DCA)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
1,2 DCA Reductase (DCAR)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Aerobic (Co)Metabolic</i>					
Soluble Methane Monooxygenase (SMMO)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Dioxygenase (TOD)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Phenol Hydroxylase (PHE)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Trichlorobenzene Dioxygenase (TCBO)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Monooxygenase 2 (RDEG)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Monooxygenase (RMO)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Ethene Monooxygenase (EtnC)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Epoxyalkane Transferase (EtnE)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Dichloromethane Dehalogenase (DCMA)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Other</i>					
Total Eubacteria (EBAC)	1.41E+05 (I)	8.59E+05 (I)	1.37E+04 (I)	2.52E+04 (I)	2.58E+05 (I)
Sulfate Reducing Bacteria (APS)	<1.00E+04	8.72E+02 (J)	<1.00E+04	<1.00E+04	<1.00E+04
Methanogens (MGN)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04

Legend:

NA = Not Analyzed
I = Inhibited

NS = Not Sampled
< = Result Not Detected

J = Estimated Gene Copies Below PQL but Above LQL

Table 2: Summary of the QuantArray®-Chlor results obtained for samples BM-247-300119-143, BM-247-310119-208, and BM-247-040219-474.

Sample Name Sample Date	BM-247-300119-143 01/30/2019	BM-247-310119-208 01/31/2019	BM-247-040219-474 02/04/2019
<i>Reductive Dechlorination</i>	cells/g	cells/g	cells/g
<i>Dehalococcoides</i> (DHC)	<1.00E+03	<1.00E+03	<1.00E+03
tceA Reductase (TCE)	<1.00E+03	<1.00E+03	<1.00E+03
BAV1 Vinyl Chloride Reductase (BVC)	<1.00E+03	<1.00E+03	<1.00E+03
Vinyl Chloride Reductase (VCR)	<1.00E+03	<1.00E+03	<1.00E+03
<i>Dehalobacter</i> spp. (DHBt)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalobacter</i> DCM (DCM)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalogenimonas</i> spp. (DHG)	<1.00E+04	<1.00E+04	<1.00E+04
cerA Reductase (CER)	<1.00E+04	<1.00E+04	<1.00E+04
trans-1,2-DCE Reductase (TDR)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Desulfitobacterium</i> spp. (DSB)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalobium chlorocoercia</i> (DECO)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Desulfuromonas</i> spp. (DSM)	<1.00E+04	<1.00E+04	<1.00E+04
PCE Reductase (PCE-1)	<1.00E+04	<1.00E+04	<1.00E+04
PCE Reductase (PCE-2)	<1.00E+04	<1.00E+04	<1.00E+04
Chloroform Reductase (CFR)	<1.00E+04	<1.00E+04	<1.00E+04
1,1 DCA Reductase (DCA)	<1.00E+04	<1.00E+04	<1.00E+04
1,2 DCA Reductase (DCAR)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Aerobic (Co)Metabolic</i>			
Soluble Methane Monooxygenase (SMMO)	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Dioxygenase (TOD)	<1.00E+04	<1.00E+04	<1.00E+04
Phenol Hydroxylase (PHE)	<1.00E+04	<1.00E+04	<1.00E+04
Trichlorobenzene Dioxygenase (TCBO)	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Monooxygenase 2 (RDEG)	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Monooxygenase (RMO)	<1.00E+04	<1.00E+04	<1.00E+04
Ethene Monooxygenase (EtnC)	<1.00E+04	<1.00E+04	<1.00E+04
Epoxyalkane Transferase (EtnE)	<1.00E+04	<1.00E+04	<1.00E+04
Dichloromethane Dehalogenase (DCMA)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Other</i>			
Total Eubacteria (EBAC)	3.61E+04 (I)	5.76E+04 (I)	2.08E+03 (I)
Sulfate Reducing Bacteria (APS)	<1.00E+04	<1.00E+04	<1.00E+04
Methanogens (MGN)	<1.00E+04	<1.00E+04	<1.00E+04

Legend:

NA = Not Analyzed
I = Inhibited

NS = Not Sampled
< = Result Not Detected

J = Estimated Gene Copies Below PQL but Above LQL

Table 3: Summary of the QuantArray®-Chlor results obtained for samples BM-247-050219-480, BM-247-050219-489, and BM-247-050219-499.

Sample Name	BM-247-050219-480	BM-247-050219-489	BM-247-050219-499
Sample Date	02/05/2019	02/05/2019	02/05/2019
<i>Reductive Dechlorination</i>	cells/g	cells/g	cells/g
<i>Dehalococcoides</i> (DHC)	<1.00E+03	<1.00E+03	<1.00E+03
tceA Reductase (TCE)	<1.00E+03	<1.00E+03	<1.00E+03
BAV1 Vinyl Chloride Reductase (BVC)	<1.00E+03	<1.00E+03	<1.00E+03
Vinyl Chloride Reductase (VCR)	<1.00E+03	<1.00E+03	<1.00E+03
<i>Dehalobacter</i> spp. (DHBt)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalobacter</i> DCM (DCM)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalogenimonas</i> spp. (DHG)	<1.00E+04	<1.00E+04	<1.00E+04
cerA Reductase (CER)	<1.00E+04	<1.00E+04	<1.00E+04
trans-1,2-DCE Reductase (TDR)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Desulfitobacterium</i> spp. (DSB)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalobium chlorocoercia</i> (DECO)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Desulfuromonas</i> spp. (DSM)	<1.00E+04	<1.00E+04	<1.00E+04
PCE Reductase (PCE-1)	<1.00E+04	<1.00E+04	<1.00E+04
PCE Reductase (PCE-2)	<1.00E+04	<1.00E+04	<1.00E+04
Chloroform Reductase (CFR)	<1.00E+04	<1.00E+04	<1.00E+04
1,1 DCA Reductase (DCA)	<1.00E+04	<1.00E+04	<1.00E+04
1,2 DCA Reductase (DCAR)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Aerobic (Co)Metabolic</i>			
Soluble Methane Monooxygenase (SMMO)	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Dioxygenase (TOD)	<1.00E+04	<1.00E+04	<1.00E+04
Phenol Hydroxylase (PHE)	<1.00E+04	<1.00E+04	<1.00E+04
Trichlorobenzene Dioxygenase (TCBO)	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Monooxygenase 2 (RDEG)	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Monooxygenase (RMO)	<1.00E+04	<1.00E+04	<1.00E+04
Ethene Monooxygenase (EtnC)	<1.00E+04	<1.00E+04	<1.00E+04
Epoxyalkane Transferase (EtnE)	<1.00E+04	<1.00E+04	<1.00E+04
Dichloromethane Dehalogenase (DCMA)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Other</i>			
Total Eubacteria (EBAC)	5.25E+04 (I)	1.54E+06 (I)	1.69E+04 (I)
Sulfate Reducing Bacteria (APS)	<1.00E+04	<1.00E+04	<1.00E+04
Methanogens (MGN)	<1.00E+04	<1.00E+04	<1.00E+04

Legend:

NA = Not Analyzed
I = Inhibited

NS = Not Sampled
< = Result Not Detected

J = Estimated Gene Copies Below PQL but Above LQL

Microbial Populations BM-S8-180119-475

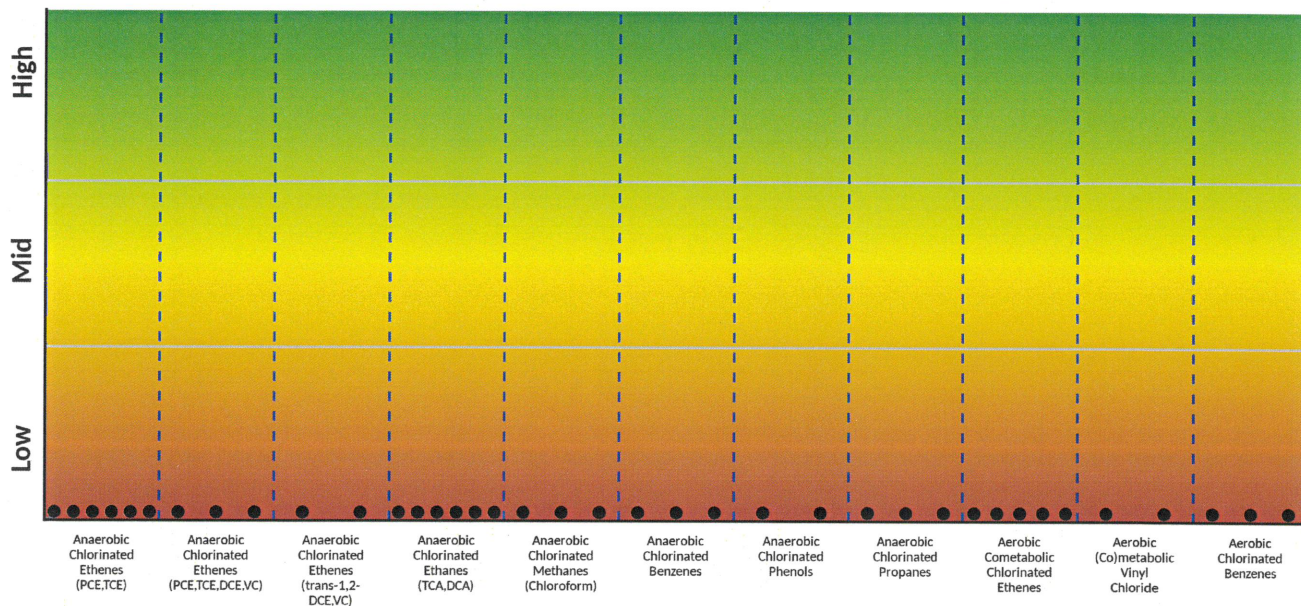


Figure 1: Microbial population summary to aid in evaluating potential pathways and biodegradation of specific contaminants.

Anaerobic - Reductive Dechlorination or Dichloroelimination		Aerobic - (Co)metabolism	
Chlorinated Ethenes (PCE, TCE)	DHC, DHBt, DSB, DSM, PCE-1, PCE-2	Chlorinated Ethenes (TCE,DCE,VC)	sMMO, TOD, PHE, RDEG, RMO
Chlorinated Ethenes (PCE, TCE, DCE, VC)	DHC, BVC, VCR	(Co)metabolic Vinyl Chloride	etnC, etnE
Chlorinated Ethenes (trans-1,2-DCE, VC)	TDR, CER	Chlorinated Benzenes	TOD, TCBO, PHE
Chlorinated Ethanes (TCA and 1,2-DCA)	DHC, DHBt, DHG, DSB ¹ , DCA, DCAR		
Chlorinated Methanes (Chloroform)	DHBt, DCM, CFR		
Chlorinated Benzenes	DHC, DHBt ² , DECO		
Chlorinated Phenols	DHC, DSB		
Chlorinated Propanes	DHC, DHG, DSB ¹		

¹*Desulfotobacterium dichloroeliminans* DCA1. ²Implicated in reductive dechlorination of dichlorobenzene and potentially chlorobenzene.

Microbial Populations BM-S8-180119-499

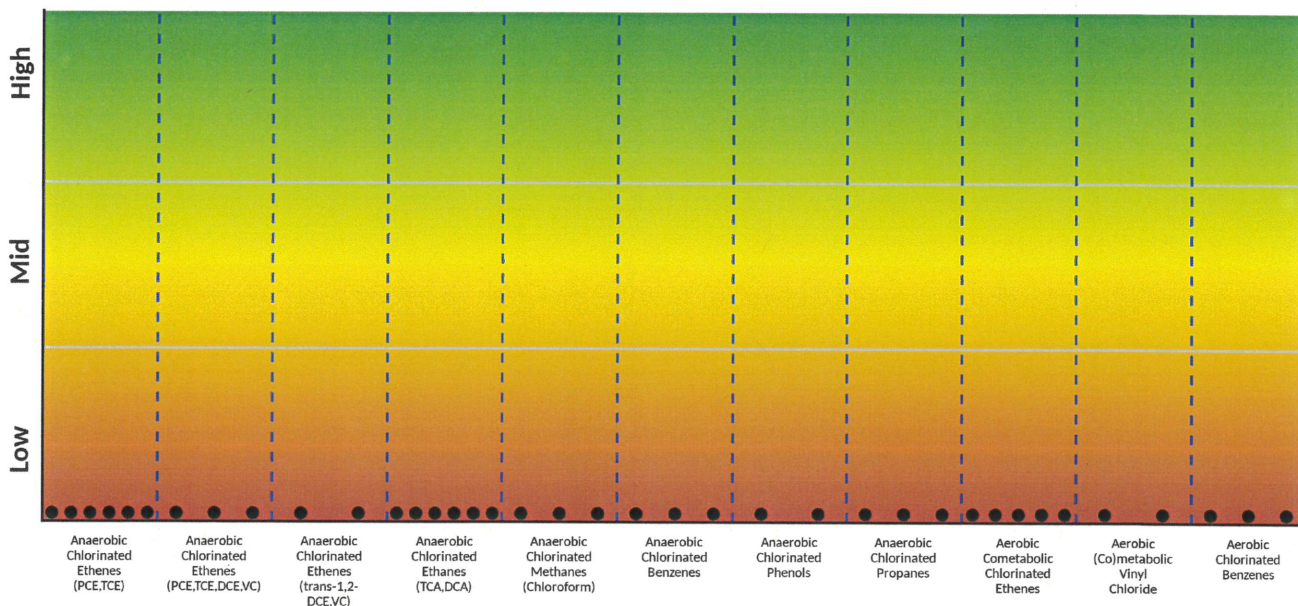


Figure 2: Microbial population summary to aid in evaluating potential pathways and biodegradation of specific contaminants.

Anaerobic - Reductive Dechlorination or Dichloroelimination		Aerobic - (Co)metabolism	
Chlorinated Ethenes (PCE, TCE)	DHC, DHBt, DSB, DSM, PCE-1, PCE-2	Chlorinated Ethenes (TCE,DCE,VC)	sMMO, TOD, PHE, RDEG, RMO
Chlorinated Ethenes (PCE, TCE, DCE, VC)	DHC, BVC, VCR	(Co)metabolic Vinyl Chloride	etnC, etnE
Chlorinated Ethenes (trans-1,2-DCE, VC)	TDR, CER	Chlorinated Benzenes	TOD, TCBO, PHE
Chlorinated Ethanes (TCA and 1,2-DCA)	DHC, DHBt, DHG, DSB ¹ , DCA, DCAR		
Chlorinated Methanes (Chloroform)	DHBt, DCM, CFR		
Chlorinated Benzenes	DHC, DHBt ² , DECO		
Chlorinated Phenols	DHC, DSB		
Chlorinated Propanes	DHC, DHG, DSB ¹		

¹ *Desulfitobacterium dichloroeliminans* DCA1. ² Implicated in reductive dechlorination of dichlorobenzene and potentially chlorobenzene.

Microbial Populations BM-S7-220119-469

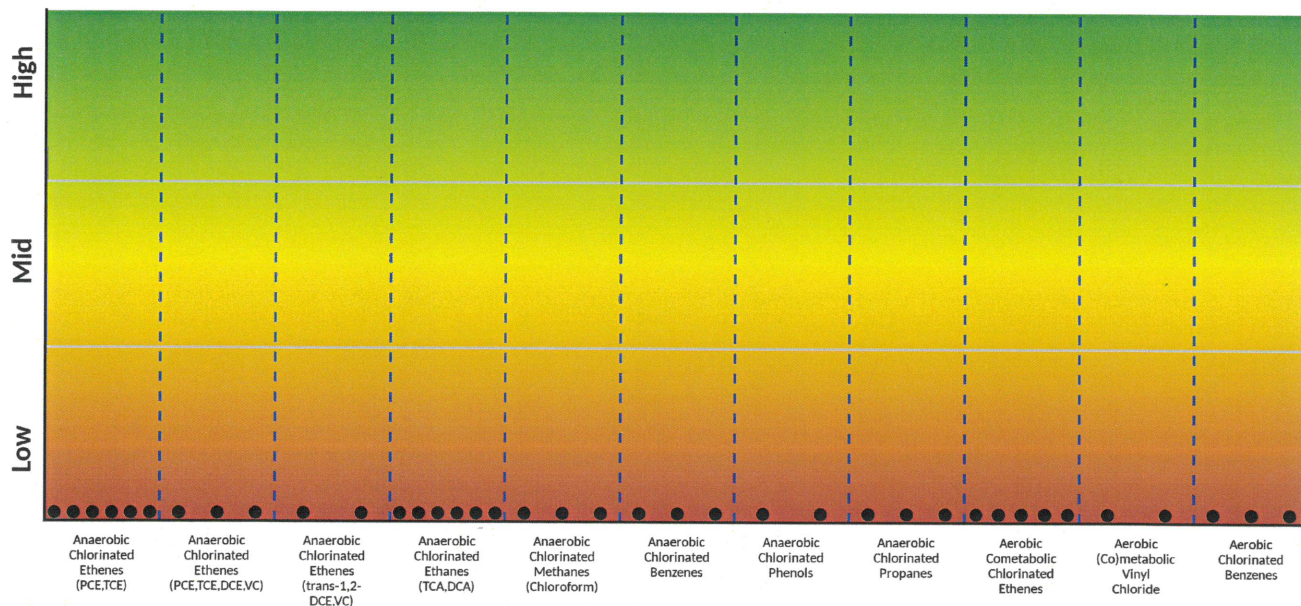


Figure 3: Microbial population summary to aid in evaluating potential pathways and biodegradation of specific contaminants.

Anaerobic - Reductive Dechlorination or Dichloroelimination		Aerobic - (Co)metabolism	
Chlorinated Ethenes (PCE, TCE)	DHC, DHBt, DSB, DSM, PCE-1, PCE-2	Chlorinated Ethenes (TCE,DCE,VC)	sMMO, TOD, PHE, RDEG, RMO
Chlorinated Ethenes (PCE, TCE, DCE, VC)	DHC, BVC, VCR	(Co)metabolic Vinyl Chloride	etnC, etnE
Chlorinated Ethenes (trans-1,2-DCE, VC)	TDR, CER	Chlorinated Benzenes	TOD, TCBO, PHE
Chlorinated Ethanes (TCA and 1,2-DCA)	DHC, DHBt, DHG, DSB ¹ , DCA, DCAR		
Chlorinated Methanes (Chloroform)	DHBt, DCM, CFR		
Chlorinated Benzenes	DHC, DHBt ² , DECO		
Chlorinated Phenols	DHC, DSB		
Chlorinated Propanes	DHC, DHG, DSB ¹		

¹Desulfotobacterium dichloroeliminans DCA1. ²Implicated in reductive dechlorination of dichlorobenzene and potentially chlorobenzene.

Microbial Populations BM-S7-220119-485

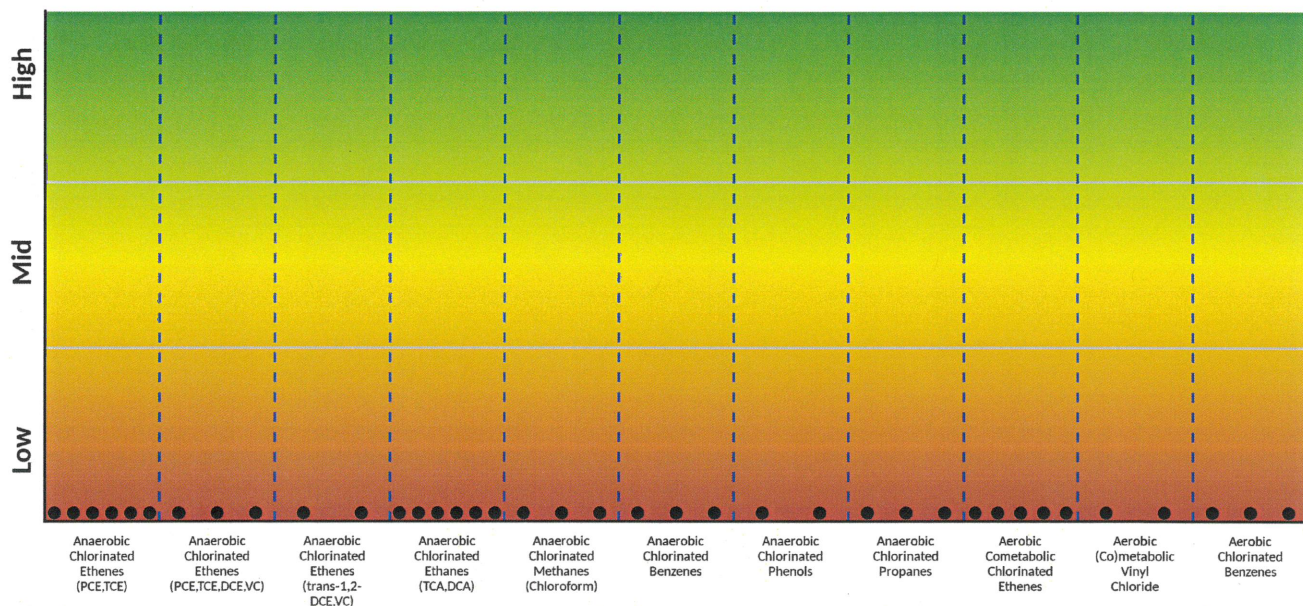


Figure 4: Microbial population summary to aid in evaluating potential pathways and biodegradation of specific contaminants.

Anaerobic - Reductive Dechlorination or Dichloroelimination

Chlorinated Ethenes (PCE, TCE)	DHC, DHBt, DSB, DSM, PCE-1, PCE-2
Chlorinated Ethenes (PCE, TCE, DCE, VC)	DHC, BVC, VCR
Chlorinated Ethenes (trans-1,2-DCE, VC)	TDR, CER
Chlorinated Ethanes (TCA and 1,2-DCA)	DHC, DHBt, DHG, DSB ¹ , DCA, DCAR
Chlorinated Methanes (Chloroform)	DHBt, DCM, CFR
Chlorinated Benzenes	DHC, DHBt ² , DECO
Chlorinated Phenols	DHC, DSB
Chlorinated Propanes	DHC, DHG, DSB ¹

Aerobic - (Co)metabolism

Chlorinated Ethenes (TCE,DCE,VC) (Co)metabolic Vinyl Chloride	sMMO, TOD, PHE, RDEG, RMO etnC, etnE
Chlorinated Benzenes	TOD, TCBO, PHE

¹ *Desulfotobacterium dichloroelimians* DCA1. ² Implicated in reductive dechlorination of dichlorobenzene and potentially chlorobenzene.

Microbial Populations BM-S7-220119-495

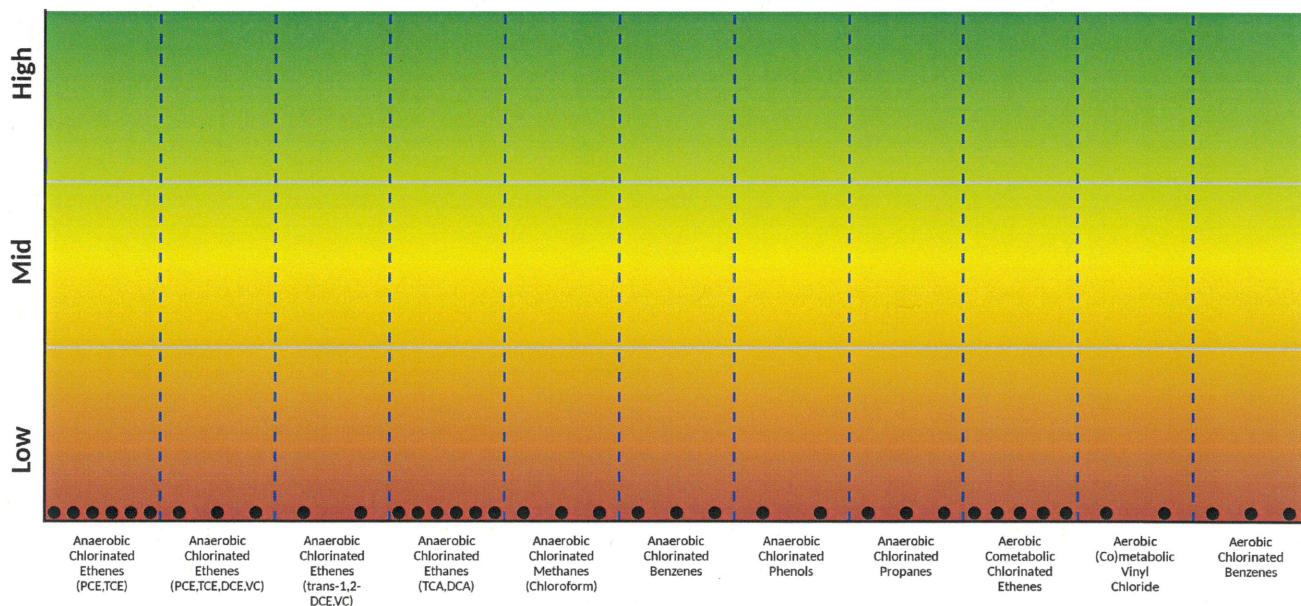


Figure 5: Microbial population summary to aid in evaluating potential pathways and biodegradation of specific contaminants.

Anaerobic - Reductive Dechlorination or Dichloroelimination			Aerobic - (Co)metabolism	
Chlorinated Ethenes (PCE, TCE)	DHC, DHBt, DSB, DSM, PCE-1, PCE-2	Chlorinated Ethenes (TCE,DCE,VC)	sMMO, TOD, PHE, RDEG, RMO	
Chlorinated Ethenes (PCE, TCE, DCE, VC)	DHC, BVC, VCR	(Co)metabolic Vinyl Chloride	etnC, etnE	
Chlorinated Ethenes (trans-1,2-DCE, VC)	TDR, CER	Chlorinated Benzenes	TOD, TCBO, PHE	
Chlorinated Ethanes (TCA and 1,2-DCA)	DHC, DHBt, DHG, DSB ¹ , DCA, DCAR			
Chlorinated Methanes (Chloroform)	DHBt, DCM, CFR			
Chlorinated Benzenes	DHC, DHBt ² , DECO			
Chlorinated Phenols	DHC, DSB			
Chlorinated Propanes	DHC, DHG, DSB ¹			

¹*Desulfotobacterium dichloroelimians* DCA1. ²Implicated in reductive dechlorination of dichlorobenzene and potentially chlorobenzene.

Microbial Populations BM-247-300119-143

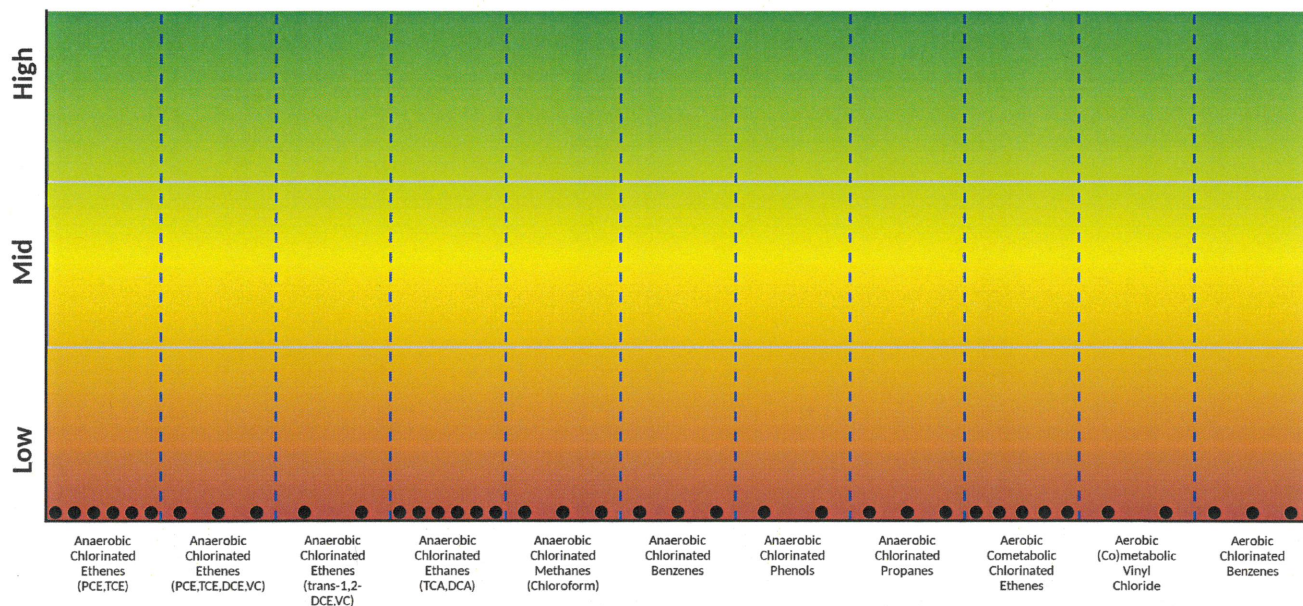


Figure 6: Microbial population summary to aid in evaluating potential pathways and biodegradation of specific contaminants.

Anaerobic - Reductive Dechlorination or Dichloroelimination		Aerobic - (Co)metabolism	
Chlorinated Ethenes (PCE, TCE)	DHC, DHBt, DSB, DSM, PCE-1, PCE-2	Chlorinated Ethenes (TCE,DCE,VC)	sMMO, TOD, PHE, RDEG, RMO
Chlorinated Ethenes (PCE, TCE, DCE, VC)	DHC, BVC, VCR	(Co)metabolic Vinyl Chloride	etnC, etnE
Chlorinated Ethenes (trans-1,2-DCE, VC)	TDR, CER	Chlorinated Benzenes	TOD, TCBO, PHE
Chlorinated Ethanes (TCA and 1,2-DCA)	DHC, DHBt, DHG, DSB ¹ , DCA, DCAR		
Chlorinated Methanes (Chloroform)	DHBt, DCM, CFR		
Chlorinated Benzenes	DHC, DHBt ² , DECO		
Chlorinated Phenols	DHC, DSB		
Chlorinated Propanes	DHC, DHG, DSB ¹		

¹ *Desulfitobacterium dichloroeliminans* DCA1. ² Implicated in reductive dechlorination of dichlorobenzene and potentially chlorobenzene.

Microbial Populations BM-247-310119-208

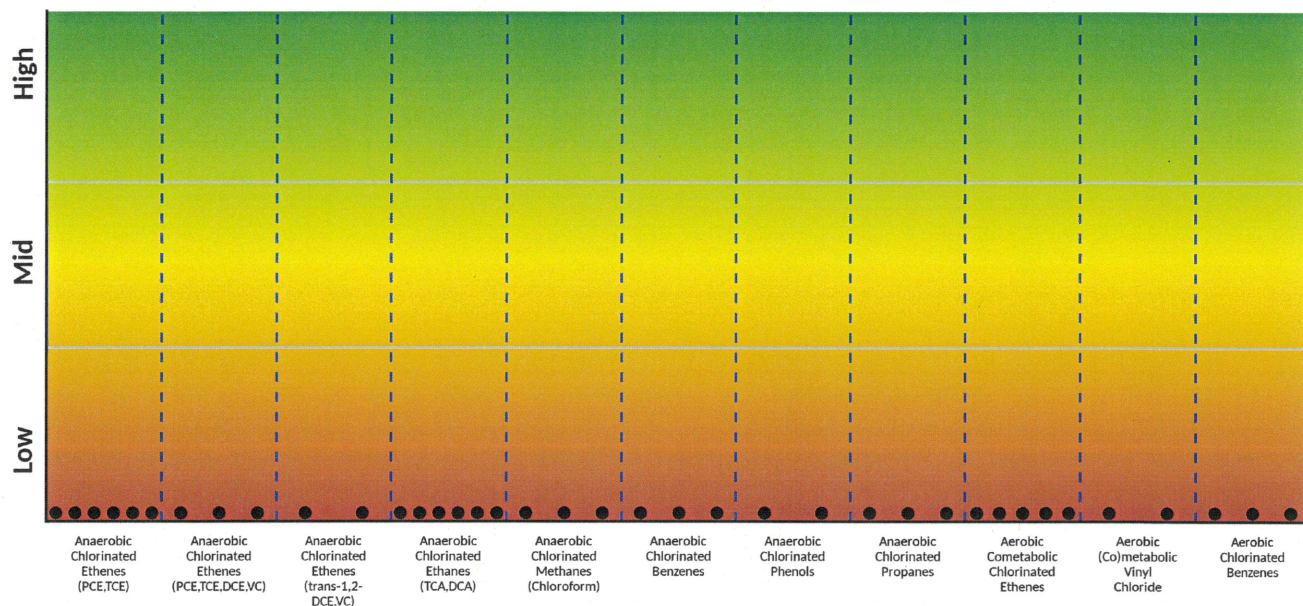


Figure 7: Microbial population summary to aid in evaluating potential pathways and biodegradation of specific contaminants.

Anaerobic - Reductive Dechlorination or Dichloroelimination			Aerobic - (Co)metabolism	
Chlorinated Ethenes (PCE, TCE)	DHC, DHBt, DSB, DSM, PCE-1, PCE-2	Chlorinated Ethenes (TCE,DCE,VC)	sMMO, TOD, PHE, RDEG, RMO	
Chlorinated Ethenes (PCE, TCE, DCE, VC)	DHC, BVC, VCR	(Co)metabolic Vinyl Chloride	etnC, etnE	
Chlorinated Ethenes (trans-1,2-DCE, VC)	TDR, CER	Chlorinated Benzenes	TOD, TCBO, PHE	
Chlorinated Ethanes (TCA and 1,2-DCA)	DHC, DHBt, DHG, DSB ¹ , DCA, DCAR			
Chlorinated Methanes (Chloroform)	DHBt, DCM, CFR			
Chlorinated Benzenes	DHC, DHBt ² , DECO			
Chlorinated Phenols	DHC, DSB			
Chlorinated Propanes	DHC, DHG, DSB ¹			

¹ *Desulfotobacterium dichloroelimians* DCA1. ² Implicated in reductive dechlorination of dichlorobenzene and potentially chlorobenzene.

Microbial Populations BM-247-040219-474

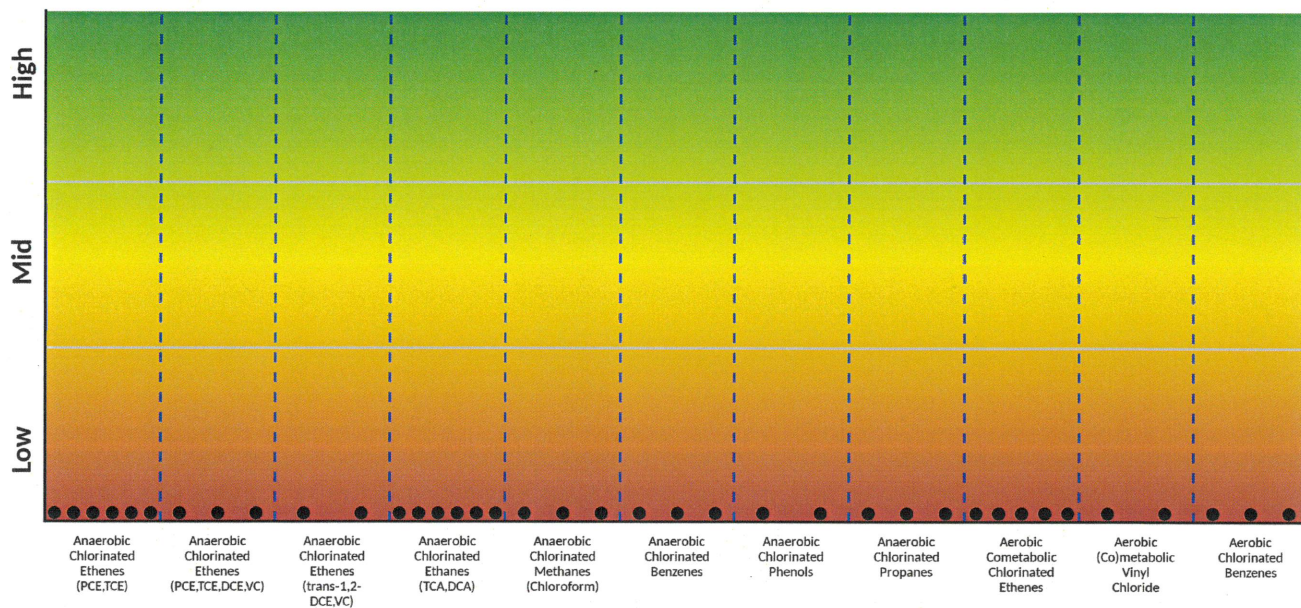


Figure 8: Microbial population summary to aid in evaluating potential pathways and biodegradation of specific contaminants.

Anaerobic - Reductive Dechlorination or Dichloroelimination		Aerobic - (Co)metabolism	
Chlorinated Ethenes (PCE, TCE)	DHC, DHBt, DSB, DSM, PCE-1, PCE-2	Chlorinated Ethenes (TCE,DCE,VC)	sMMO, TOD, PHE, RDEG, RMO
Chlorinated Ethenes (PCE, TCE, DCE, VC)	DHC, BVC, VCR	(Co)metabolic Vinyl Chloride	etnC, etnE
Chlorinated Ethenes (trans-1,2-DCE, VC)	TDR, CER	Chlorinated Benzenes	TOD, TCBO, PHE
Chlorinated Ethanes (TCA and 1,2-DCA)	DHC, DHBt, DHG, DSB ¹ , DCA, DCAR		
Chlorinated Methanes (Chloroform)	DHBt, DCM, CFR		
Chlorinated Benzenes	DHC, DHBt ² , DECO		
Chlorinated Phenols	DHC, DSB		
Chlorinated Propanes	DHC, DHG, DSB ¹		

¹ *Desulfotobacterium dichloroeliminans* DCA1. ² Implicated in reductive dechlorination of dichlorobenzene and potentially chlorobenzene.

Microbial Populations BM-247-050219-480

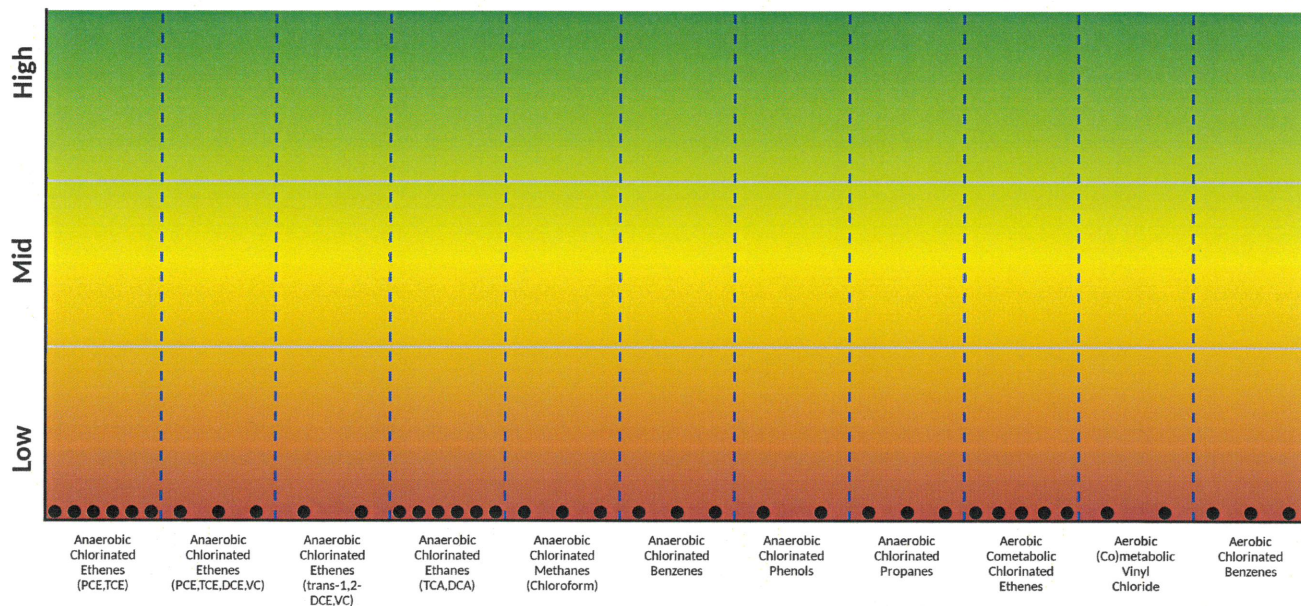


Figure 9: Microbial population summary to aid in evaluating potential pathways and biodegradation of specific contaminants.

Anaerobic - Reductive Dechlorination or Dichloroelimination		Aerobic - (Co)metabolism	
Chlorinated Ethenes (PCE, TCE)	DHC, DHBt, DSB, DSM, PCE-1, PCE-2	Chlorinated Ethenes (TCE,DCE,VC)	sMMO, TOD, PHE, RDEG, RMO
Chlorinated Ethenes (PCE, TCE, DCE, VC)	DHC, BVC, VCR	(Co)metabolic Vinyl Chloride	etnC, etnE
Chlorinated Ethenes (trans-1,2-DCE, VC)	TDR, CER	Chlorinated Benzenes	TOD, TCBO, PHE
Chlorinated Ethanes (TCA and 1,2-DCA)	DHC, DHBt, DHG, DSB ¹ , DCA, DCAR		
Chlorinated Methanes (Chloroform)	DHBt, DCM, CFR		
Chlorinated Benzenes	DHC, DHBt ² , DECO		
Chlorinated Phenols	DHC, DSB		
Chlorinated Propanes	DHC, DHG, DSB ¹		

¹*Desulfotobacterium dichloroeliminans* DCA1. ²Implicated in reductive dechlorination of dichlorobenzene and potentially chlorobenzene.

Microbial Populations BM-247-050219-489

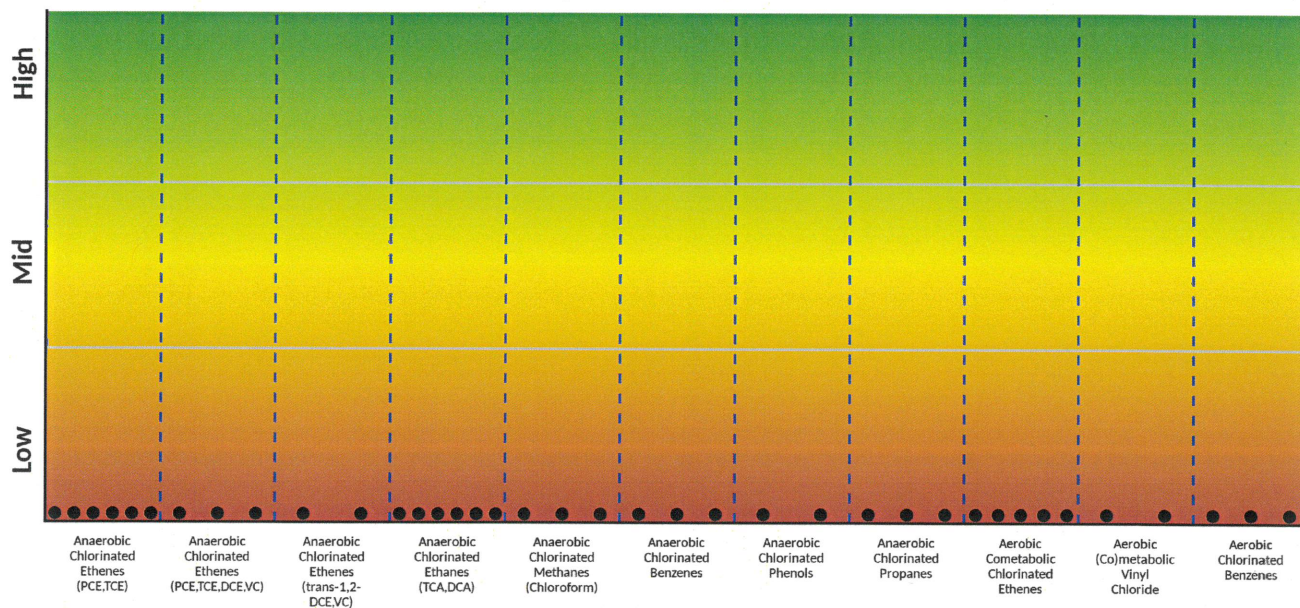


Figure 10: Microbial population summary to aid in evaluating potential pathways and biodegradation of specific contaminants.

Anaerobic - Reductive Dechlorination or Dichloroelimination		Aerobic - (Co)metabolism	
Chlorinated Ethenes (PCE, TCE)	DHC, DHBt, DSB, DSM, PCE-1, PCE-2	Chlorinated Ethenes (TCE,DCE,VC)	sMMO, TOD, PHE, RDEG, RMO
Chlorinated Ethenes (PCE, TCE, DCE, VC)	DHC, BVC, VCR	(Co)metabolic Vinyl Chloride	etnC, etnE
Chlorinated Ethenes (trans-1,2-DCE, VC)	TDR, CER	Chlorinated Benzenes	TOD, TCBO, PHE
Chlorinated Ethanes (TCA and 1,2-DCA)	DHC, DHBt, DHG, DSB ¹ , DCA, DCAR		
Chlorinated Methanes (Chloroform)	DHBt, DCM, CFR		
Chlorinated Benzenes	DHC, DHBt ² , DECO		
Chlorinated Phenols	DHC, DSB		
Chlorinated Propanes	DHC, DHG, DSB ¹		

¹*Desulfotobacterium dichloroeliminans* DCA1. ²Implicated in reductive dechlorination of dichlorobenzene and potentially chlorobenzene.

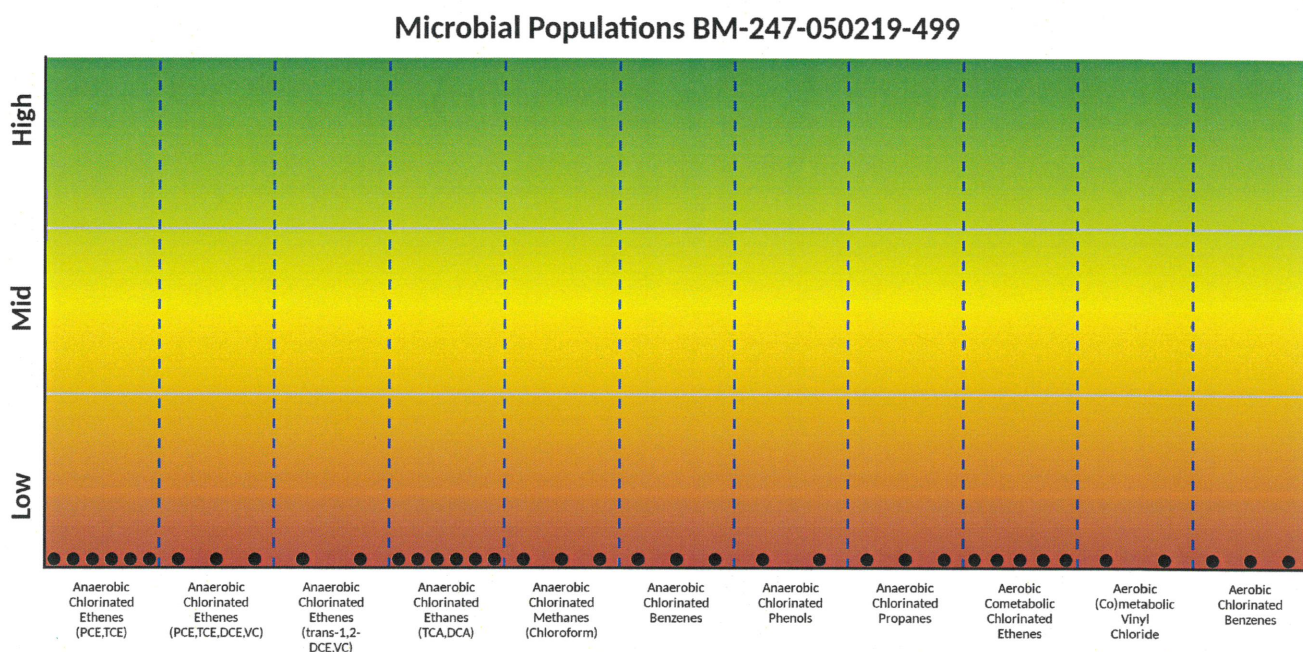


Figure 11: Microbial population summary to aid in evaluating potential pathways and biodegradation of specific contaminants.

Anaerobic - Reductive Dechlorination or Dichloroelimination

Chlorinated Ethenes (PCE, TCE)	DHC, DHBt, DSB, DSM, PCE-1, PCE-2
Chlorinated Ethenes (PCE, TCE, DCE, VC)	DHC, BVC, VCR
Chlorinated Ethenes (trans-1,2-DCE, VC)	TDR, CER
Chlorinated Ethanes (TCA and 1,2-DCA)	DHC, DHBt, DHG, DSB ¹ , DCA, DCAR
Chlorinated Methanes (Chloroform)	DHBt, DCM, CFR
Chlorinated Benzenes	DHC, DHBt ² , DECO
Chlorinated Phenols	DHC, DSB
Chlorinated Propanes	DHC, DHG, DSB ¹

Aerobic - (Co)metabolism

Chlorinated Ethenes (TCE,DCE,VC) (Co)metabolic Vinyl Chloride	sMMO, TOD, PHE, RDEG, RMO etnC, etnE
Chlorinated Benzenes	TOD, TCBO, PHE

¹*Desulfotobacterium dichloroeliminans* DCA1. ²Implicated in reductive dechlorination of dichlorobenzene and potentially chlorobenzene.

Table 4: Summary of the QuantArray®-Chlor results for microorganisms responsible for reductive dechlorination for samples BM-S8-180119-475, BM-S8-180119-499, BM-S7-220119-469, BM-S7-220119-485, and BM-S7-220119-495.

Sample Name	BM-S8-180119-475	BM-S8-180119-499	BM-S7-220119-469	BM-S7-220119-485	BM-S7-220119-495
Sample Date	01/18/2019	01/18/2019	01/22/2019	01/22/2019	01/22/2019
Reductive Dechlorination	cells/g	cells/g	cells/g	cells/g	cells/g
<i>Dehalococcoides</i> (DHC)	<1.00E+03	<1.00E+03	<1.00E+03	<1.00E+03	<1.00E+03
tceA Reductase (TCE)	<1.00E+03	<1.00E+03	<1.00E+03	<1.00E+03	<1.00E+03
BAV1 Vinyl Chloride Reductase (BVC)	<1.00E+03	<1.00E+03	<1.00E+03	<1.00E+03	<1.00E+03
Vinyl Chloride Reductase (VCR)	<1.00E+03	<1.00E+03	<1.00E+03	<1.00E+03	<1.00E+03
<i>Dehalobacter</i> spp. (DHBt)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalobacter</i> DCM (DCM)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalogenimonas</i> spp. (DHG)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Desulfotobacterium</i> spp. (DSB)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalobium chlorocoercia</i> (DECO)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
<i>Desulfuromonas</i> spp. (DSM)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04

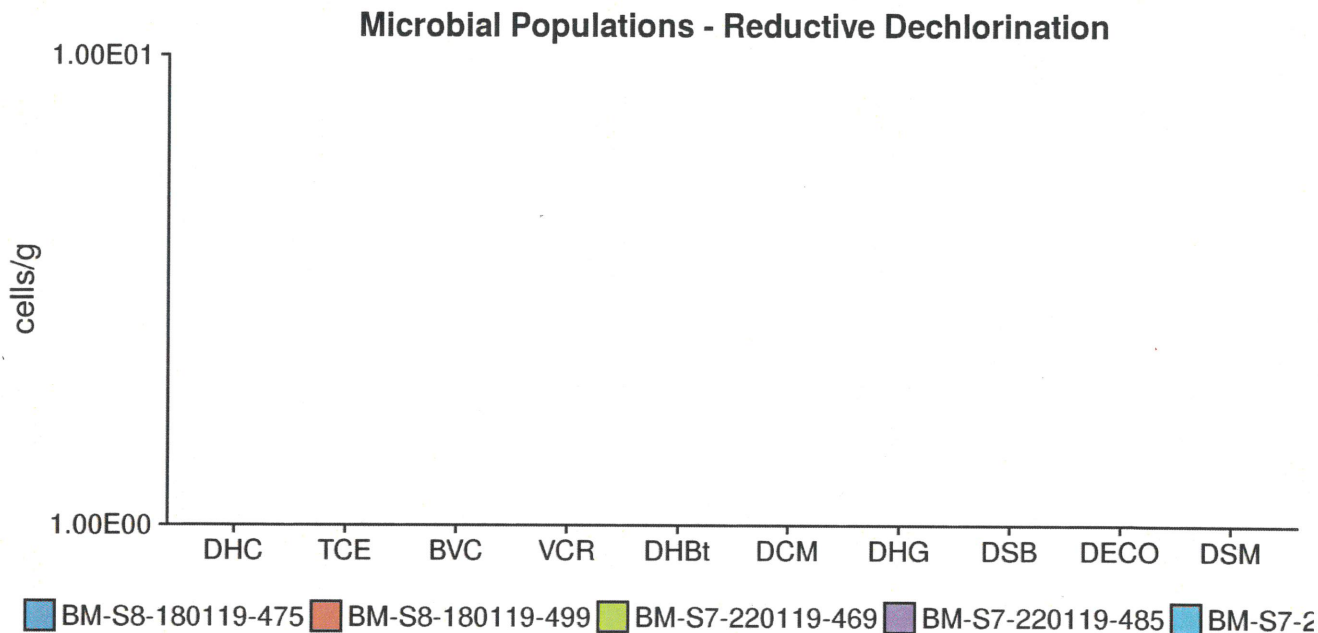


Figure 12: Comparison - microbial populations involved in reductive dechlorination.

Table 5: Summary of the QuantArray®-Chlor results for microorganisms responsible for reductive dechlorination for samples BM-S8-180119-475, BM-S8-180119-499, BM-S7-220119-469, BM-S7-220119-485, and BM-S7-220119-495.

Sample Name	BM-S8-180119-475	BM-S8-180119-499	BM-S7-220119-469	BM-S7-220119-485	BM-S7-220119-495
Sample Date	01/18/2019	01/18/2019	01/22/2019	01/22/2019	01/22/2019
Reductive Dechlorination	cells/g	cells/g	cells/g	cells/g	cells/g
Chloroform Reductase (CFR)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
1,1 DCA Reductase (DCA)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
1,2 DCA Reductase (DCAR)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
PCE Reductase (PCE-1)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
PCE Reductase (PCE-2)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Dehalogenimonas trans-1,2-DCE Reductase (TDR)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Dehalogenimonas cerA Reductase (CER)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04

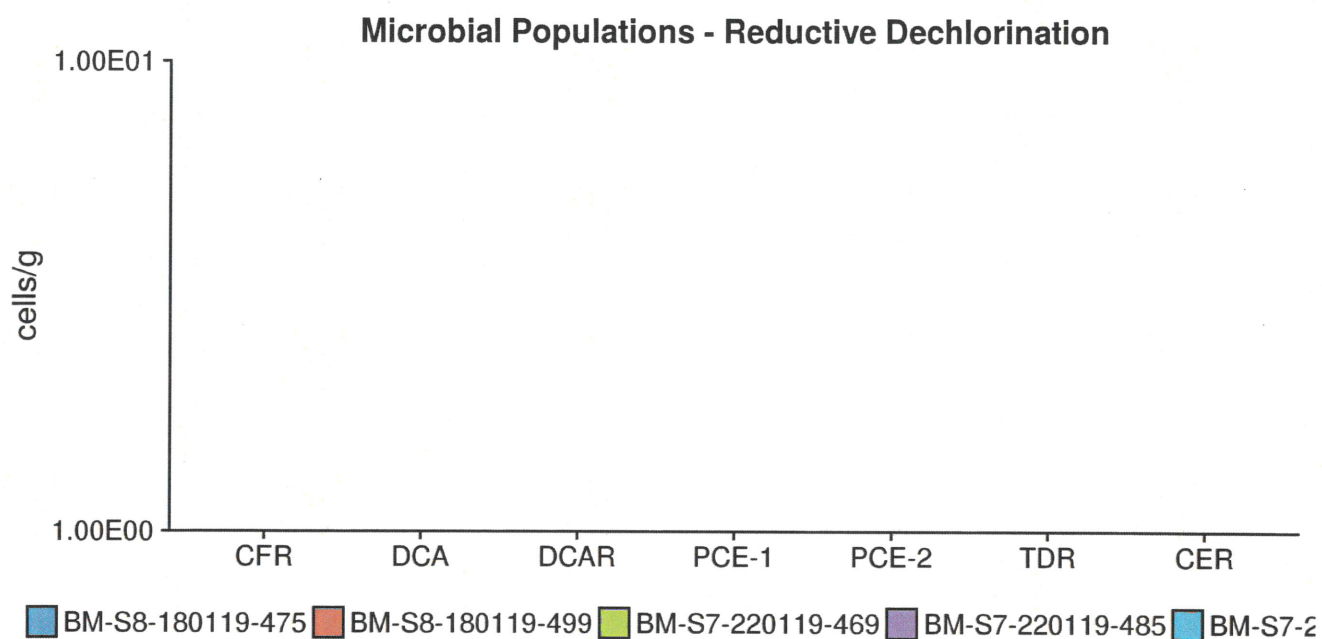


Figure 13: Comparison - microbial populations involved in reductive dechlorination.

Table 6: Summary of the QuantArray®-Chlor results for microorganisms responsible for reductive dechlorination for samples BM-247-300119-143, BM-247-310119-208, and BM-247-040219-474.

Sample Name	BM-247-300119-143	BM-247-310119-208	BM-247-040219-474
Sample Date	01/30/2019	01/31/2019	02/04/2019
Reductive Dechlorination	cells/g	cells/g	cells/g
<i>Dehalococcoides</i> (DHC)	<1.00E+03	<1.00E+03	<1.00E+03
tceA Reductase (TCE)	<1.00E+03	<1.00E+03	<1.00E+03
BAV1 Vinyl Chloride Reductase (BVC)	<1.00E+03	<1.00E+03	<1.00E+03
Vinyl Chloride Reductase (VCR)	<1.00E+03	<1.00E+03	<1.00E+03
<i>Dehalobacter</i> spp. (DHBt)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalobacter</i> DCM (DCM)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalogenimonas</i> spp. (DHG)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Desulfitobacterium</i> spp. (DSB)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalobium chlorocoercia</i> (DECO)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Desulfuromonas</i> spp. (DSM)	<1.00E+04	<1.00E+04	<1.00E+04

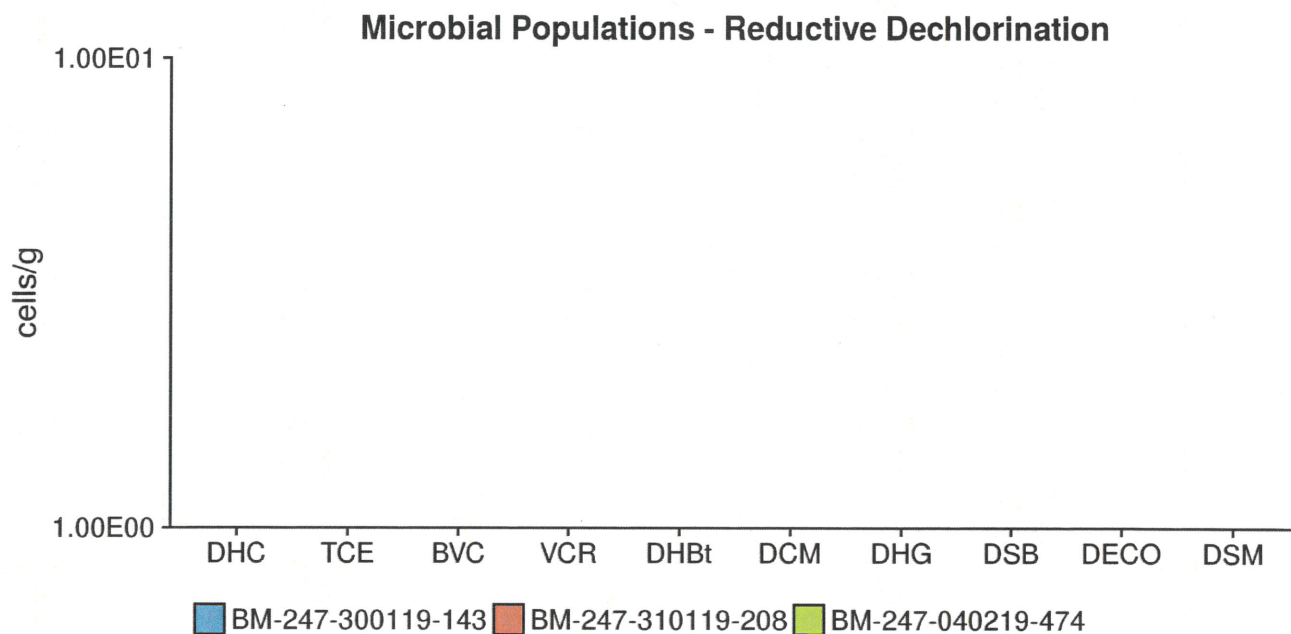


Figure 14: Comparison - microbial populations involved in reductive dechlorination.

Table 7: Summary of the QuantArray®-Chlor results for microorganisms responsible for reductive dechlorination for samples BM-247-300119-143, BM-247-310119-208, and BM-247-040219-474.

Sample Name	BM-247-300119-143	BM-247-310119-208	BM-247-040219-474
Sample Date	01/30/2019	01/31/2019	02/04/2019
<i>Reductive Dechlorination</i>	cells/g	cells/g	cells/g
Chloroform Reductase (CFR)	<1.00E+04	<1.00E+04	<1.00E+04
1,1 DCA Reductase (DCA)	<1.00E+04	<1.00E+04	<1.00E+04
1,2 DCA Reductase (DCAR)	<1.00E+04	<1.00E+04	<1.00E+04
PCE Reductase (PCE-1)	<1.00E+04	<1.00E+04	<1.00E+04
PCE Reductase (PCE-2)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalogenimonas trans</i> -1,2-DCE Reductase (TDR)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalogenimonas cerA</i> Reductase (CER)	<1.00E+04	<1.00E+04	<1.00E+04

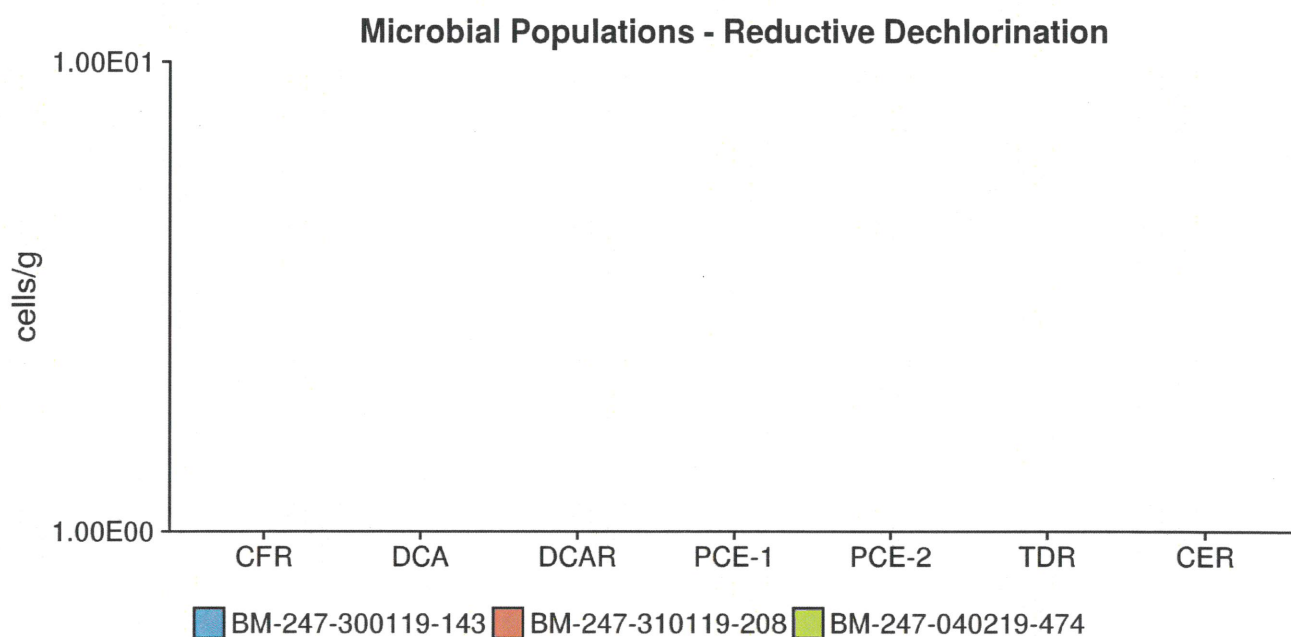


Figure 15: Comparison - microbial populations involved in reductive dechlorination.

Table 8: Summary of the QuantArray®-Chlor results for microorganisms responsible for reductive dechlorination for samples BM-247-050219-480, BM-247-050219-489, and BM-247-050219-499.

Sample Name	BM-247-050219-480	BM-247-050219-489	BM-247-050219-499
Sample Date	02/05/2019	02/05/2019	02/05/2019
Reductive Dechlorination	cells/g	cells/g	cells/g
<i>Dehalococcoides</i> (DHC)	<1.00E+03	<1.00E+03	<1.00E+03
tceA Reductase (TCE)	<1.00E+03	<1.00E+03	<1.00E+03
BAV1 Vinyl Chloride Reductase (BVC)	<1.00E+03	<1.00E+03	<1.00E+03
Vinyl Chloride Reductase (VCR)	<1.00E+03	<1.00E+03	<1.00E+03
<i>Dehalobacter</i> spp. (DHBt)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalobacter</i> DCM (DCM)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalogenimonas</i> spp. (DHG)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Desulfitobacterium</i> spp. (DSB)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalobium chlorocoercia</i> (DECO)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Desulfuromonas</i> spp. (DSM)	<1.00E+04	<1.00E+04	<1.00E+04

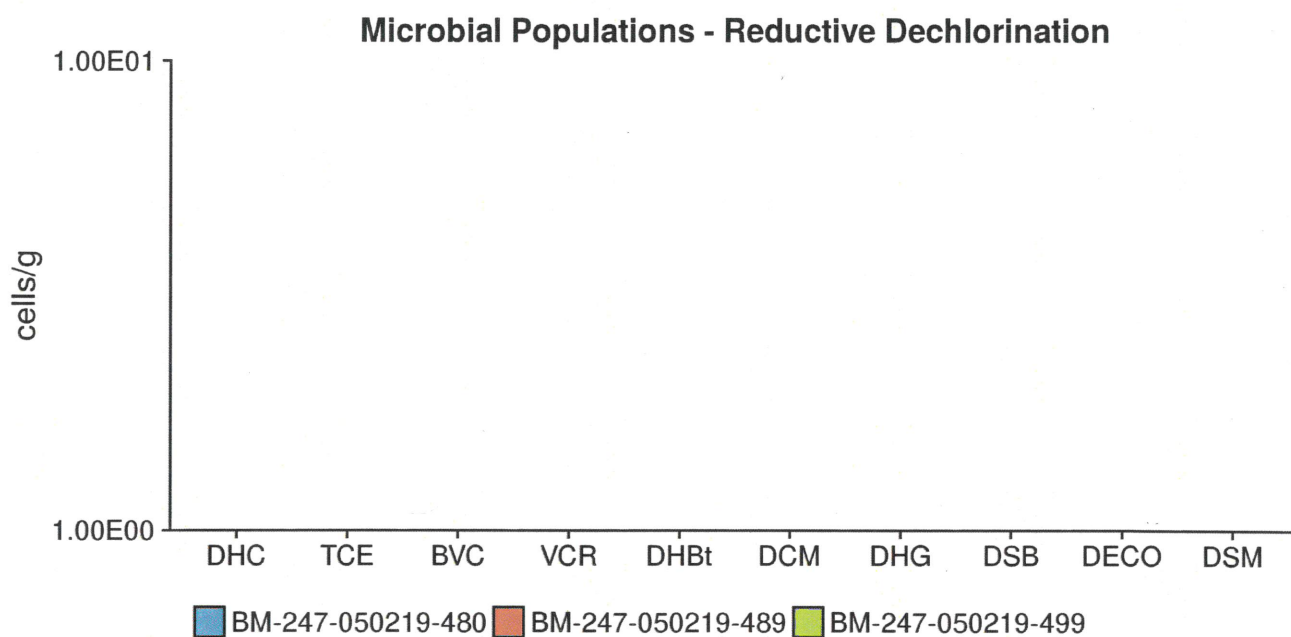


Figure 16: Comparison - microbial populations involved in reductive dechlorination.

Table 9: Summary of the QuantArray®-Chlor results for microorganisms responsible for reductive dechlorination for samples BM-247-050219-480, BM-247-050219-489, and BM-247-050219-499.

Sample Name	BM-247-050219-480	BM-247-050219-489	BM-247-050219-499
Sample Date	02/05/2019	02/05/2019	02/05/2019
Reductive Dechlorination	cells/g	cells/g	cells/g
Chloroform Reductase (CFR)	<1.00E+04	<1.00E+04	<1.00E+04
1,1 DCA Reductase (DCA)	<1.00E+04	<1.00E+04	<1.00E+04
1,2 DCA Reductase (DCAR)	<1.00E+04	<1.00E+04	<1.00E+04
PCE Reductase (PCE-1)	<1.00E+04	<1.00E+04	<1.00E+04
PCE Reductase (PCE-2)	<1.00E+04	<1.00E+04	<1.00E+04
Dehalogenimonas trans-1,2-DCE Reductase (TDR)	<1.00E+04	<1.00E+04	<1.00E+04
Dehalogenimonas cerA Reductase (CER)	<1.00E+04	<1.00E+04	<1.00E+04

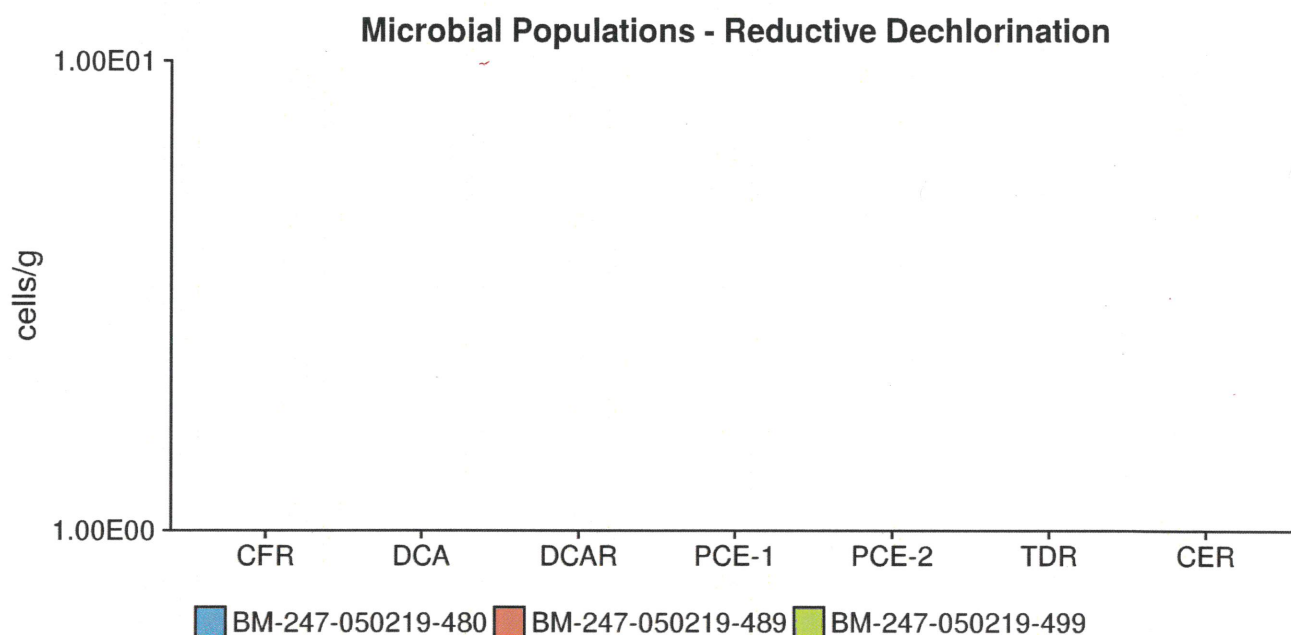


Figure 17: Comparison - microbial populations involved in reductive dechlorination.

Table 10: Summary of the QuantArray®-Chlor results for microorganisms responsible for aerobic (co)metabolism for samples BM-S8-180119-475, BM-S8-180119-499, BM-S7-220119-469, BM-S7-220119-485, and BM-S7-220119-495.

Sample Name	BM-S8-180119-475	BM-S8-180119-499	BM-S7-220119-469	BM-S7-220119-485	BM-S7-220119-495
Sample Date	01/18/2019	01/18/2019	01/22/2019	01/22/2019	01/22/2019
<i>Aerobic (Co)Metabolic</i>	cells/g	cells/g	cells/g	cells/g	cells/g
Soluble Methane Monooxygenase (SMMO)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Dioxygenase (TOD)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Phenol Hydroxylase (PHE)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Trichlorobenzene Dioxygenase (TCBO)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Monooxygenase 2 (RDEG)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Monooxygenase (RMO)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Ethene Monooxygenase (EtnC)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Epoxyalkane Transferase (EtnE)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04
Dichloromethane Dehalogenase (DCMA)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04

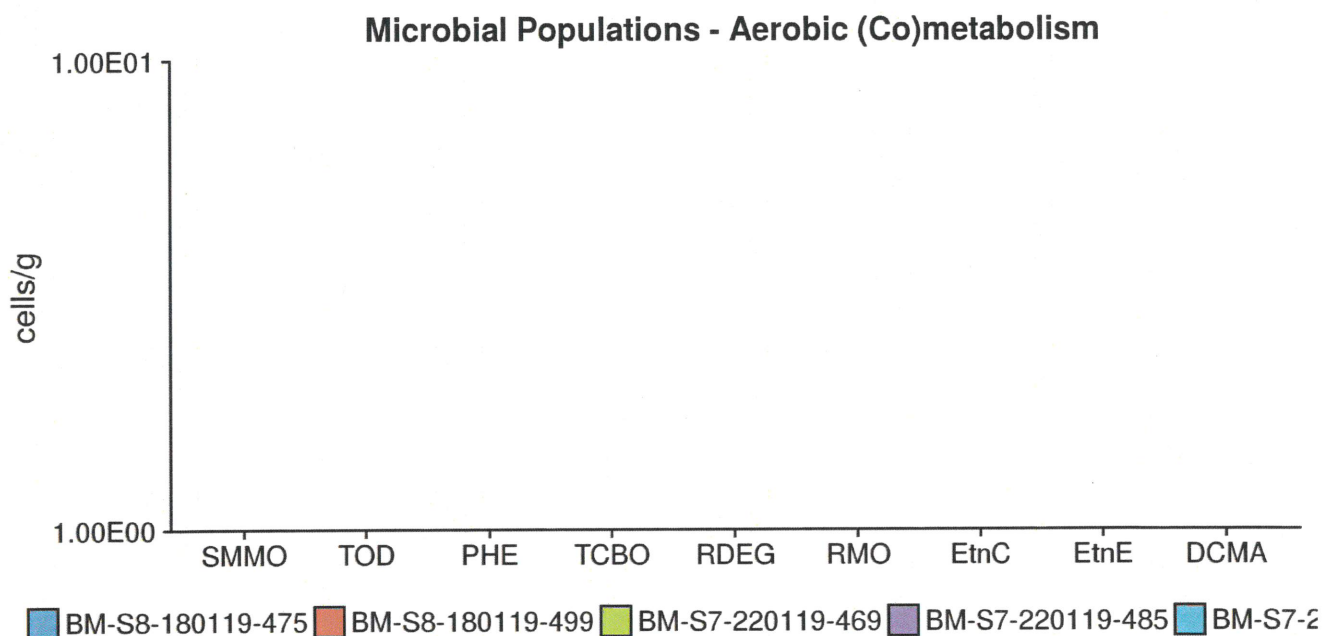


Figure 18: Comparison - microbial populations involved in aerobic (co)metabolism.

Table 11: Summary of the QuantArray®-Chlor results for microorganisms responsible for aerobic (co)metabolism for samples BM-247-300119-143, BM-247-310119-208, and BM-247-040219-474.

Sample Name	BM-247-300119-143	BM-247-310119-208	BM-247-040219-474
Sample Date	01/30/2019	01/31/2019	02/04/2019
<i>Aerobic (Co)Metabolic</i>	cells/g	cells/g	cells/g
Soluble Methane Monooxygenase (SMMO)	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Dioxygenase (TOD)	<1.00E+04	<1.00E+04	<1.00E+04
Phenol Hydroxylase (PHE)	<1.00E+04	<1.00E+04	<1.00E+04
Trichlorobenzene Dioxygenase (TCBO)	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Monooxygenase 2 (RDEG)	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Monooxygenase (RMO)	<1.00E+04	<1.00E+04	<1.00E+04
Ethene Monooxygenase (EtnC)	<1.00E+04	<1.00E+04	<1.00E+04
Epoxyalkane Transferase (EtnE)	<1.00E+04	<1.00E+04	<1.00E+04
Dichloromethane Dehalogenase (DCMA)	<1.00E+04	<1.00E+04	<1.00E+04

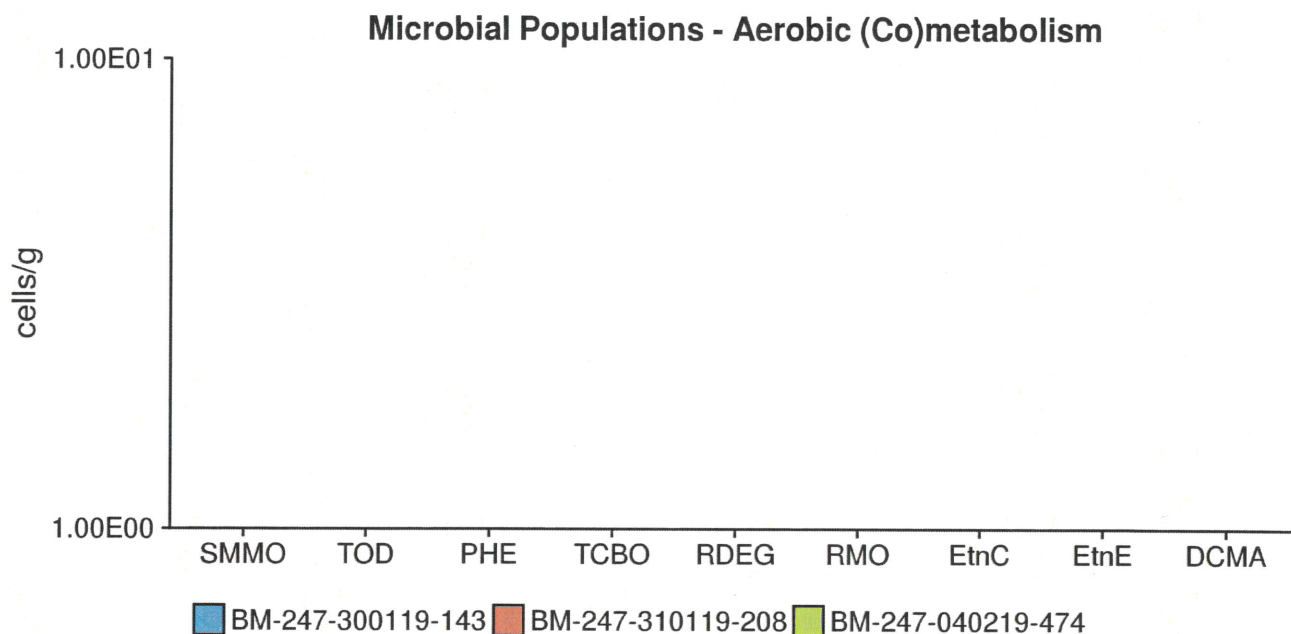


Figure 19: Comparison - microbial populations involved in aerobic (co)metabolism.

Table 12: Summary of the QuantArray®-Chlor results for microorganisms responsible for aerobic (co)metabolism for samples BM-247-050219-480, BM-247-050219-489, and BM-247-050219-499.

Sample Name	BM-247-050219-480	BM-247-050219-489	BM-247-050219-499
Sample Date	02/05/2019	02/05/2019	02/05/2019
<i>Aerobic (Co)Metabolic</i>	cells/g	cells/g	cells/g
Soluble Methane Monooxygenase (SMMO)	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Dioxygenase (TOD)	<1.00E+04	<1.00E+04	<1.00E+04
Phenol Hydroxylase (PHE)	<1.00E+04	<1.00E+04	<1.00E+04
Trichlorobenzene Dioxygenase (TCBO)	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Monooxygenase 2 (RDEG)	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Monooxygenase (RMO)	<1.00E+04	<1.00E+04	<1.00E+04
Ethene Monooxygenase (EtnC)	<1.00E+04	<1.00E+04	<1.00E+04
Epoxyalkane Transferase (EtnE)	<1.00E+04	<1.00E+04	<1.00E+04
Dichloromethane Dehalogenase (DCMA)	<1.00E+04	<1.00E+04	<1.00E+04

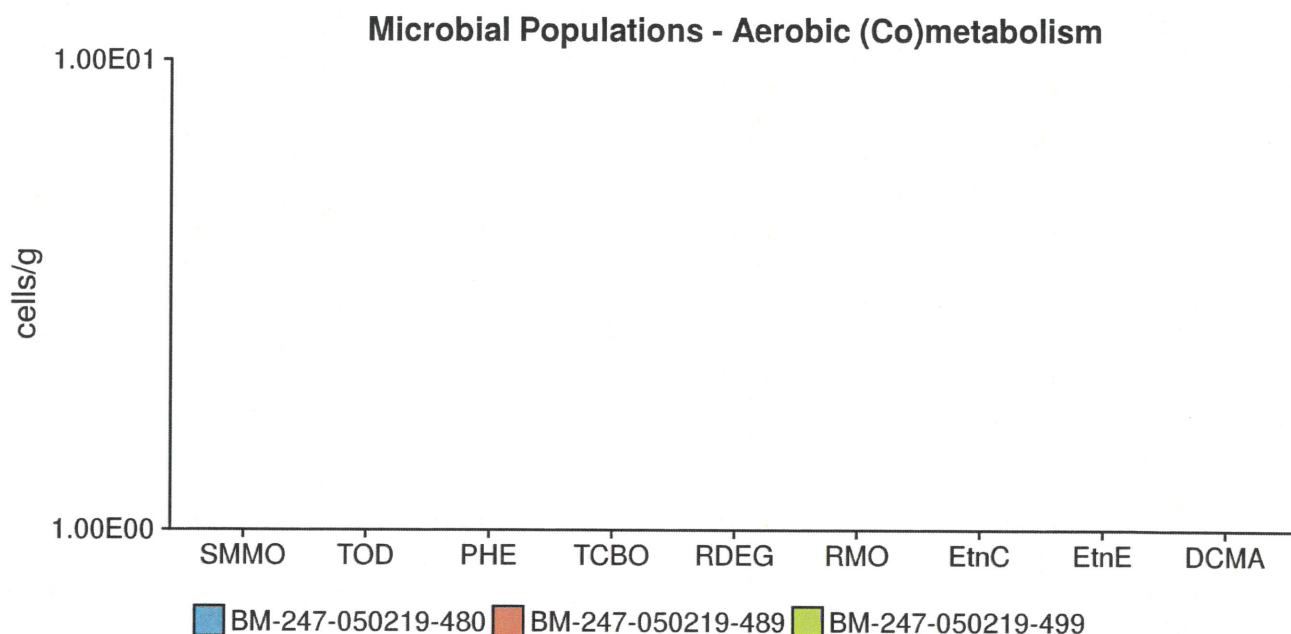


Figure 20: Comparison - microbial populations involved in aerobic (co)metabolism.

Table 13: Summary of the QuantArray® results for total bacteria and other populations for samples BM-S8-180119-475, BM-S8-180119-499, BM-S7-220119-469, BM-S7-220119-485, and BM-S7-220119-495.

Sample Name	BM-S8-180119-475	BM-S8-180119-499	BM-S7-220119-469	BM-S7-220119-485	BM-S7-220119-495
Sample Date	01/18/2019	01/18/2019	01/22/2019	01/22/2019	01/22/2019
Other	cells/g	cells/g	cells/g	cells/g	cells/g
Total Eubacteria (EBAC)	1.41E+05 (I)	8.59E+05 (I)	1.37E+04 (I)	2.52E+04 (I)	2.58E+05 (I)
Sulfate Reducing Bacteria (APS)	<1.00E+04	8.72E+02 (J)	<1.00E+04	<1.00E+04	<1.00E+04
Methanogens (MGN)	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04	<1.00E+04

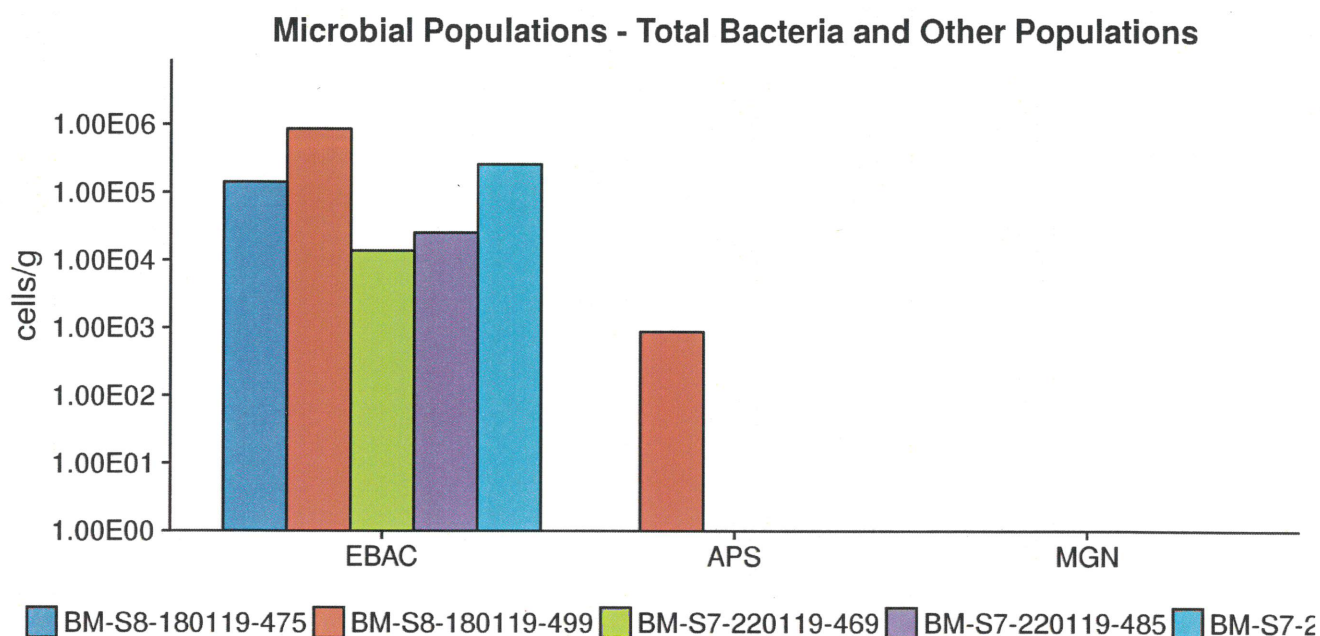


Figure 21: Comparison - microbial populations.

Table 14: Summary of the QuantArray® results for total bacteria and other populations for samples BM-247-300119-143, BM-247-310119-208, and BM-247-040219-474.

Sample Name	BM-247-300119-143	BM-247-310119-208	BM-247-040219-474
Sample Date	01/30/2019	01/31/2019	02/04/2019
Other	cells/g	cells/g	cells/g
Total Eubacteria (EBAC)	3.61E+04 (I)	5.76E+04 (I)	2.08E+03 (I)
Sulfate Reducing Bacteria (APS)	<1.00E+04	<1.00E+04	<1.00E+04
Methanogens (MGN)	<1.00E+04	<1.00E+04	<1.00E+04

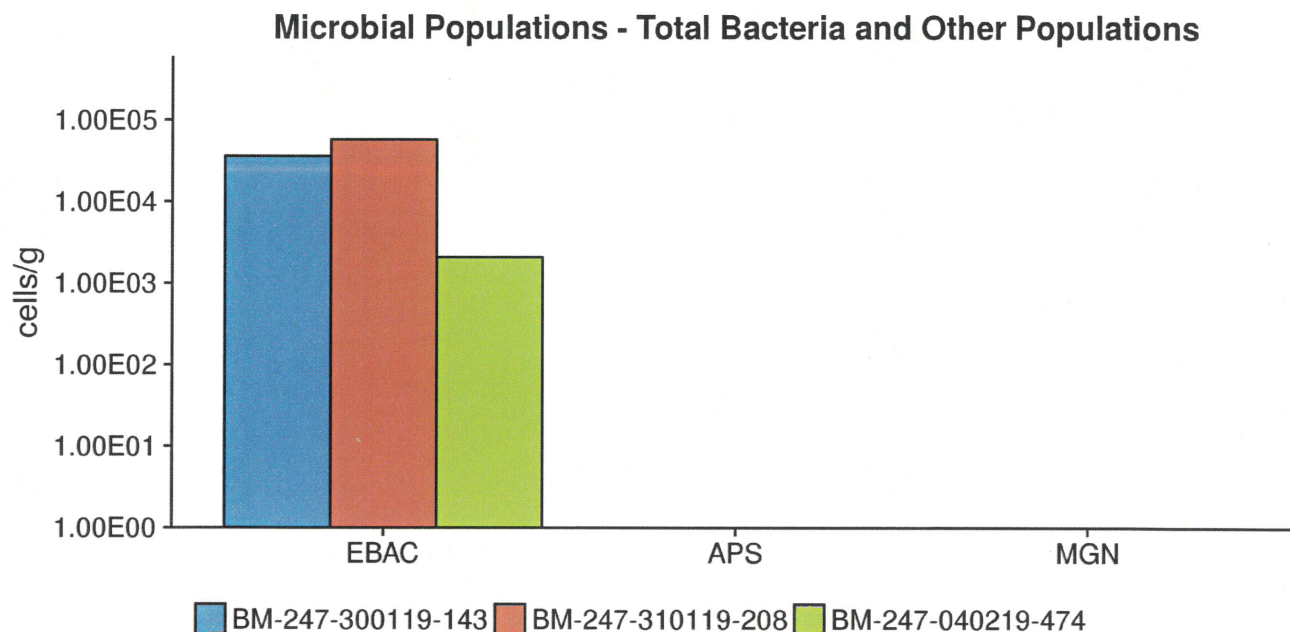


Figure 22: Comparison - microbial populations.

Table 15: Summary of the QuantArray® results for total bacteria and other populations for samples BM-247-050219-480, BM-247-050219-489, and BM-247-050219-499.

Sample Name	BM-247-050219-480	BM-247-050219-489	BM-247-050219-499
Sample Date	02/05/2019	02/05/2019	02/05/2019
Other	cells/g	cells/g	cells/g
Total Eubacteria (EBAC)	5.25E+04 (I)	1.54E+06 (I)	1.69E+04 (I)
Sulfate Reducing Bacteria (APS)	<1.00E+04	<1.00E+04	<1.00E+04
Methanogens (MGN)	<1.00E+04	<1.00E+04	<1.00E+04

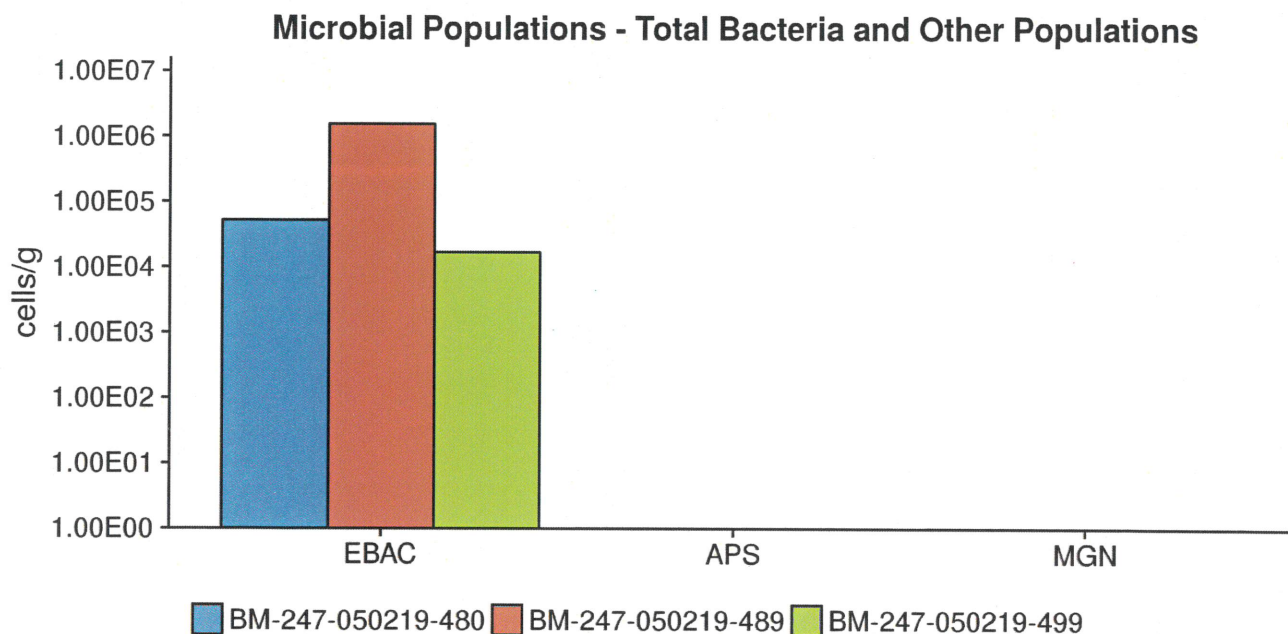


Figure 23: Comparison - microbial populations.

Interpretation

The overall purpose of the QuantArray®-Chlor is to give site managers the ability to simultaneously yet economically evaluate the potential for biodegradation of a spectrum of common chlorinated contaminants through a multitude of anaerobic and aerobic (co)metabolic pathways in order to provide a clearer and more comprehensive view of contaminant biodegradation. The following discussion describes the interpretation of results in general terms and is meant to serve as a guide.

Reductive Dechlorination - Chlorinated Ethenes: While a number of bacterial cultures including *Dehalococcoides*, *Dehalobacter*, *Desulfotobacterium*, and *Desulfuromonas* spp. capable of utilizing PCE and TCE as growth-supporting electron acceptors have been isolated [1–5], *Dehalococcoides* may be the most important because they are the only bacterial group that has been isolated to date which is capable of complete reductive dechlorination of PCE to ethene [6]. In fact, the presence of *Dehalococcoides* has been associated with complete reductive dechlorination to ethene at sites across North America and Europe [7], and Lu et al. [8] have proposed using a *Dehalococcoides* concentration of 1×10^4 cells/mL as a screening criterion to identify sites where biological reductive dechlorination is predicted to proceed at “generally useful” rates.

At chlorinated ethene sites, any “stall” leading to the accumulation of daughter products, especially vinyl chloride, would be a substantial concern. While *Dehalococcoides* concentrations greater than 1×10^4 cells/mL correspond to ethene production and useful rates of dechlorination, the range of chlorinated ethenes degraded varies by strain within the *Dehalococcoides* genus [6, 9], and the presence of co-contaminants and competitors can have complex impacts on the halo-respiring microbial community [10–15]. Therefore, QuantArray®-Chlor also provides quantification of a suite of reductive dehalogenase genes (PCE, TCE, BVC, VCR, CER, and TDR) to more definitively confirm the potential for reductive dechlorination of all chlorinated ethene compounds including vinyl chloride.

Perhaps most importantly, QuantArray®-Chlor quantifies TCE reductase (TCE) and both known vinyl chloride reductase genes (BVC, VCR) from *Dehalococcoides* to conclusively evaluate the potential for complete reductive dechlorination of chlorinated ethenes to non-toxic ethene [16–18]. In addition, the analysis also includes quantification of reductive dehalogenase genes from *Dehalogenimonas* spp. capable of reductive dechlorination of chlorinated ethenes. More specifically, these are the trans-1,2-DCE dehalogenase gene (TDR) from strain WBC-2 [19] and the vinyl chloride reductase gene (CER) from GP, the only known organisms other than *Dehalococcoides* capable of vinyl chloride reduction [20]. Finally, PCE reductase genes responsible for sequential reductive dechlorination of PCE to cis-DCE by *Sulfurospirillum* and *Geobacter* spp. are also quantified. In mixed cultures, evidence increasingly suggests that partial dechlorinators like *Sulfurospirillum* and *Geobacter* may be responsible for the majority of reductive dechlorination of PCE to TCE and cis-DCE while *Dehalococcoides* functions more as cis-DCE and vinyl chloride reducing specialists [10, 21].

Reductive Dechlorination - Chlorinated Ethanes: Under anaerobic conditions, chlorinated ethanes are susceptible to reductive dechlorination by several groups of halo-respiring bacteria including *Dehalobacter*, *Dehalogenimonas*, and *Dehalococcoides*. While the reported range of chlorinated ethanes utilized varies by genus, species, and sometimes at the strain level, several general observations can be made regarding biodegradation pathways and daughter product formation. *Dehalobacter* spp. have been isolated that are capable of sequential reductive dechlorination of 1,1,1-TCA through 1,1-DCA to chloroethane [13]. Biodegradation of 1,1,2-TCA by several halo-respiring bacteria including *Dehalobacter* and *Dehalogenimonas* spp. proceeds via dichloroelimination producing vinyl chloride [22–24]. Similarly, 1,2-DCA biodegradation by *Dehalobacter*, *Dehalogenimonas*, and *Dehalococcoides* occurs via dichloroelimination producing ethene. While not utilized by many *Desulfotobacterium* isolates, at least one strain, *Desulfotobacterium dichloroeliminans* strain DCA1, is also capable of dichloroelimination of 1,2-DCA [25]. The 1,2-dichloroethane reductive dehalogenase gene (DCAR) from members of *Desulfotobacterium* and *Dehalobacter* is known to dechlorinate 1,2-DCA to ethene, while the 1,1-dichloroethane reductive dehalogenase (DCA) targets the gene responsible for 1,1-DCA dechlorination in some strains of *Dehalobacter*. In addition to chloroform, chloroform reductase (CFR) has also been shown to be responsible for reductive dechlorination of 1,1,1-TCA [26].

Reductive Dechlorination - Chlorinated Methanes: Chloroform is a common co-contaminant at chlorinated solvent sites and can inhibit reductive dechlorination of chlorinated ethenes. Grostern et al. demonstrated that a *Dehalobacter* population was capable of reductive dechlorination of chloroform to produce dichloromethane [27]. The *cfrA* gene encodes the reductase which catalyzes this initial step in chloroform biodegradation [26]. Justicia-Leon et al. have since shown that dichloromethane can support growth of a distinct group of *Dehalobacter* strains via fermentation [28]. The *Dehalobacter* DCM assay targets the 16S rRNA gene of these strains.

Reductive Dechlorination - Chlorinated Benzenes: Chlorinated benzenes are an important class of industrial solvents and chemical intermediates in the production of drugs, dyes, herbicides, and insecticides. The physical-chemical properties of chlorinated benzenes as well as susceptibility to biodegradation are functions of their degree of chlorination and the positions of chlorine substituents. Under anaerobic conditions, reductive dechlorination of higher chlorinated benzenes including hexachlorobenzene (HCB),

pentachlorobenzene (PeCB), tetrachlorobenzene (TeCB) isomers, and trichlorobenzene (TCB) isomers has been well documented [29], although biodegradation of individual compounds and isomers varies between isolates. For example, *Dehalococcoides* strain CBDB1 reductively dechlorinates HCB, PeCB, all three TeCB isomers, 1,2,3-TCB, and 1,2,4-TCB [9, 30]. *Dehalobium chlorocoercia* DF-1 has been shown to be capable of reductive dechlorination of HCB, PeCB, and 1,2,3,5-TeCB [31]. The dichlorobenzene (DCB) isomers and chlorobenzene (CB) were considered relatively recalcitrant under anaerobic conditions. However, new evidence has demonstrated reductive dechlorination of DCBs to CB and CB to benzene [32] with corresponding increases in concentrations of *Dehalobacter* spp. [33].

Reductive Dechlorination - Chlorinated Phenols: Pentachlorophenol (PCP) was one of the most widely used biocides in the U.S. and despite residential use restrictions, is still extensively used industrially as a wood preservative. Along with PCP, the tetrachlorophenol and trichlorophenol isomers were also used as fungicides in wood preserving formulations. 2,4-Dichlorophenol and 2,4,5-TCP were used as chemical intermediates in herbicide production (e.g. 2,4-D) and chlorophenols are known byproducts of chlorine bleaching in the pulp and paper industry. While the range of compounds utilized varies by strain, some *Dehalococcoides* isolates are capable of reductive dechlorination of PCP and other chlorinated phenols. For example, *Dehalococcoides* strain CBDB1 is capable of utilizing PCP, all three tetrachlorophenol (TeCP) congeners, all six trichlorophenol (TCP) congeners, and 2,3-dichlorophenol (2,3-DCP). PCP dechlorination by strain CBDB1 produces a mixture of 3,5-DCP, 3,4-DCP, 2,4-DCP, 3-CP, and 4-CP [34]. In the same study, however, *Dehalococcoides* strain 195 dechlorinated a more narrow spectrum of chlorophenols which included 2,3-DCP, 2,3,4-TCP, and 2,3,6-TCP, but no other TCPs or PCP. Similar to *Dehalococcoides*, some species and strains of *Desulfitobacterium* are capable of utilizing PCP and other chlorinated phenols. *Desulfitobacterium hafniense* PCP-1 is capable of reductive dechlorination of PCP to 3-CP [35]. However, the ability to biodegrade PCP is not universal among *Desulfitobacterium* isolates. *Desulfitobacterium* sp. strain PCE1 and *D. chlororespirans* strain Co23, for example, can utilize some TCP and DCP isomers, but not PCP for growth [2, 36].

Reductive Dechlorination - Chlorinated Propanes: *Dehalogenimonas* is a recently described bacterial genus of the phylum Chloroflexi which also includes the well-known chloroethene-respiring *Dehalococcoides* [23]. The *Dehalogenimonas* isolates characterized to date are also halo-respiring bacteria, but utilize a rather unique range of chlorinated compounds as electron acceptors including chlorinated propanes (1,2,3-TCP and 1,2-DCP) and a variety of other vicinally chlorinated alkanes including 1,1,2,2-tetrachloroethane, 1,1,2-trichloroethane, and 1,2-dichloroethane [23].

Aerobic - Chlorinated Ethene Cometabolism: Under aerobic conditions, several different types of bacteria including methane-oxidizing bacteria (methanotrophs), and many benzene, toluene, ethylbenzene, xylene, and (BTEX)-utilizing bacteria can cometabolize or co-oxidize TCE, DCE, and vinyl chloride [37]. In general, cometabolism of chlorinated ethenes is mediated by monooxygenase enzymes with "relaxed" specificity that oxidize a primary (growth supporting) substrate (e.g. methane) and co-oxidize the chlorinated compound (e.g. TCE). QuantArray[®]-Chlor provides quantification of a suite of genes encoding oxygenase enzymes capable of co-oxidation of chlorinated ethenes including soluble methane monooxygenase (sMMO). Soluble methane monooxygenases co-oxidize a broad range of chlorinated compounds [38–41] including TCE, *cis*-DCE, and vinyl chloride. Furthermore, soluble methane monooxygenases are generally believed to support greater rates of aerobic cometabolism [40]. QuantArray[®]-Chlor also quantifies aromatic oxygenase genes encoding ring hydroxylating toluene monooxygenase genes (RMO, RDEG), toluene dioxygenase (TOD) and phenol hydroxylases (PHE) capable of TCE co-oxidation [42–46]. TCE or a degradation product has been shown to induce expression of toluene monooxygenases in some laboratory studies [43, 47] raising the possibility of TCE cometabolism with an alternative (non-aromatic) growth substrate. Moreover, while a number of additional factors must be considered, recent research under ESTCP Project 201584 has shown positive correlations between concentrations of monooxygenase genes (soluble methane monooxygenase, ring hydroxylating monooxygenases, and phenol hydroxylase) and the rate of TCE degradation [48].

Aerobic - Chlorinated Ethane Cometabolism: While less widely studied than cometabolism of chlorinated ethenes, some chlorinated ethanes are also susceptible to co-oxidation. As mentioned previously, soluble methane monooxygenases (sMMO) exhibit very relaxed specificity. In laboratory studies, sMMO has been shown to co-oxidize a number of chlorinated ethanes including 1,1,1-TCA and 1,2-DCA [38, 40].

Aerobic - Vinyl Chloride Cometabolism: Beginning in the early 1990s, numerous microcosm studies demonstrated aerobic oxidation of vinyl chloride under MNA conditions without the addition of exogenous primary substrates. Since then, strains of

Mycobacterium, *Nocardioides*, *Pseudomonas*, *Ochrobactrum*, and *Ralstonia* species have been isolated which are capable of aerobic growth on both ethene and vinyl chloride (see Mattes et al. [49] for a review). The initial steps in the pathway are the monooxygenase (*etnABCD*) catalyzed conversion of ethene and vinyl chloride to their respective epoxyalkanes (epoxyethane and chlorooxirane), followed by epoxyalkane:CoM transferase (*etnE*) mediated conjugation and breaking of the epoxide [50].

Aerobic - Chlorinated Benzenes: In general, chlorobenzenes with four or less chlorine groups are susceptible to aerobic biodegradation and can serve as growth-supporting substrates. Toluene dioxygenase (TOD) has a relatively relaxed substrate specificity and mediates the incorporation of both atoms of oxygen into the aromatic ring of benzene and substituted benzenes (toluene and chlorobenzene). Comparison of TOD levels in background and source zone samples from a CB-impacted site suggested that CBs promoted growth of TOD-containing bacteria [51]. In addition, aerobic biodegradation of some trichlorobenzene and even tetrachlorobenzene isomers is initiated by a group of related trichlorobenzene dioxygenase genes (TCBO). Finally, phenol hydroxylases catalyze the continued oxidation and in some cases, the initial oxidation of a variety of monoaromatic compounds. In an independent study, significant increases in numbers of bacteria containing PHE genes corresponded to increases in biodegradation of DCB isomers [51].

Aerobic - Chlorinated Methanes: Many aerobic methylotrophic bacteria, belonging to diverse genera (*Hyphomicrobium*, *Methylobacterium*, *Methylophilus*, *Pseudomonas*, *Paracoccus*, and *Alibacter*) have been isolated which are capable of utilizing dichloromethane (DCM) as a growth substrate. The DCM metabolic pathway in methylotrophic bacteria is initiated by a dichloromethane dehalogenase (DCMA) gene. DCMA is responsible for aerobic biodegradation of dichloromethane by methylotrophs by first producing formaldehyde which is then further oxidized [52]. As discussed in previous sections, soluble methane monooxygenase (sMMO) exhibits relaxed specificity and co-oxidizes a broad spectrum of chlorinated hydrocarbons. In addition to chlorinated ethenes, sMMO has been shown to co-oxidize chloroform in laboratory studies [38, 41].

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SITE LOGIC Report

Abiotic Potential

Contact: Pamela J. Moss

Phone:

Address: EA Engineering- Greenwood Village, CO
7995 E. Prentice Ave, Suite 206E
Greenwood Village, CO 80111

Email: pmoss@eaest.com

MI Identifier: 053QA

Report Date: 2-8-2019

Project Name: KAFB-Bulk Fuels Fac Vadose Zone

Comments:

Overview

Although not always fully considered, abiotic degradation can be a substantial or even the primary process for chlorinated hydrocarbon destruction at sites undergoing or transitioning to monitored natural attenuation (MNA). A variety of iron-bearing minerals including iron sulfides (mackinawite and pyrite), iron oxides (magnetite), green rust, and iron-bearing clays are capable of complete or nearly complete degradation of PCE, TCE, and carbon tetrachloride (He et al. 2009). Some iron-bearing minerals also catalyze the degradation of chlorinated ethanes and the lesser chlorinated ethenes cis-dichloroethene (DCE) and vinyl chloride. While the quantities and types will vary, these reactive iron minerals are frequently identified in subsurface environments under iron reducing and sulfate reducing conditions.

Brown et al. (2007) recommend four avenues for evaluating the role of abiotic processes in contaminant attenuation. First, examining contaminant concentrations along the flow path - decreasing parent compound concentrations with no evidence of accumulation of chlorinated transformation products like cis-DCE and vinyl chloride suggest abiotic degradation. Performing compound specific isotope analysis (CSIA) or monitoring for products unique to abiotic reactions such as acetylene can also provide a strong line of evidence. Microcosm studies with native sediment and killed controls can also be performed. Finally, Brown et al. (2007) suggest performing mineralogical analyses on aquifer sediment to characterize reactive minerals such as magnetite or iron monosulfides

Magnetic Susceptibility - Magnetite

Magnetite (Fe_3O_4) is a mixed valence iron mineral shown to react with PCE, TCE, and carbon tetrachloride. Furthermore, Ferrey et al. (2004) conclusively linked the observed degradation of cis-DCE at a former ammunition plant to magnetite in the subsurface. No direct chemical test is available for quantification of magnetite. However, magnetite is the most abundant mineral in natural sediments that exhibits magnetic behavior. Therefore, magnetic susceptibility provides an inexpensive and valuable estimate of the quantity of magnetite in environmental samples.

X-ray Diffraction (XRD) - Mackinawite, Pyrite, Magnetite and Green Rust

XRD is one of the primary techniques used to identify unknown crystalline materials. Most minerals are crystalline and will scatter X-rays in a regular, characteristic manner dependent on their crystal structure.

- Mackinawite is the most reactive of the iron-bearing minerals and a crystalline form (tetragonal FeS) can be detected by XRD. Mackinawite will transform PCE and TCE primarily by elimination to acetylene. Carbon tetrachloride is transformed mainly to chloroform but carbon dioxide, formate, and carbon disulfide have also been detected. Finally, the more heavily chlorinated ethanes including hexachloroethane, pentachloroethane, and tetrachloroethanes react to form chlorinated ethenes which can be further degraded.
- Pyrite (FeS_2) catalyzes beta elimination transforming PCE, TCE, and cis-DCE to acetylene and ethene. Vinyl chloride is transformed to ethene and ethane. Pyrite is also capable of degradation of carbon tetrachloride potentially forming a number of products including chloroform, carbon dioxide, carbon disulfide, and formate depending on reaction conditions.
- While not quantitative like the magnetic susceptibility test, XRD can also detect magnetite when present at between 2% and 5% on a weight basis.
- Green rusts have been reported to transform a number of common chlorinated contaminants including cis-DCE, vinyl chloride, trichloroethanes, and tetrachloroethanes. While special sample care to prevent oxidation would be needed, XRD can be used to detect green rust.

Percent Clay

Clays have large surface areas, balanced by exchangeable cations, which can bind a large number of both organic and inorganic molecules impacting their availability and reactivity in the subsurface. While less well studied than the other iron-bearing minerals, various phyllosilicate clays have been shown to be capable of degradation of PCE, TCE, cis-DCE, vinyl chloride, and carbon tetrachloride

Sample Location	BM-S8-180119-475	BM-S8-180119-499	BM-S7-220119-469
Date Sampled	01/18/2019	01/18/2019	01/22/2019

Magnetic Susceptibility Analysis

Magnetic Susceptibility (m3/kg)	$3.0E-6 \pm 8.0E-8$	$2.1E-6 \pm 10.0E-8$	$2.0E-6 \pm 9.8E-8$
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Sample Location	BM-S7-220119-485	BM-S7-220119-495	BM-247-300119-143
Date Sampled	01/22/2019	01/22/2019	01/30/2019

Magnetic Susceptibility Analysis

Magnetic Susceptibility (m3/kg)	$2.8E-6 \pm 1.3E-7$	$1.7E-6 \pm 3.6E-8$	$8.8E-7 \pm 1.1E-8$
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Sample Location	BM-247-310119-208	BM-247-040219-474	BM-247-050219-480
Date Sampled	01/31/2019	02/04/2019	02/05/2019

Magnetic Susceptibility Analysis

Magnetic Susceptibility (m3/kg)	$2.0E-6 \pm 9.3E-8$	$2.3E-6 \pm 5.3E-8$	$1.5E-6 \pm 2.8E-9$
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Sample Location	BM-247-050219-489	BM-247-050219-499
Date Sampled	02/05/2019	02/05/2019

Magnetic Susceptibility Analysis

Magnetic Susceptibility (m3/kg)	$1.5E-6 \pm 1.8E-8$	$2.6E-6 \pm 4.2E-8$
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* Analysis performed in triplicate and results reported as the mean followed by +/- standard deviation.

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Abiotic Reactions of Chlorinated Compounds with Iron Bearing Minerals and Zero Valent Iron (ZVI). Summaries for iron bearing minerals are based on He et al. (2009) and references therein. He et al. available at <http://nepis.epa.gov/>. Summary of ZVI based on Liu et al. (2005) and Song et al. (2005).

Contaminant	Mineral	Degradation	Reported Degradation Intermediates and Products ¹
PCE	FeS	Yes	Acetylene, TCE, cis-DCE, 1,1-DCE, ethene
	Pyrite	Yes	TCE, acetylene, ethene
	Magnetite	Yes	Unknown ²
TCE	³ GR(SO ₄)	Reports Differ	
	phyllosilicate clays	Yes	TCE, 1,1-DCE, vinyl chloride, 1,1,2-TCA, 1,1-DCA, chloroacetylene, acetylene, ethene, ethane
	ZVI	Yes	Ethene and ethane
cis-DCE	FeS	Yes	Acetylene, cis-DCE, vinyl chloride, 1,1-DCE
	Pyrite	Yes	Acetylene, ethene, cis-DCE, (organic acids with DO present)
	Magnetite	Yes	Unknown ¹
	GR(SO ₄), GR(CO ₃)	No	Only observed degradation when Cu(II) added
	phyllosilicate clays	Yes	cDCE, vinyl chloride, acetylene, ethene, ethane
	ZVI	Yes	Ethane, ethene, acetylene with minor amounts of DCE, VC depending on conditions
		No	None detected
Vinyl chloride	FeS	No	None detected
	Pyrite	Yes	Acetylene, ethene
	Magnetite	Yes	Unknown ²
	GR(SO ₄)	Yes	
	phyllosilicate clays	Yes	
	ZVI	Yes	Primarily acetylene and ethene but also much lesser amounts of ethane and VC and traces of methane, propane, propene, butane and butene
Vinyl chloride	FeS	Unknown	
	Pyrite	Yes	Ethene, ethane
	Magnetite	Yes	Unknown ²
	GR(SO ₄)	Yes	
	phyllosilicate clays	Yes	
	ZVI	Yes	Ethene, ethane, (no evidence of acetylene)

Contaminant	Mineral	Degradation	Reported Degradation Intermediates and Products ¹
1,1-DCA	FeS	Not Significant	None detected
1,1-DCA	GR(SO ₄)	Low conversion	Ethene and ethane (w/ Cu or Ag)
1,1-DCA	ZVI	Yes (low)	Ethane
1,2-DCA	FeS	Not Significant	None detected
1,2-DCA	FeS (Biogenic)	Yes	Not monitored
1,2-DCA	GR(SO ₄)	No	
1,2-DCA	ZVI	No	
1,1,1-TCA	FeS	Yes	1,1-DCA, ethene, 2-butyne
1,1,1-TCA	GR(SO ₄)	Yes	1,1-DCA, CA, ethene ethane
1,1,1-TCA	ZVI	Yes	1,1-DCA, ethane
1,1,2-TCA	FeS	Rate not significant	Small amounts of 1,1-DCE and vinyl chloride but rate not significant
1,1,2-TCA	GR(SO ₄)	Yes	Vinyl chloride, 1,1-DCE, ethene, ethane
1,1,2-TCA	ZVI	Yes	Ethane
1,1,1,2-TeCA	FeS	Yes	1,1-DCE
1,1,1,2-TeCA	GR(SO ₄)	Yes	1,1-DCE and minor (<1%) vinyl chloride, ethene, ethane
1,1,1,2-TeCA	phyllosilicate clays	Yes	1,1-DCE
1,1,1,2-TeCA	ZVI	Yes	TCE, 1,1-DCE
1,1,2,2-TeCA	FeS	Yes	TCE, cis-DCE, trans-DCE, acetylene
1,1,2,2-TeCA	GR(SO ₄)	Yes	TCE (major), cis-DCE, trans-DCE
1,1,2,2-TeCA	phyllosilicate clays	Yes	TCE
1,1,2,2-TeCA	ZVI	Yes	TCE, trans-DCE, cis-DCE
Carbon Tetrachloride	FeS	Yes	Chloroform, carbon disulfide, possibly methane, ethene, ethane
CT	Pyrite	Yes	Chloroform, CO ₂ , carbon disulfide, formate (highly dependent on conditions)
CT	Magnetite	Yes	Chloroform, carbon monoxide, methane, formate (highly dependent on conditions)
CT	GR(SO ₄)	Yes	Chloroform and hexachloroethane; Chloroform, DCM, methane, ethene
CT	phyllosilicate clays	Yes	Chloroform
CT	ZVI	Yes	Chloroform, dichloromethane, methane (depending on conditions)

Notes: GR(SO₄) sulfate green rust. GR(CO₃) carbonate green rust. ZVI zero valent iron

¹Compilation of reported degradation products. Mass recovery of products typically low --- additional undetected and unreported products are likely. Reported reaction products or proportions of reaction products were often a function of environmental conditions.

²No published studies that identify the transformation products of PCE, TCE, cis-DCE or vinyl chloride with magnetite. Ferrey et al (2004) analyzed for products of cis-DCE dechlorination including vinyl chloride, ethene, and ethane and did not find them. If Fe²⁺ sorbed to magnetite stabilizes carbene ions, the ultimate degradation product of cis-DCE on magnetite would be CO₂.

SITE LOGIC Report

Abiotic Potential

Contact: Pamela J. Moss

Phone:

Address: EA Engineering- Greenwood Village, CO
7995 E. Prentice Ave, Suite 206E
Greenwood Village, CO 80111

Email: pmoss@eaest.com

MI Identifier: 053QA

Report Date: 1-29-2019

Project Name: KAFB-Bulk Fuels Fac Vadose Zone

Comments:

Overview

Although not always fully considered, abiotic degradation can be a substantial or even the primary process for chlorinated hydrocarbon destruction at sites undergoing or transitioning to monitored natural attenuation (MNA). A variety of iron-bearing minerals including iron sulfides (mackinawite and pyrite), iron oxides (magnetite), green rust, and iron-bearing clays are capable of complete or nearly complete degradation of PCE, TCE, and carbon tetrachloride (He et al. 2009). Some iron-bearing minerals also catalyze the degradation of chlorinated ethanes and the lesser chlorinated ethenes cis-dichloroethene (DCE) and vinyl chloride. While the quantities and types will vary, these reactive iron minerals are frequently identified in subsurface environments under iron reducing and sulfate reducing conditions.

Brown et al. (2007) recommend four avenues for evaluating the role of abiotic processes in contaminant attenuation. First, examining contaminant concentrations along the flow path - decreasing parent compound concentrations with no evidence of accumulation of chlorinated transformation products like cis-DCE and vinyl chloride suggest abiotic degradation. Performing compound specific isotope analysis (CSIA) or monitoring for products unique to abiotic reactions such as acetylene can also provide a strong line of evidence. Microcosm studies with native sediment and killed controls can also be performed. Finally, Brown et al. (2007) suggest performing mineralogical analyses on aquifer sediment to characterize reactive minerals such as magnetite or iron monosulfides

Magnetic Susceptibility - Magnetite

Magnetite (Fe_3O_4) is a mixed valence iron mineral shown to react with PCE, TCE, and carbon tetrachloride. Furthermore, Ferrey et al. (2004) conclusively linked the observed degradation of cis-DCE at a former ammunition plant to magnetite in the subsurface. No direct chemical test is available for quantification of magnetite. However, magnetite is the most abundant mineral in natural sediments that exhibits magnetic behavior. Therefore, magnetic susceptibility provides an inexpensive and valuable estimate of the quantity of magnetite in environmental samples.

X-ray Diffraction (XRD) - Mackinawite, Pyrite, Magnetite and Green Rust

XRD is one of the primary techniques used to identify unknown crystalline materials. Most minerals are crystalline and will scatter X-rays in a regular, characteristic manner dependent on their crystal structure.

- Mackinawite is the most reactive of the iron-bearing minerals and a crystalline form (tetragonal FeS) can be detected by XRD. Mackinawite will transform PCE and TCE primarily by elimination to acetylene. Carbon tetrachloride is transformed mainly to chloroform but carbon dioxide, formate, and carbon disulfide have also been detected. Finally, the more heavily chlorinated ethanes including hexachloroethane, pentachloroethane, and tetrachloroethanes react to form chlorinated ethenes which can be further degraded.
- Pyrite (FeS_2) catalyzes beta elimination transforming PCE, TCE, and cis-DCE to acetylene and ethene. Vinyl chloride is transformed to ethene and ethane. Pyrite is also capable of degradation of carbon tetrachloride potentially forming a number of products including chloroform, carbon dioxide, carbon disulfide, and formate depending on reaction conditions.
- While not quantitative like the magnetic susceptibility test, XRD can also detect magnetite when present at between 2% and 5% on a weight basis.
- Green rusts have been reported to transform a number of common chlorinated contaminants including cis-DCE, vinyl chloride, trichloroethanes, and tetrachloroethanes. While special sample care to prevent oxidation would be needed, XRD can be used to detect green rust.

Percent Clay

Clays have large surface areas, balanced by exchangeable cations, which can bind a large number of both organic and inorganic molecules impacting their availability and reactivity in the subsurface. While less well studied than the other iron-bearing minerals, various phyllosilicate clays have been shown to be capable of degradation of PCE, TCE, cis-DCE, vinyl chloride, and carbon tetrachloride

Sample Location	BM-S8-180119-475	BM-S8-180119-499	BM-S7-220119-469
Date Sampled	01/18/2019	01/18/2019	01/22/2019

Magnetic Susceptibility Analysis

Magnetic Susceptibility (m3/kg)	3.0E-6 ± 8.0E-8	2.1E-6 ± 10.0E-8	2.0E-6 ± 9.8E-8
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Sample Location	BM-S7-220119-485	BM-S7-220119-495
Date Sampled	01/22/2019	01/22/2019

Magnetic Susceptibility Analysis

Magnetic Susceptibility (m3/kg)	2.8E-6 ± 1.3E-7	1.7E-6 ± 3.6E-8
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* Analysis performed in triplicate and results reported as the mean followed by +/- standard deviation.

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Contaminant	Mineral	Degradation	Reported Degradation Intermediates and Products ¹
PCE	FeS	Yes	Acetylene, TCE, cis-DCE, 1,1-DCE, ethene
	Pyrite	Yes	TCE, acetylene, ethene
	Magnetite	Yes	Unknown ²
	³ GR(SO ₄)	Reports Differ	
	phyllosilicate clays	Yes	TCE, 1,1-DCE, vinyl chloride, 1,1,2-TCA, 1,1-DCA, chloroacetylene, acetylene, ethene, ethane
TCE	ZVI	Yes	Ethene and ethane
	FeS	Yes	Acetylene, cis-DCE, vinyl chloride, 1,1-DCE
	Pyrite	Yes	Acetylene, ethene, cis-DCE, (organic acids with DO present)
	Magnetite	Yes	Unknown ¹
	GR(SO ₄), GR(CO ₃)	No	Only observed degradation when Cu(II) added
cis-DCE	phyllosilicate clays	Yes	cDCE, vinyl chloride, acetylene, ethene, ethane
	ZVI	Yes	Ethane, ethene, acetylene with minor amounts of DCE, VC depending on conditions
	FeS	No	None detected
	Pyrite	Yes	Acetylene, ethene
	Magnetite	Yes	Unknown ²
Vinyl chloride	GR(SO ₄)	Yes	
	phyllosilicate clays	Yes	
	ZVI	Yes	Primarily acetylene and ethene but also much lesser amounts of ethane and VC and traces of methane, propane, propene, butane and butene
	FeS	Unknown	
	Pyrite	Yes	Ethene, ethane
	Magnetite	Yes	Unknown ²
	GR(SO ₄)	Yes	
	phyllosilicate clays	Yes	
	ZVI	Yes	Ethene, ethane, (no evidence of acetylene)

Contaminant	Mineral	Degradation	Reported Degradation Intermediates and Products ¹
1,1-DCA	FeS	Not Significant	None detected
1,1-DCA	GR(SO ₄)	Low conversion	Ethene and ethane (w/ Cu or Ag)
1,1-DCA	ZVI	Yes (low)	Ethane
1,2-DCA	FeS	Not Significant	None detected
1,2-DCA	FeS (Biogenic)	Yes	Not monitored
1,2-DCA	GR(SO ₄)	No	
1,2-DCA	ZVI	No	
1,1,1-TCA	FeS	Yes	1,1-DCA, ethene, 2-butyne
1,1,1-TCA	GR(SO ₄)	Yes	1,1-DCA, CA, ethene ethane
1,1,1-TCA	ZVI	Yes	1,1-DCA, ethane
1,1,2-TCA	FeS	Rate not significant	Small amounts of 1,1-DCE and vinyl chloride but rate not significant
1,1,2-TCA	GR(SO ₄)	Yes	Vinyl chloride, 1,1-DCE, ethene, ethane
1,1,2-TCA	ZVI	Yes	Ethane
1,1,1,2-TeCA	FeS	Yes	1,1-DCE
1,1,1,2-TeCA	GR(SO ₄)	Yes	1,1-DCE and minor (<1%) vinyl chloride, ethene, ethane
1,1,1,2-TeCA	phylosilicate clays	Yes	1,1-DCE
1,1,1,2-TeCA	ZVI	Yes	TCE, 1,1-DCE
1,1,2,2-TeCA	FeS	Yes	TCE, cis-DCE, trans-DCE, acetylene
1,1,2,2-TeCA	GR(SO ₄)	Yes	TCE (major), cis-DCE, trans-DCE
1,1,2,2-TeCA	phylosilicate clays	Yes	TCE
1,1,2,2-TeCA	ZVI	Yes	TCE, trans-DCE, cis-DCE
Carbon Tetrachloride	FeS	Yes	Chloroform, carbon disulfide, possibly methane, ethene, ethane
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SITE LOGIC Report

Abiotic Potential

Contact: Pamela J. Moss
Address: EA Engineering
7995 E. Prentice Ave
Suite 206E
Greenwood Village, CO 80111

Phone: 303-590-9143

Email: pmoss@eaest.com

MI Identifier: 053QA

Report Date: March 21, 2019

Project: KAFB-Bulk Fuels Fac Vadose Zone, 627350DM02

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Results

Table 1. Summary of X-Ray Diffraction and XRF results.

Sample Location	BM-S8-180119-475	BM-S8-180119-499	BM-S7-220119-469	BM-S7-220119-485
Sample Type	Solid	Solid	Solid	Solid
Sample Date	1/18/2019	1/18/2019	1/22/2019	1/22/2019
MI ID	053QA1	053QA2	053QA3	053QA4
X-Ray Diffraction Results				
Mineral Constituent	Relative Abundance (%)			
Quartz	46	51	52	42
Oligoclase	17	21	23	29
Microcline	27	23	17.5	17
Calcite	1	ND	ND	ND
Dolomite	ND	ND	ND	ND
Magnetite	2	2	2	3
Rutile	0.5	0.5	0.5	0.5
Heulandite	ND	ND	ND	ND
Kaolinite	<0.5	<0.5	<0.5	<0.5
Chlorite	ND	ND	ND	ND
Illite/Mica	0.5	<0.5	<0.5	0.5
Montmorillonite	6	2.5	5	8
Mixed-layered Illite/Smectite	ND	ND	ND	ND
Total	100	100	100	100
% Illite Layers in ML Illite/Smectite	ND	ND	ND	ND
X-ray Fluorescence Results				
Component	Results (mass %)			
MgO	0.3798	ND	0.2366	0.2725
Al ₂ O ₃	8.7800	7.6620	7.6373	9.3256
SiO ₂	74.3213	78.1395	77.8895	72.2035
P ₂ O ₅	1.0963	1.0902	1.0144	1.1023
Cl	ND	0.0317	ND	ND
SO ₃	0.0647	ND	0.0335	ND
K ₂ O	5.3489	5.4289	5.2647	5.6372
CaO	3.5964	2.8848	2.7783	3.8011
TiO ₂	0.6119	0.5193	0.5266	0.6671
MnO	0.1094	0.0839	0.0911	0.1336
Fe ₂ O ₃	5.2429	4.0196	4.4032	6.6252
ZnO	ND	ND	ND	0.0233
Rb ₂ O	0.0268	0.0218	0.0244	0.0332
SrO	0.1060	0.0928	0.0771	0.1027
Y ₂ O ₃	ND	0.0007	ND	ND
ZrO ₂	0.0294	0.0249	0.0233	0.0728
BaO	0.2861	ND	ND	ND

Legend: J = Estimated result below PQL but above LQL ND = Result not detected < = Below detection limit

Table 2. Summary of X-Ray Diffraction and XRF results continued.

Sample Location	BM-S7-220119-495	BM-247-300119-143	BM-247-310119-208	BM-247-040219-474
Sample Type	Solid	Solid	Solid	Solid
Sample Date	1/22/2019	1/30/2019	1/31/2019	2/4/2019
MI ID	053QA5	053QA6	053QA7	053QA8
X-Ray Diffraction Results				
Mineral Constituent	Relative Abundance (%)			
Quartz	53	46.5	52	60.5
Oligase	25	19	27	21
Microcline	17.5	9	17	14
Calcite	0.5	3	ND	ND
Dolomite	ND	ND	ND	ND
Magnetite	2	3	2	2
Rutile	ND	0.5	ND	0.5
Heulanite	ND	ND	ND	<0.5
Kaolinite	ND	1	<0.5	0.5
Chlorite	ND	ND	<0.5	<0.5
Illite/Mica	<0.5	1	0.5	0.5
Montmorillonite	2	ND	ND	ND
Mixed-layered Illite/Smectite	ND	17	1.5	1
Total	100	100	100	100
% Illite Layers in ML Illite/Smectite	ND	10%	10%	10%
X-ray Fluorescence Results				
Component	Results (mass %)			
MgO	0.2168	0.6637	ND	0.2061
Al ₂ O ₃	7.8973	10.3432	7.7642	6.8164
SiO ₂	77.8303	67.3376	79.5446	79.7081
P ₂ O ₅	1.0206	0.9950	0.8449	0.8362
Cl	ND	0.0141	0.0187	ND
SO ₃	0.0736	0.0933	ND	0.0679
K ₂ O	5.6476	5.7476	5.2114	4.7075
CaO	2.7388	4.8093	2.0557	2.3251
TiO ₂	0.3938	1.2686	0.4004	0.3771
MnO	0.0752	0.1189	0.0930	0.0792
Fe ₂ O ₃	3.9675	8.0953	3.9716	4.7442
ZnO	ND	0.0275	ND	ND
Rb ₂ O	0.0213	0.0428	0.0213	0.0180
SrO	0.0903	0.1061	0.0534	0.0756
Y ₂ O ₃	0.0027	0.0982	0.0012	ND
ZrO ₂	0.0242	0.1721	0.0197	0.0387
BaO	ND	ND	ND	ND
Cr ₂ O ₃	ND	0.0667	ND	ND

Legend: J = Estimated result below PQL but above LQL ND = Result not detected < = Below detection limit

Table 3. Summary of X-Ray Diffraction and XRF results continued.

Sample Location	BM-247-050219-480	BM-247-050219-489	BM-247-050219-499
Sample Type	Solid	Solid	Solid
Sample Date	2/5/2019	2/5/2019	2/5/2019
MI ID	053QA9	053QA10	053QA11
X-Ray Diffraction Results			
Mineral Constituent	Relative Abundance (%)		
Quartz	61	61	36
Oligoclase	20	20	37
Microcline	13	15	20
Calcite	0.5	0.5	0.5
Dolomite	0.5	ND	ND
Magnetite	2	1.5	2
Rutile	ND	ND	0.5
Heulandite	ND	ND	ND
Kaolinite	0.5	<0.5	<0.5
Chlorite	ND	ND	ND
Illite/Mica	0.5	0.5	0.5
Montmorillonite	ND	ND	3.5
Mixed-layered Illite/Smectite	2	1.5	ND
Total	100	100	100
% Illite Layers in ML Illite/Smectite	10%	10%	10%
X-ray Fluorescence Results			
Component	Results (mass %)		
MgO	0.2604	ND	0.2376
Al ₂ O ₃	6.6194	7.5282	9.8080
SiO ₂	81.0614	80.2252	72.5567
P ₂ O ₅	0.9519	0.9672	1.0858
Cl	ND	ND	ND
SO ₃	ND	0.0586	0.0628
K ₂ O	4.4687	4.8606	6.5113
CaO	2.2370	2.5981	3.4109
TiO ₂	0.4547	0.3901	0.5894
MnO	0.0603	0.0654	0.0882
Fe ₂ O ₃	3.7480	3.1866	5.4745
ZnO	ND	ND	ND
Rb ₂ O	0.0194	0.0192	0.0288
SrO	0.0728	0.0807	0.1146
Y ₂ O ₃	0.0019	ND	0.0081
ZrO ₂	0.0440	0.0201	0.0233
BaO	ND	ND	ND
Cr ₂ O ₃	ND	ND	ND

Legend: J = Estimated result below PQL but above LQL ND = Result not detected < = Below detection limit

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Abiotic Reactions of Chlorinated Compounds with Iron Bearing Minerals and Zero Valent Iron (ZVI). Summaries for iron bearing minerals are based on He et al. (2009) and references therein. He et al. available at <http://nepis.epa.gov/>. Summary of ZVI based on Liu et al. (2005) and Song et al. (2005).

Contaminant	Mineral	Degradation	Reported Degradation Intermediates and Products ¹
PCE	FeS	Yes	Acetylene, TCE, cis-DCE, 1,1-DCE, ethene
	Pyrite	Yes	TCE, acetylene, ethene
	Magnetite	Yes	Unknown ²
	³ GR(SO ₄)	Reports Differ	
	phyllosilicate clays	Yes	TCE, 1,1-DCE, vinyl chloride, 1,1,2-TCA, 1,1-DCA, chloroacetylene, acetylene, ethene, ethane
	ZVI	Yes	Ethene and ethane
TCE	FeS	Yes	Acetylene, cis-DCE, vinyl chloride, 1,1-DCE
	Pyrite	Yes	Acetylene, ethene, cis-DCE, (organic acids with DO present)
	Magnetite	Yes	Unknown ¹
	GR(SO ₄), GR(CO ₃)	No	Only observed degradation when Cu(II) added
	phyllosilicate clays	Yes	cDCE, vinyl chloride, acetylene, ethene, ethane
	ZVI	Yes	Ethane, ethene, acetylene with minor amounts of DCE, VC depending on conditions
cis-DCE	FeS	No	None detected
	Pyrite	Yes	Acetylene, ethene
	Magnetite	Yes	Unknown ²
	GR(SO ₄)	Yes	
	phyllosilicate clays	Yes	
	ZVI	Yes	Primarily acetylene and ethene but also much lesser amounts of ethane and VC and traces of methane, propane, propene, butane and butene
Vinyl chloride	FeS	Unknown	
	Pyrite	Yes	Ethene, ethane
	Magnetite	Yes	Unknown ²
	GR(SO ₄)	Yes	
	phyllosilicate clays	Yes	
	ZVI	Yes	Ethene, ethane, (no evidence of acetylene)

Contaminant	Mineral	Degradation	Reported Degradation Intermediates and Products ¹
1,1-DCA	FeS	Not Significant	None detected
1,1-DCA	GR(SO ₄)	Low conversion	Ethene and ethane (w/ Cu or Ag)
1,1-DCA	ZVI	Yes (low)	Ethane
1,2-DCA	FeS	Not Significant	None detected
1,2-DCA	FeS (Biogenic)	Yes	Not monitored
1,2-DCA	GR(SO ₄)	No	
1,2-DCA	ZVI	No	
1,1,1-TCA	FeS	Yes	1,1-DCA, ethene, 2-butyne
1,1,1-TCA	GR(SO ₄)	Yes	1,1-DCA, CA, ethene ethane
1,1,1-TCA	ZVI	Yes	1,1-DCA, ethane
1,1,2-TCA	FeS	Rate not significant	Small amounts of 1,1-DCE and vinyl chloride but rate not significant
1,1,2-TCA	GR(SO ₄)	Yes	Vinyl chloride, 1,1-DCE, ethene, ethane
1,1,2-TCA	ZVI	Yes	Ethane
1,1,1,2-TeCA	FeS	Yes	1,1-DCE
1,1,1,2-TeCA	GR(SO ₄)	Yes	1,1-DCE and minor (<1%) vinyl chloride, ethene, ethane
1,1,1,2-TeCA	phyllosilicate clays	Yes	1,1-DCE
1,1,1,2-TeCA	ZVI	Yes	TCE, 1,1-DCE
1,1,2,2-TeCA	FeS	Yes	TCE, cis-DCE, trans-DCE, acetylene
1,1,2,2-TeCA	GR(SO ₄)	Yes	TCE (major), cis-DCE, trans-DCE
1,1,2,2-TeCA	phyllosilicate clays	Yes	TCE
1,1,2,2-TeCA	ZVI	Yes	TCE, trans-DCE, cis-DCE
Carbon Tetrachloride	FeS	Yes	Chloroform, carbon disulfide, possibly methane, ethene, ethane
CT	Pyrite	Yes	Chloroform, CO ₂ , carbon disulfide, formate (highly dependent on conditions)
CT	Magnetite	Yes	Chloroform, carbon monoxide, methane, formate (highly dependent on conditions)
CT	GR(SO ₄)	Yes	Chloroform and hexachloroethane; Chloroform, DCM, methane, ethene
CT	phyllosilicate clays	Yes	Chloroform
CT	ZVI	Yes	Chloroform, dichloromethane, methane (depending on conditions)

Notes: GR(SO₄) sulfate green rust. GR(CO₃) carbonate green rust. ZVI zero valent iron

¹Compilation of reported degradation products. Mass recovery of products typically low - additional undetected and unreported products are likely. Reported reaction products or proportions of reaction products were often a function of environmental conditions.

²No published studies that identify the transformation products of PCE, TCE, cis-DCE or vinyl chloride with magnetite. Ferrey et al (2004) analyzed for products of cis-DCE dechlorination including vinyl chloride, ethene, and ethane and did not find them. If Fe²⁺ sorbed to magnetite stabilizes carbene ions, the ultimate degradation product of cis-DCE on magnetite would be CO₂.

7735 2159 5696

COI-S7-MI-20190122



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Name: P. Moss
Company: EA Engineering, Science & Technology, Inc.
Address: 7995 E Prentice Ave.
Greenwood Village CO 80111
email: p.moss@easest.com
Phone: 303-590-9143
Fax:

INVOICE TO: (For Invoices paid by a third party it is imperative that all information be provided)

Name: Dallas Accounts Payable via Email below
Company: EA Engineering, Science & Technology, Inc.
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email: dallasap@easest.com
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Fax:

10515 Research Dr
Knoxville, TN 37932
865-573-8188
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Project Manager:

Project Name: Kirtland Vadose Zone Coring
Project No.: 62735DM02.1017

Purchase Order No. 18091

Subcontract No.

MI Quote No. Q2018628.0001

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Sample Information				Analyses										CENSUS: Please select the target organism/gene									
MI ID (Laboratory Use Only)	Sample Name	Date Sampled	Time Sampled	Matrix	Total Number of Containers	PLFA	NGS	QuantArray Chlor	QuantArray Petro	DHC (Dehalococoides)	Magnetic suscept./XRD/ERD	Other:	Other:	Other:									
053QA	3 BM-S7-220119-469	01/22/19	1055	S	1																		
4	BM-S7-220119-485	01/22/19	1420	S	1																		
5	BM-S7-220119-495	01/22/19	1620	S	1																		
JR																							
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Sample Information					Analyses				CENSUS: Please select the target organism/gene																														
MI ID (Laboratory Use Only)	Sample Name	Date Sampled	Time Sampled	Matrix	Total Number of Containers	PLFA	NGS	QuantArray Chlor	QuantArray Petro	DHC (Dehalococoides)	Magnetic suscept./XRD/IERD	Other:																											
0530A1	BM-58-180119-475	01/18/19	1045	S	1	—	—	X	—	—	X																												
2	BM-58-180119-499	01/18/19	1505	S	1	—	—	X	—	—	X																												
JRL																																							

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7735 2154 5680 Coc1-297-BM-20190130



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email: p.moss@east.com
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email: dallasap@east.com
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Fax:

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Knoxville, TN 37932
865-573-8188
www.microbe.com

Project Manager:

Project Name: Bernie Bockisch
Project No.: Kirtland Vadose Zone Coring
62735DM02.1017

Purchase Order No. 18091

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MI Quote No. Q2018628.0001

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Sample Information			Analyses										CENSUS: Please select the target organism/gene									
MI ID (Laboratory Use Only)	Sample Name	Date Sampled	Time Sampled	Matrix	Total Number of Containers	PLFA	NGS	QuantArray Chlor	QuantArray Petro	DHC (Dehalococoides)	Magnetic suscept./XRD/ERD	Other:	Other:	Other:								
053QA Le	BM-143-300119-143	01/30/19	1540	S	1			X			X											
JRL																						
(The rest of the table is empty and crossed out with a large X)																						

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Date: 01/30/19 1800
Received by: J. Stettin
Date: 1/31/19 1110

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Address: 7995 E Prentice Ave.

email: p.moss@eaast.com
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Project Name: Kirtland Vadose Zone Coring
Project No.: 62735DM02.1017

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Company: EA Engineering, Science & Technology, Inc.
Address: 7995 E Prentice Ave.
Greenwood Village CO 80111
Email: p.moss@eaeast.com
Phone: 303-590-9143
Fax:

Project Manager:	Bernie Bockisch
Project Name:	Kirtland Vadose Zone Coring
Project No.:	62735DM02.1017

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SITE LOGIC Report

Next Generation Sequencing (NGS) Report

Contact: Pamela J. Moss

Phone: (303) 590-9144

Address: Greenwood Village, CO
7995 E. Prentice Ave
Suite 206E
Greenwood Village, CO 80111

Email: pmoss@eaest.com

MI Identifier: 046PK

Report Date: 12/18/2018

Project: Kirtland Vandose Zone Coring
Comments:

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Sample Overview

Table 1: Sample information for 046PK.

MI Identifier	Sample Name	Sample Date	Reads Passing Quality Filtering	% Reads Classified to Genus
046PK-1	NGS239-091118-01	11/9/2018	578,080	89.2%
046PK-2	NGS239-091118-02	11/9/2018	732,567	86.3%

Table 2: Genus diversity indices for 046PK. Please refer to the Interpretation section for more information on what these diversity indices mean.

MI Identifier	Sample Name	Shannon	Simpson	Chao1 Predicted Genera	Total Genera Observed
046PK-1	NGS239-091118-01	3.2	0.9	600	525
046PK-2	NGS239-091118-02	3.8	0.94	630	567

Results for NGS239-091118-01

Table 3: Sequencing Statistics for NGS239-091118-01

Total Reads	Reads Passing Quality Filtering	% Reads Passing Quality Filtering
578,080	578,080	100.0%

Table 4: Classification Rate Summary NGS239-091118-01

Taxonomic Level	Reads Classified to Taxonomic Level	% Total Reads Classified to Taxonomic Level
Kingdom	575,871	99.6%
Phylum	568,392	98.3%
Class	560,237	96.9%
Order	546,246	94.5%
Family	538,077	93.1%
Genus	515,453	89.2%
Species	325,098	56.2%

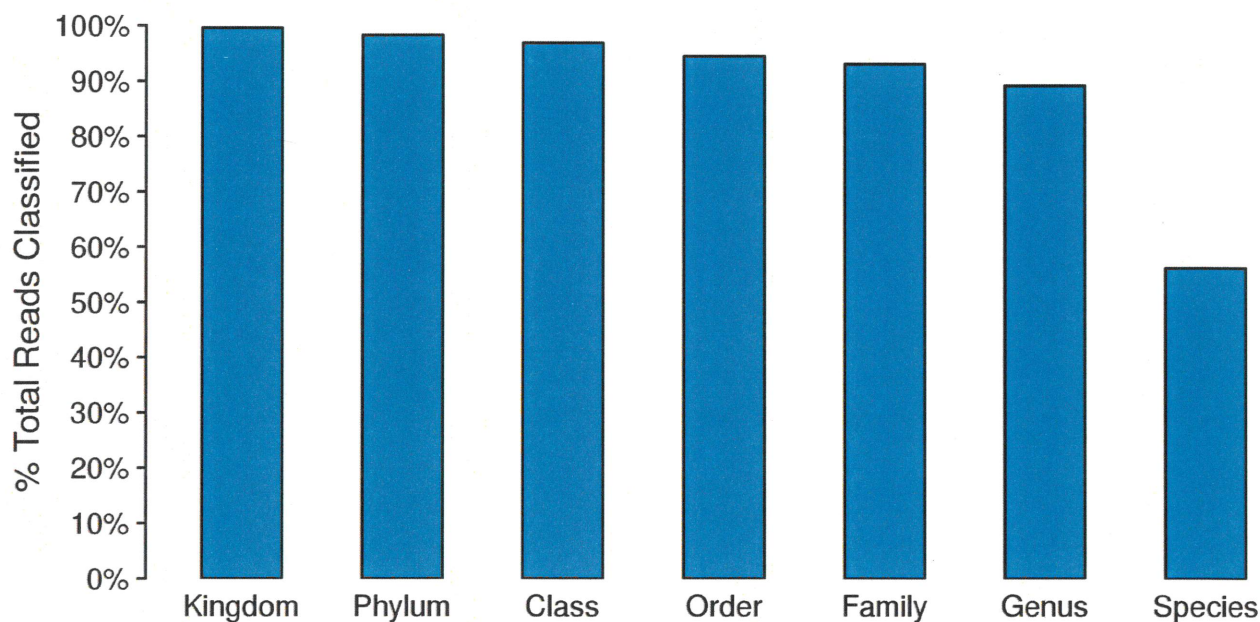


Figure 1: Classification Rate by Taxonomic Level for NGS239-091118-01

NGS239-091118-01 Classification Results by Taxonomic Level

Table 5: Top phyla classification results for NGS239-091118-01. Additional identified phyla are included in the accompanying Excel data file.

Phylum	Reads	Percent
Proteobacteria	311,509	54.8%
Bacteroidetes	162,612	28.6%
Firmicutes	69,121	12.2%
Actinobacteria	6,822	1.2%
Chlamydiae	6,149	1.1%
Tenericutes	3,544	0.6%
Cyanobacteria	1,635	0.3%
Spirochaetes	1,574	0.3%

NGS239-091118-01 Top Phyla

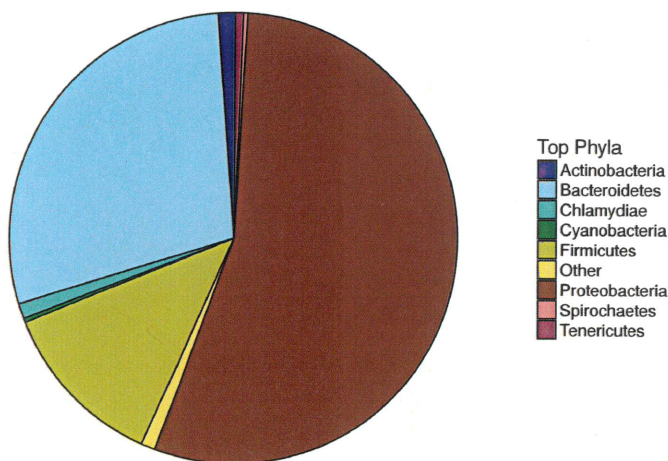


Figure 2: Top Phylum Classification Results for NGS239-091118-01

Table 6: Top genera classification results for sample NGS239-091118-01. Additional identified genera are included in the accompanying Excel data file.

Genus	Reads	Percent	Description
Runella	128,638	25.0%	Species of this genus are metabolically diverse and generally aerobic. Acids are produced either fermentatively or oxidatively depending on the species. Two species have been isolated from phosphorus-removing activated sludge.
Sulfuritalea	60,920	11.8%	Genus of facultatively anaerobic sulfur-oxidizing bacteria (SOB). Thiosulfate and elemental sulfur are oxidized with hydrogen serving as the electron donor while nitrate is reduced to nitrogen. Growth can also occur heterotrophically with certain carboxylates and organic acids as carbon and energy sources. A recent study has shown that in nitrate-depleted waters, arsenate can also be utilized as an electron acceptor. Species have been isolated from freshwater environments.
Ectothiorhodospira	48,905	9.5%	Genus of purple sulfur photoautotrophic or photoheterotrophic bacteria. Hydrogen or reduced sulfur compounds are utilized as electron donors under anoxic conditions in the light. Sulfur globules are produced from the oxidation of sulfide to sulfate. Salt is required for growth. Ectothiorhodospira has been found in alkaline lakes and estuaries exposed to light and sometimes in soil.
Hydrocarboniphaga	37,244	7.2%	Genus of aerobic heterotrophs. Alkanes are used as the sole carbon and energy sources with ammonium salts or nitrate utilized as nitrogen sources. Hydrocarboniphaga is found in sediment.
Weissella	30,938	6.0%	Genus of facultatively anaerobic chemoorganotrophs. Glucose and other carbohydrates are fermented to a variety of end-products including ethanol, carbon dioxide and acetate. Lactic acid producing bacteria and slime is produced by some species from sucrose fermentation. Vitamins are required for growth. Weissella is found in fermented food and clinical samples but has also been found in petroleum samples.
Novosphingobium	23,645	4.6%	Strictly aerobic genus of chemoorganotrophs. A sub-genus of Sphingomonas. Some species reduce nitrate and degrade aromatic compounds with one species capable of polychlorophenol degradation. Novosphingobium are ubiquitous in nature.
Tindallia	11,982	2.3%	Genus of strictly anaerobic halotolerant alkaliphiles. Growth can be chemoorganoheterotrophic or dissimilatory (utilizing low molecular weight compounds). Metabolism can be fermentative or respiratory. Tindallia is an acetogenic ammonifier, converting amino acids to acetate. This genus is found in soda lakes and other high-alkalinity sediments.
Flectobacillus	11,676	2.3%	This aerobic genus is chemoorganotrophic but not methylotrophic and has a strictly respiratory metabolism with oxygen as the terminal electron acceptor. This genus is naturally found in eutrophic ponds and freshwater lakes.
Halochromatium	11,624	2.3%	Genus of purple sulfur bacteria. Elemental sulfur (as well as hydrogen) is used as an electron donor to form sulfate under anoxic conditions in the light. Growth under oxic conditions in the dark is also possible. Requiring salt for growth, Halochromatium is found in sediments and water with high salt concentrations.

Continued on next page

Table 6: Top genera classification results for sample NGS239-091118-01. Additional identified genera are included in the accompanying Excel data file.

Genus	Reads	Percent	Description
Dechloromonas	7,872	1.5%	This facultatively anaerobic genus utilize a variety of electron acceptors including chlorate, perchlorate, and oxygen. Organic acids are used as electron donors although sulfide and Fe(II) are also used for chlorate reduction. Some Dechloromonas species perform nitrate reduction coupled with hydrocarbon degradation.
Rubrivivax	7,617	1.5%	Rubrivivax is a genus of purple non-sulfur bacteria. Under anaerobic conditions in the light, Rubrivivax grows as a photoheterotroph with a variety of organic substrates or photoautotrophically using hydrogen and carbon dioxide. In the dark, chemoautotrophic growth with CO or aerobic heterotrophic growth with carbohydrates, sugars and other organics are also possible. One species, R. benzoatilyticus can degrade benzoates.
Thiobacillus	7,289	1.4%	Genus of sulfur oxidizing aerobes or facultative anaerobes. Reduced sulfur compounds are oxidized to sulfate. Thiobacillus generally grow autotrophically but some species grow chemoorganotrophically or chemolithotrophically. All species are carbon dioxide fixers. Some species are facultative denitrifiers with T. denitrificans capable of complete denitrification. Thiobacillus spp are found in a variety of environments where high levels of oxidizable sulfur compounds are present such as sulfur springs or sewage treatment plants.
Clostridium	5,651	1.1%	Obligately anaerobic genus possessing a fermentative type of metabolism. Some species are aerotolerant. Growth requirements vary greatly between species, utilizing a range of compounds. Fermentation end products are usually a combination of alcohols and organic acids. Some species are acetogens.
Candidatus Rhab- dochlamydia	5,397	1.0%	These obligate intracellular bacteria infect arthropods.
Chryseobacterium	5,390	1.0%	Most species are chemoorganotrophic obligate aerobes and have a strict respiratory metabolism with oxygen as the terminal electron acceptor. Some species have demonstrated anaerobic respiration with nitrate or fumarate serving as the terminal electron acceptor. Slightly psychrotolerant, Chryseobacterium utilize several types of carbohydrates as carbon and energy sources.
Azospirillum	5,135	1.0%	Microaerobic genus of nitrogen-fixing chemoorganotrophs. Metabolism is generally respiratory but can be fermentative also. Carboxylates and certain carbohydrates are utilized as carbon and energy sources. Some strains exhibit autotrophic growth facultatively with hydrogen. These plant growth-promoting rhizobacteria have been isolated from soil or plant material.
Mycobacterium	4,402	0.9%	Genus of mostly obligately aerobic, acid-fast-bacteria. Species are highly adaptable, utilizing many types of organic compounds for growth and amino acids, and ammonia as nitrogen sources. Carbon dioxide is required. Mycobacterium species are particularly useful in bioremediation, having been shown to degrade polycyclic aromatic hydrocarbons and pentachlorophenol as well as mineralizing vinyl chloride. Members of this genus are ubiquitous in nature and are found free-living in soil and water and in mammalian and clinical samples.

Continued on next page

Table 6: Top genera classification results for sample NGS239-091118-01. Additional identified genera are included in the accompanying Excel data file.

Genus	Reads	Percent	Description
Propionispora	4,249	0.8%	Strictly anaerobic genus of fermenters. Endospores are formed. Only certain sugar alcohols and fructose are utilized for fermentation. Propionic acid, acetic acid, carbon dioxide and a small amount of molecular hydrogen are the major end products.
Sphingobium	3,641	0.7%	Strictly aerobic genus of chemoorganotrophs isolated from a wastewaters and variety of contaminated soils. While substrates differ greatly between species, strains of some characterized species are capable of utilizing aromatic compounds such as phenanthrene and chlorinated aromatics including chlorophenols.
Mycoplasma	3,457	0.7%	Pleomorphic genus of aerobic or facultatively anaerobic chemoorganotrophs. Sugars or arginine are used as the main energy sources and sterols are required for growth. Mycoplasma is obligately commensal being found in eukaryotic cells.
Bdellovibrio	3,289	0.6%	Aerobic genus of predatory bacteria. Bdellovibrio invade and consume gram-negative bacteria and are sometimes associated with biofilms. This organism degrades the biofilms of <i>S. aureus</i> , in particular, by obtaining nutrients and secreting proteolytic enzymes which then continues the breakdown cycle. Bdellovibrio is found in soil, sewage and fresh and salt water.
Methyloversatilis	3,266	0.6%	Genus of facultative methylotrophic bacteria. In addition to C-1 carbon compounds, many other multi-carbon compounds such as sugars and organic acids are utilized for growth. Species have been isolated from soil and hot springs.
Plesiocystis	3,149	0.6%	Strictly aerobic genus of Myxobacteria known for fruiting body formation. In general, Myxobacteria move by gliding and grow in cell communities called swarms. Myxobacteria are thought to survive from the degradation products of biopolymers such as proteins, cellulose, and other cellular constituents. The type species, <i>P. pacifica</i> produces a haloalkane dehalogenase enzyme useful in bioremediation efforts. This halophile is found in marine environments.
Burkholderia	3,119	0.6%	A genus of metabolically versatile bacteria capable of utilizing a broad spectrum of carbon and energy sources. Many are strict aerobes, but some species are also capable of nitrate reduction. Some species have been shown to degrade aromatic hydrocarbons.
Sphingobacterium	3,059	0.6%	This aerobic genus of chemoorganotrophic bacteria is metabolically diverse. No special growth factors are required. Carbohydrates are utilized as substrates producing acids. Some species have been known to degrade uncommon compounds such as herbicide, KDN, and pentachlorophenol.
Chondromyces	3,000	0.6%	These slime-producing bacteria which feed on insoluble organic compounds, referred to as myxobacteria can create a variety of biologically active substances, including anti-fungal metabolites. Some species of Chondromyces require a symbiont for growth. In general, Myxobacteria move by gliding, are known to produce fruiting bodies during starvation conditions, are social - growing in cell communities called swarms. Myxobacteria are generally thought to live from the degradation products of biopolymers such as proteins, cellulose, and other cellular constituents.

Continued on next page

Table 6: Top genera classification results for sample NGS239-091118-01. Additional identified genera are included in the accompanying Excel data file.

Genus	Reads	Percent	Description
Kushneria	2,964	0.6%	Genus of halophilic, strictly aerobic chemoorganotrophs. A variety of carbohydrates, sugars and sugar alcohols are utilized as carbon sources and Na ⁺ is required for growth. Species of Kushneria have been isolated from hypersaline marine environments.

NGS239-091118-01 Top Genera

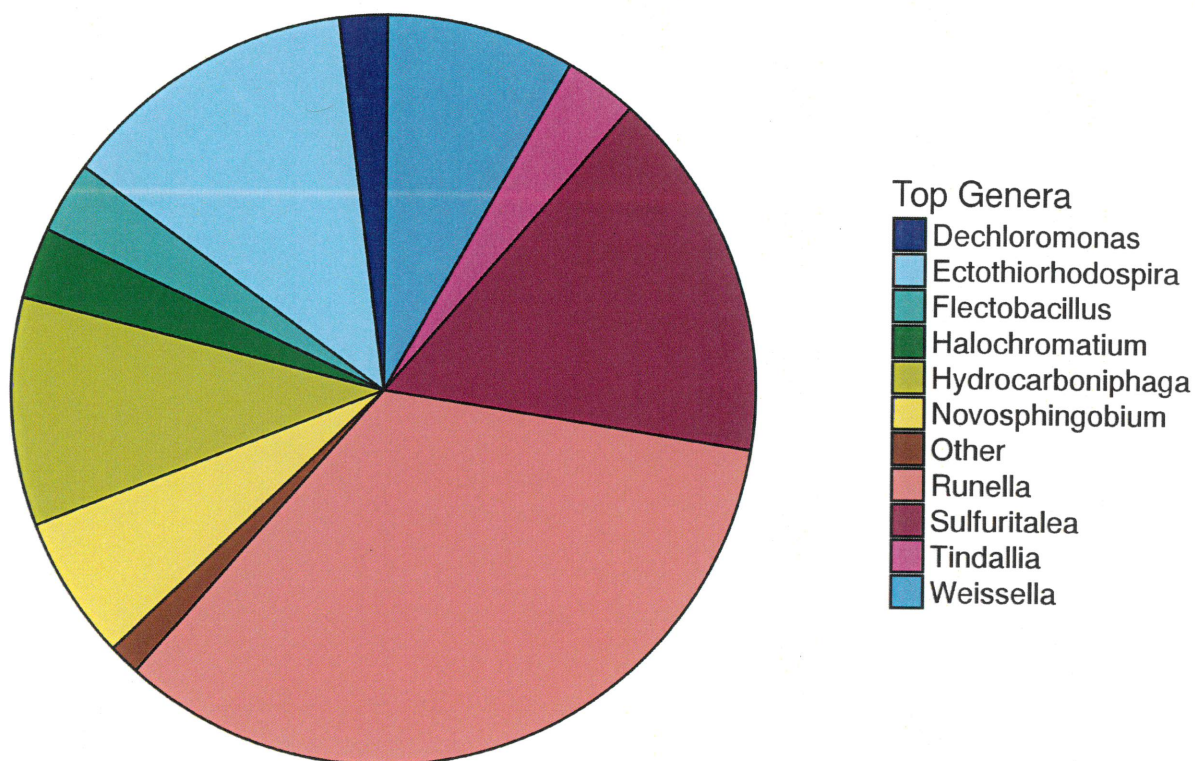


Figure 3: Top Genus Classification Results for NGS239-091118-01

Results for NGS239-091118-02

Table 7: Sequencing Statistics for NGS239-091118-02

Total Reads	Reads Passing Quality Filtering	% Reads Passing Quality Filtering
732,567	732,567	100.0%

Table 8: Classification Rate Summary NGS239-091118-02

Taxonomic Level	Reads Classified to Taxonomic Level	% Total Reads Classified to Taxonomic Level
Kingdom	730,819	99.8%
Phylum	706,572	96.5%
Class	694,269	94.8%
Order	670,376	91.5%
Family	657,657	89.8%
Genus	632,145	86.3%
Species	345,063	47.1%

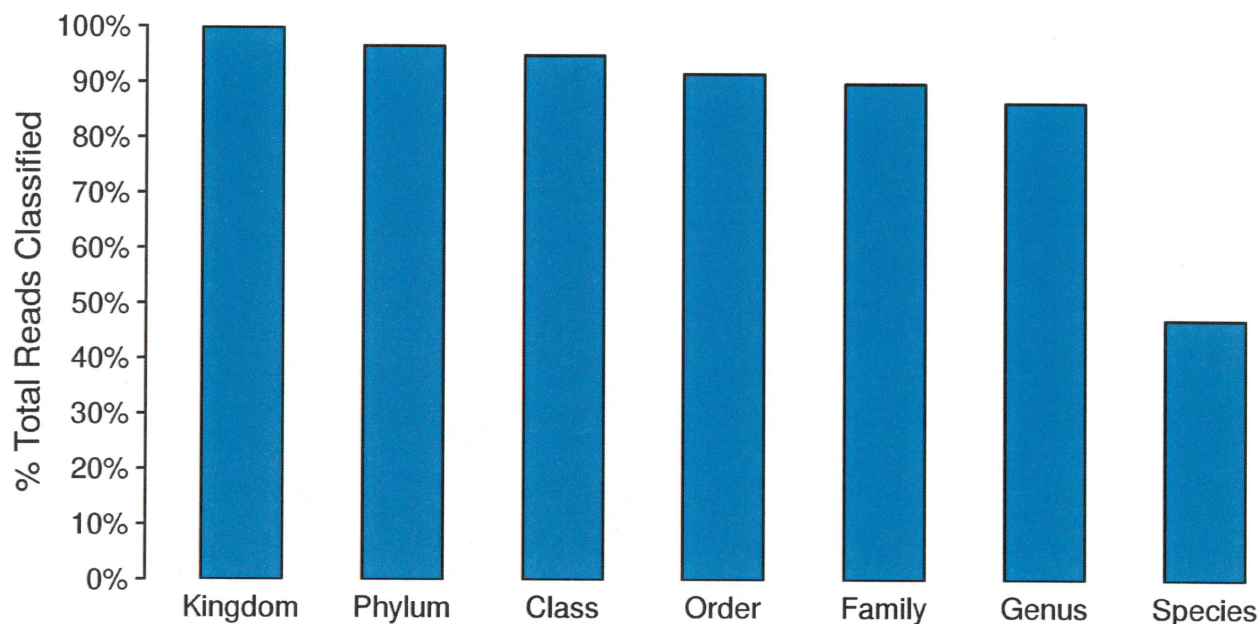


Figure 4: Classification Rate by Taxonomic Level for NGS239-091118-02

NGS239-091118-02 Classification Results by Taxonomic Level

Table 9: Top phyla classification results for NGS239-091118-02. Additional identified phyla are included in the accompanying Excel data file.

Phylum	Reads	Percent
Proteobacteria	345,452	48.9%
Firmicutes	143,748	20.3%
Bacteroidetes	126,282	17.9%
Actinobacteria	17,587	2.5%
Chlamydiae	14,840	2.1%
Cyanobacteria	12,811	1.8%
Tenericutes	8,010	1.1%
Chloroflexi	7,500	1.1%

NGS239-091118-02 Top Phyla

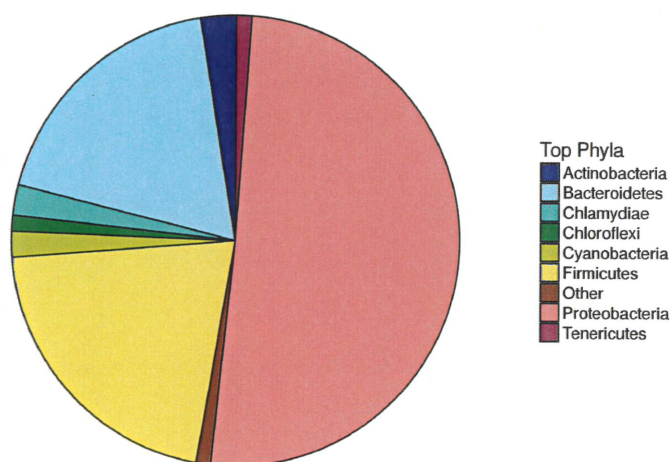


Figure 5: Top Phylum Classification Results for NGS239-091118-02

Table 10: Top genera classification results for sample NGS239-091118-02. Additional identified genera are included in the accompanying Excel data file.

Genus	Reads	Percent	Description
Runella	104,129	16.5%	Species of this genus are metabolically diverse and generally aerobic. Acids are produced either fermentatively or oxidatively depending on the species. Two species have been isolated from phosphorus-removing activated sludge.
Weissella	66,168	10.5%	Genus of facultatively anaerobic chemoorganotrophs. Glucose and other carbohydrates are fermented to a variety of end-products including ethanol, carbon dioxide and acetate. Lactic acid producing bacteria and slime is produced by some species from sucrose fermentation. Vitamins are required for growth. Weissella is found in fermented food and clinical samples but has also been found in petroleum samples.
Ectothiorhodospira	47,896	7.6%	Genus of purple sulfur photoautotrophic or photoheterotrophic bacteria. Hydrogen or reduced sulfur compounds are utilized as electron donors under anoxic conditions in the light. Sulfur globules are produced from the oxidation of sulfide to sulfate. Salt is required for growth. Ectothiorhodospira has been found in alkaline lakes and estuaries exposed to light and sometimes in soil.
Hydrocarboniphaga	41,906	6.6%	Genus of aerobic heterotrophs. Alkanes are used as the sole carbon and energy sources with ammonium salts or nitrate utilized as nitrogen sources. Hydrocarboniphaga is found in sediment.
Sulfuritalea	32,643	5.2%	Genus of facultatively anaerobic sulfur-oxidizing bacteria (SOB). Thiosulfate and elemental sulfur are oxidized with hydrogen serving as the electron donor while nitrate is reduced to nitrogen. Growth can also occur heterotrophically with certain carboxylates and organic acids as carbon and energy sources. A recent study has shown that in nitrate-depleted waters, arsenate can also be utilized as an electron acceptor. Species have been isolated from freshwater environments.
Fusibacter	30,034	4.8%	Genus of mostly anaerobic sulfur-reducing chemoorganotrophs. Some species are microaerophilic. A limited amount of carbohydrates are fermented to acetate, CO ₂ , H ₂ and butyrate. Sulfide is produced from the reduction of thiosulfate and sulfur. Fusibacter species have been found in a variety of environments including in springs and wastewater, oil producing wells and a corroded kerosene tank.
Novosphingobium	28,594	4.5%	Strictly aerobic genus of chemoorganotrophs. A sub-genus of Sphingomonas. Some species reduce nitrate and degrade aromatic compounds with one species capable of polychlorophenol degradation. Novosphingobium are ubiquitous in nature.
Xanthobacter	14,466	2.3%	Genus of aerobic slime-producing bacteria. Xanthobacter are hydrogen oxidizing bacteria. Chemoorganoheterotrophic growth can also occur with alcohols and other organic compounds as carbon sources but carbohydrate use is limited. Chemolithotrophic growth can occur when minerals are present. Characterized species are capable of CO ₂ and nitrogen fixation. Xanthobacter is ubiquitous in nature, appearing free in freshwater, marine sediments, soil and in roots of plants.

Continued on next page

Table 10: Top genera classification results for sample NGS239-091118-02. Additional identified genera are included in the accompanying Excel data file.

Genus	Reads	Percent	Description
Halochromatium	10,529	1.7%	Genus of purple sulfur bacteria. Elemental sulfur (as well as hydrogen) is used as an electron donor to form sulfate under anoxic conditions in the light. Growth under oxic conditions in the dark is also possible. Requiring salt for growth, Halochromatium is found in sediments and water with high salt concentrations.
Alkaliphilus	10,074	1.6%	Genus of strictly anaerobic alkaliphilic heterotrophic bacteria possessing both a respiratory and fermentative type of metabolism. Proteins, such as peptone, tryptone and yeast extracts are used as carbon and energy sources. A variety of electron acceptors are utilized such as elemental sulfur and thiosulfate, fumarate, crotonate, Fe(III), Co(III), or Cr (VI). Formate and acetate are the main fermentation products.
Leptolyngbya	9,197	1.5%	Genus of thin, filamentous cyanobacteria (photosynthetic bacteria) which are common and ubiquitous in soils, freshwater, marine, thermal or mineral springs, and rocks. Leptolyngbya are classified by the formation of cylindrical trichomes and can form sheaths or mucous-like membranes.
Clostridium	9,156	1.4%	Obligately anaerobic genus possessing a fermentative type of metabolism. Some species are aerotolerant. Growth requirements vary greatly between species, utilizing a range of compounds. Fermentation end products are usually a combination of alcohols and organic acids. Some species are acetogens.
Mycobacterium	9,076	1.4%	Genus of mostly obligately aerobic, acid-fast-bacteria. Species are highly adaptable, utilizing many types of organic compounds for growth and amino acids, and ammonia as nitrogen sources. Carbon dioxide is required. Mycobacterium species are particularly useful in bioremediation, having been shown to degrade polycyclic aromatic hydrocarbons and pentachlorophenol as well as mineralizing vinyl chloride. Members of this genus are ubiquitous in nature and are found free-living in soil and water and in mammalian and clinical samples.
Burkholderia	7,979	1.3%	A genus of metabolically versatile bacteria capable of utilizing a broad spectrum of carbon and energy sources. Many are strict aerobes, but some species are also capable of nitrate reduction. Some species have been shown to degrade aromatic hydrocarbons.
Mycoplasma	7,925	1.3%	Pleomorphic genus of aerobic or facultatively anaerobic chemoorganotrophs. Sugars or arginine are used as the main energy sources and sterols are required for growth. Mycoplasma is obligately commensal being found in eukaryotic cells.
Thiobacillus	7,222	1.1%	Genus of sulfur oxidizing aerobes or facultative anaerobes. Reduced sulfur compounds are oxidized to sulfate. Thiobacillus generally grow autotrophically but some species grow chemoorganotrophically or chemolithotrophically. All species are carbon dioxide fixers. Some species are facultative denitrifiers with T. denitrificans capable of complete denitrification. Thiobacillus spp are found in a variety of environments where high levels of oxidizable sulfur compounds are present such as sulfur springs or sewage treatment plants.

Continued on next page

Table 10: Top genera classification results for sample NGS239-091118-02. Additional identified genera are included in the accompanying Excel data file.

Genus		Reads	Percent	Description
Rhodobacter		6,821	1.1%	Rhodobacter includes species which possess an extensive range of metabolic capabilities. This genus of mesophilic bacteria grows photoheterotrophically or chemotrophically in the light under microaerobic conditions by fermentative or denitrifying metabolism. Rhodobacter can also grow photoautotrophically utilizing sulfur compounds as electron donors.
Flectobacillus		6,159	1.0%	This aerobic genus is chemoorganotrophic but not methylotrophic and has a strictly respiratory metabolism with oxygen as the terminal electron acceptor. This genus is naturally found in eutrophic ponds and freshwater lakes.
Candidatus tochlamydia	Pro-	6,073	1.0%	Protochlamydia is an endosymbiont, found intracellularly in the amoeba Acanthamoeba.
Treponema		5,808	0.9%	Genus of anaerobic to microaerophilic chemoorganotrophic homoacetogens. Many carbohydrates and amino acids are utilized as carbon and energy sources. Acetate is produced from carbon dioxide and hydrogen. Treponema has been isolated from mammalian fauna and from wood-feeding insects.
Candidatus dochlamydia	Rhab-	5,635	0.9%	These obligate intracellular bacteria infect arthropods.
Rubrivivax		5,137	0.8%	Rubrivivax is a genus of purple non-sulfur bacteria. Under anaerobic conditions in the light, Rubrivivax grows as a photoheterotroph with a variety of organic substrates or photoautotrophically using hydrogen and carbon dioxide. In the dark, chemoautotrophic growth with CO or aerobic heterotrophic growth with carbohydrates, sugars and other organics are also possible. One species, R. benzoatilyticus can degrade benzoates.
Methyloversatilis		4,812	0.8%	Genus of facultative methylotrophic bacteria. In addition to C-1 carbon compounds, many other multi-carbon compounds such as sugars and organic acids are utilized for growth. Species have been isolated from soil and hot springs.
Bdellovibrio		4,537	0.7%	Aerobic genus of predatory bacteria. Bdellovibrio invade and consume gram-negative bacteria and are sometimes associated with biofilms. This organism degrades the biofilms of S. aureus, in particular, by obtaining nutrients and secreting proteolytic enzymes which then continues the breakdown cycle. Bdellovibrio is found in soil, sewage and fresh and salt water.
Longilinea		4,517	0.7%	Genus of strictly anaerobic chemoheterotrophs (fermenters). Yeast extract is required for growth and in co-culture with methanogens, growth is stimulated. Longilinea species form multicellular filaments and have been isolated from rice paddy soil.
Kushneria		4,187	0.7%	Genus of halophilic, strictly aerobic chemoorganotrophs. A variety of carbohydrates, sugars and sugar alcohols are utilized as carbon sources and Na ⁺ is required for growth. Species of Kushneria have been isolated from hypersaline marine environments.
Ignavibacterium		4,037	0.6%	Moderately thermophilic genus of strictly anaerobic green sulfur bacteria. Some carbohydrates are fermented. No photosynthetic growth reported. Ignavibacterium has been found in terrestrial hot spring in microbial mats.

Continued on next page

Table 10: Top genera classification results for sample NGS239-091118-02. Additional identified genera are included in the accompanying Excel data file.

Genus	Reads	Percent	Description
Caloramator	3,474	0.5%	Fermentative genus of strictly anaerobic thermophilic chemoorganotrophs. Carbohydrates and amino acids are utilized as carbon sources. Acetate, lactate and ethanol are the main fermentation products. H ₂ and CO ₂ are also formed but is species-dependent. Caloramator have been isolated from hot springs and other heated environments.
Methylophaga	3,384	0.5%	Methylophaga are a restricted type of methylotroph mineralizing one carbon (C1) compounds such as methanol but do not grow on methane. In addition to methanol, most species also grow on methylamine. Some species also grow on dimethylsulfide with some strains also capable of chemolithoheterotrophic growth oxidizing sulfide to thiosulfate.

NGS239-091118-02 Top Genera

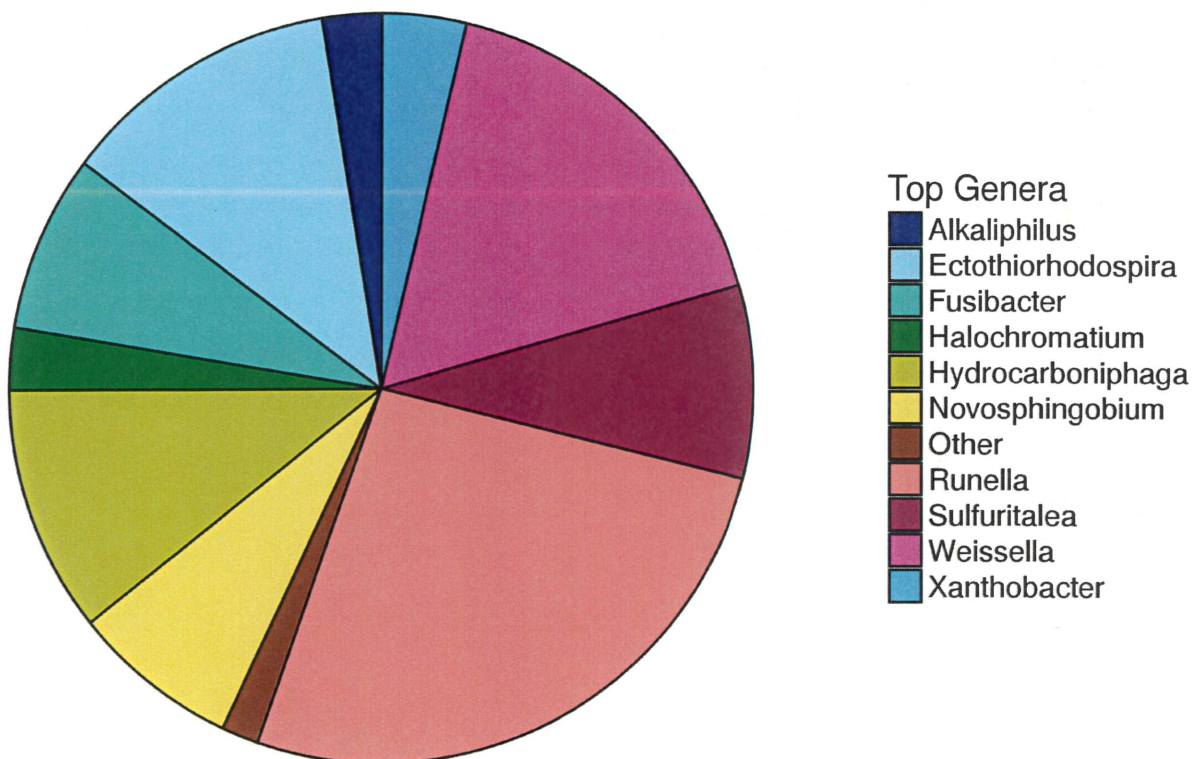


Figure 6: Top Genus Classification Results for NGS239-091118-02

Interpretation

Diversity Indices

The Shannon diversity index is a quantitative measurement that characterizes how many different genera are present in the sample and takes into account the distribution of the number of organisms classified to each genus present in the sample (commonly referred to as species evenness) [1, 2]. Shannon's diversity index increases in value as the number of genera increases and as the number of organisms present per genera becomes even. Simpson's index measures the probability that two individuals selected randomly from the sample would belong to different genera: the greater the value, the greater the sample diversity. The Chao1 index is an excellent indicator of species richness and is based on the number of reads when one (singleton) or two (doubleton) operational taxonomic units (OTUs) are observed. This value is the predicted number of genera based on the number of singletons and doubletons. The total genera observed is presented here, but does not include reads unclassified at genus species.

Principal Coordinate Analysis

Principal coordinate analysis (PCoA) is an excellent tool for visualizing differences in microbial communities between samples [3]. Unlike more traditional methods such as principal component analysis (PCA), PCoA calculates complex functions for the axes rather than dimensional scaling used in PCA. Therefore, PCoA is able to better demonstrate dissimilarities that may be nuanced in PCA tests. PCoA accomplishes this by using a dissimilarity matrix to assign each sample a location in dimensional space, then changes the coordinate system to display the data in two dimensions.

Hierarchical Clustering Dendrogram

Hierarchical clustering is accomplished by comparing dissimilarities between the samples using complete agglomeration of the Bray-Curtis dissimilarity. This groups samples which are the least dissimilar together. The length of the branches indicate the amount of dissimilarity between samples. Therefore, shorter branches are more similar. The stacked bar chart below each leaf of the tree represents the relative abundance of genus-level classifications.

References

1. Gotelli, N. J. & Colwell, R. K. Quantifying biodiversity: procedures and pitfalls in the measurement and comparison of species richness. *Ecology Letters* **4**, 379–391 (2001).
2. Hill, M. O. Diversity and evenness: a unifying notation and its consequences. *Ecology* **54**, 427–432 (1973).
3. Buttigieg, P. L. & Ramette, A. A guide to statistical analysis in microbial ecology: a community-focused, living review of multivariate data analyses. *FEMS Microbiology Ecology* **90**, 543–550. ISSN: 1574-6941 (2014).

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SITE LOGIC Report

QuantArray®-Chlor Study

Contact: Pamela J. Moss

Phone: (303) 590-9144

Address: EA Engineering- Greenwood
Village, CO
7995 E. Prentice Ave
Suite 206E
Greenwood Village, CO 80111

Email: pmoss@eaest.com

MI Identifier: 028PL

Report Date: 01/02/2019

Project: KAFB-Bulk Fuels Fac Vadose Zone, 627350DM02
Comments: Please note the total bacterial biomass for samples
BM-V2-131218-215, BM-V1-171218-122, and BM-V1-
171218-161 was low, and the samples may have PCR
inhibition.

NOTICE: This report is intended only for the addressee shown above and may contain confidential or privileged information. If the recipient of this material is not the intended recipient or if you have received this in error, please notify Microbial Insights, Inc. immediately. The data and other information in this report represent only the sample(s) analyzed and are rendered upon condition that it is not to be reproduced without approval from Microbial Insights, Inc. Thank you for your cooperation.

The QuantArray®-Chlor Approach

Quantification of *Dehalococcoides*, the only known bacterial group capable of complete reductive dechlorination of PCE and TCE to ethene, has become an indispensable component of assessment, remedy selection, and performance monitoring at sites impacted by chlorinated solvents. While undeniably a key group of halorespiring bacteria, *Dehalococcoides* are not the only bacteria of interest in the subsurface because reductive dechlorination is not the only potential biodegradation pathway operative at contaminated sites, and chlorinated ethenes are not always the primary contaminants of concern. The QuantArray®-Chlor not only includes a variety of halorespiring bacteria (*Dehalococcoides*, *Dehalobacter*, *Dehalogenimonas*, etc.) to assess the potential for reductive dechlorination of chloroethenes, chloroethanes, chlorobenzenes, chlorophenols, and chloroform, but also provides quantification of functional genes involved in aerobic (co)metabolic pathways for biodegradation of chlorinated solvents and even competing biological processes. Thus, the QuantArray®-Chlor will give site managers the ability to simultaneously yet economically evaluate the potential for biodegradation of a spectrum of common chlorinated contaminants through a multitude of anaerobic and aerobic (co) metabolic pathways to give a much more clear and comprehensive view of contaminant biodegradation.

The QuantArray®-Chlor is used to quantify specific microorganisms and functional genes to evaluate the following:

Anaerobic Reductive Dechlorination

Quantification of important halorespiring bacteria (e.g. *Dehalococcoides*, *Dehalobacter*, *Dehalogenimonas*, *Desulfitobacterium* spp.) and key functional genes (e.g. vinyl chloride reductases, TCE reductase, chloroform reductase) responsible for reductive dechlorination of a broad spectrum of chlorinated solvents.

Aerobic Cometabolism

Several different types of bacteria including methanotrophs and some toluene/phenol utilizing bacteria can co-oxidize TCE, DCE, and vinyl chloride. The QuantArray®-Chlor quantifies functional genes like soluble methane monooxygenase encoding enzymes capable of co-oxidation of chlorinated ethenes.

Aerobic (Co)metabolism of Vinyl Chloride

Ethene oxidizing bacteria are capable of cometabolism of vinyl chloride. In some cases, ethenotrophs can also utilize vinyl chloride as a growth supporting substrate. The QuantArray®-Chlor targets key functional genes in ethene metabolism.

How do QuantArrays® work?

The QuantArray®-Chlor in many respects is a hybrid technology combining the highly parallel detection of microarrays with the accurate and precise quantification provided by qPCR into a single platform. The key to highly parallel qPCR reactions is the nanoliter fluidics platform for low volume, solution phase qPCR reactions.

How are QuantArray® results reported?

One of the primary advantages of the QuantArray®-Chlor is the simultaneous quantification of a broad spectrum of different microorganisms and key functional genes involved in a variety of pathways for chlorinated hydrocarbon biodegradation. However, highly parallel quantification combined with the various metabolic and cometabolic capabilities of different target organisms can complicate data presentation. Therefore, in addition to Summary Tables, QuantArray® results will be presented as Microbial Population Summary and Comparison Figures to aid in data interpretation and subsequent evaluation of site management activities.

Types of Tables and Figures:

Microbial Population Summary

Figure presenting the concentrations of QuantArray®-Chlor target populations (e.g. *Dehalococcoides*) and functional genes (e.g. vinyl chloride reductase) relative to typically observed values.

Summary Tables

Tables of target population concentrations grouped by biodegradation pathway and contaminant type.

Comparison Figures

Depending on the project, sample results can be presented to compare changes over time or examine differences in microbial populations along a transect of the dissolved plume.

Results

Table 1: Summary of the QuantArray®-Chlor results obtained for samples BM-S1-051218-480, BM-S1-051218-489, and BM-V2-131218-144.

Sample Name Sample Date	BM-S1-051218-480 12/05/2018	BM-S1-051218-489 12/05/2018	BM-V2-131218-144 12/13/2018
<i>Reductive Dechlorination</i>	cells/g	cells/g	cells/g
<i>Dehalococcoides</i> (DHC)	<1.00E+03	<1.00E+03	<1.00E+03
tceA Reductase (TCE)	<1.00E+03	<1.00E+03	<1.00E+03
BAV1 Vinyl Chloride Reductase (BVC)	<1.00E+03	<1.00E+03	<1.00E+03
Vinyl Chloride Reductase (VCR)	<1.00E+03	<1.00E+03	<1.00E+03
<i>Dehalobacter</i> spp. (DHBt)	<1.00E+04	3.35E+04	<1.00E+04
<i>Dehalobacter</i> DCM (DCM)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalogenimonas</i> spp. (DHG)	<1.00E+04	<1.00E+04	<1.00E+04
cerA Reductase (CER)	<1.00E+04	<1.00E+04	<1.00E+04
trans-1,2-DCE Reductase (TDR)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Desulfitobacterium</i> spp. (DSB)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalobium chlorocoercia</i> (DECO)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Desulfuromonas</i> spp. (DSM)	5.31E+03 (J)	<1.00E+04	<1.00E+04
PCE Reductase (PCE-1)	<1.00E+04	<1.00E+04	<1.00E+04
PCE Reductase (PCE-2)	<1.00E+04	<1.00E+04	<1.00E+04
Chloroform Reductase (CFR)	<1.00E+04	<1.00E+04	<1.00E+04
1,1 DCA Reductase (DCA)	<1.00E+04	<1.00E+04	<1.00E+04
1,2 DCA Reductase (DCAR)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Aerobic (Co)Metabolic</i>			
Soluble Methane Monooxygenase (SMMO)	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Dioxygenase (TOD)	<1.00E+04	<1.00E+04	<1.00E+04
Phenol Hydroxylase (PHE)	1.04E+07	<1.00E+04	5.18E+03 (J)
Trichlorobenzene Dioxygenase (TCBO)	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Monooxygenase 2 (RDEG)	2.86E+06	<1.00E+04	<1.00E+04
Toluene Monooxygenase (RMO)	4.53E+06	<1.00E+04	<1.00E+04
Ethene Monooxygenase (EtnC)	<1.00E+04	<1.00E+04	<1.00E+04
Epoxyalkane Transferase (EtnE)	<1.00E+04	<1.00E+04	<1.00E+04
Dichloromethane Dehalogenase (DCMA)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Other</i>			
Total Eubacteria (EBAC)	1.96E+07	3.45E+05	5.30E+06
Sulfate Reducing Bacteria (APS)	<1.00E+04	5.66E+05	<1.00E+04
Methanogens (MGN)	<1.00E+04	<1.00E+04	<1.00E+04

Legend:

NA = Not Analyzed
I = Inhibited

NS = Not Sampled
< = Result Not Detected

J = Estimated Gene Copies Below PQL but Above LQL

Table 2: Summary of the QuantArray®-Chlor results obtained for samples BM-V2-141218-215, BM-V1-171218-122, and BM-V1-171218-161.

Sample Name Sample Date	BM-V2-141218-215 12/14/2018	BM-V1-171218-122 12/17/2018	BM-V1-171218-161 12/17/2018
<i>Reductive Dechlorination</i>	cells/g	cells/g	cells/g
<i>Dehalococcoides</i> (DHC)	<1.00E+03	<2.00E+03	<2.00E+03
tceA Reductase (TCE)	<1.00E+03	<2.00E+03	<2.00E+03
BAV1 Vinyl Chloride Reductase (BVC)	<1.00E+03	<2.00E+03	<2.00E+03
Vinyl Chloride Reductase (VCR)	<1.00E+03	<2.00E+03	<2.00E+03
<i>Dehalobacter</i> spp. (DHBt)	<1.00E+04	<2.00E+04	<2.00E+04
<i>Dehalobacter</i> DCM (DCM)	<1.00E+04	<2.00E+04	<2.00E+04
<i>Dehalogenimonas</i> spp. (DHG)	<1.00E+04	<2.00E+04	<2.00E+04
cerA Reductase (CER)	<1.00E+04	<2.00E+04	<2.00E+04
trans-1,2-DCE Reductase (TDR)	<1.00E+04	<2.00E+04	<2.00E+04
<i>Desulfotobacterium</i> spp. (DSB)	<1.00E+04	<2.00E+04	<2.00E+04
<i>Dehalobium chlorocoercia</i> (DECO)	<1.00E+04	<2.00E+04	<2.00E+04
<i>Desulfuromonas</i> spp. (DSM)	<1.00E+04	<2.00E+04	<2.00E+04
PCE Reductase (PCE-1)	<1.00E+04	<2.00E+04	<2.00E+04
PCE Reductase (PCE-2)	<1.00E+04	<2.00E+04	<2.00E+04
Chloroform Reductase (CFR)	<1.00E+04	<2.00E+04	<2.00E+04
1,1 DCA Reductase (DCA)	<1.00E+04	<2.00E+04	<2.00E+04
1,2 DCA Reductase (DCAR)	<1.00E+04	<2.00E+04	<2.00E+04
<i>Aerobic (Co)Metabolic</i>			
Soluble Methane Monooxygenase (SMMO)	<1.00E+04	<2.00E+04	<2.00E+04
Toluene Dioxygenase (TOD)	<1.00E+04	<2.00E+04	<2.00E+04
Phenol Hydroxylase (PHE)	<1.00E+04	<2.00E+04	<2.00E+04
Trichlorobenzene Dioxygenase (TCBO)	<1.00E+04	<2.00E+04	<2.00E+04
Toluene Monooxygenase 2 (RDEG)	<1.00E+04	<2.00E+04	<2.00E+04
Toluene Monooxygenase (RMO)	<1.00E+04	<2.00E+04	<2.00E+04
Ethene Monooxygenase (EtnC)	<1.00E+04	<2.00E+04	<2.00E+04
Epoxyalkane Transferase (EtnE)	<1.00E+04	<2.00E+04	<2.00E+04
Dichloromethane Dehalogenase (DCMA)	<1.00E+04	<2.00E+04	<2.00E+04
<i>Other</i>			
Total Eubacteria (EBAC)	2.31E+04 (I)	9.96E+04 (I)	2.94E+04 (I)
Sulfate Reducing Bacteria (APS)	<1.00E+04	<2.00E+04	<2.00E+04
Methanogens (MGN)	<1.00E+04	<2.00E+04	<2.00E+04

Legend:

NA = Not Analyzed
I = Inhibited

NS = Not Sampled
< = Result Not Detected

J = Estimated Gene Copies Below PQL but Above LQL

Microbial Populations BM-S1-051218-480

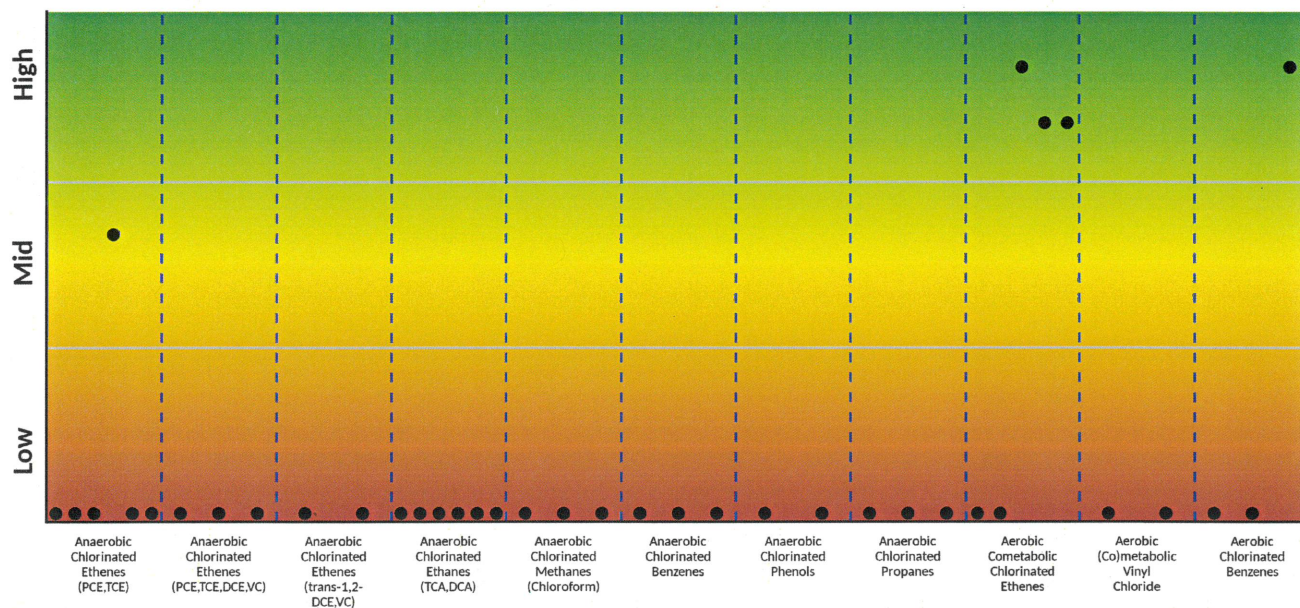


Figure 1: Microbial population summary to aid in evaluating potential pathways and biodegradation of specific contaminants.

Anaerobic - Reductive Dechlorination or Dichloroelimination		Aerobic - (Co)metabolism	
Chlorinated Ethenes (PCE, TCE)	DHC, DHBt, DSB, DSM, PCE-1, PCE-2	Chlorinated Ethenes (TCE, DCE, VC)	sMMO, TOD, PHE, RDEG, RMO
Chlorinated Ethenes (PCE, TCE, DCE, VC)	DHC, BVC, VCR	(Co)metabolic Vinyl Chloride	etnC, etnE
Chlorinated Ethenes (trans-1,2-DCE, VC)	TDR, CER	Chlorinated Benzenes	TOD, TCBO, PHE
Chlorinated Ethanes (TCA and 1,2-DCA)	DHC, DHBt, DHG, DSB ¹ , DCA, DCAR		
Chlorinated Methanes (Chloroform)	DHBt, DCM, CFR		
Chlorinated Benzenes	DHC, DHBt ² , DECO		
Chlorinated Phenols	DHC, DSB		
Chlorinated Propanes	DHC, DHG, DSB ¹		

¹Desulfotobacterium dichloroelimans DCA1. ²Implicated in reductive dechlorination of dichlorobenzene and potentially chlorobenzene.

Microbial Populations BM-S1-051218-489

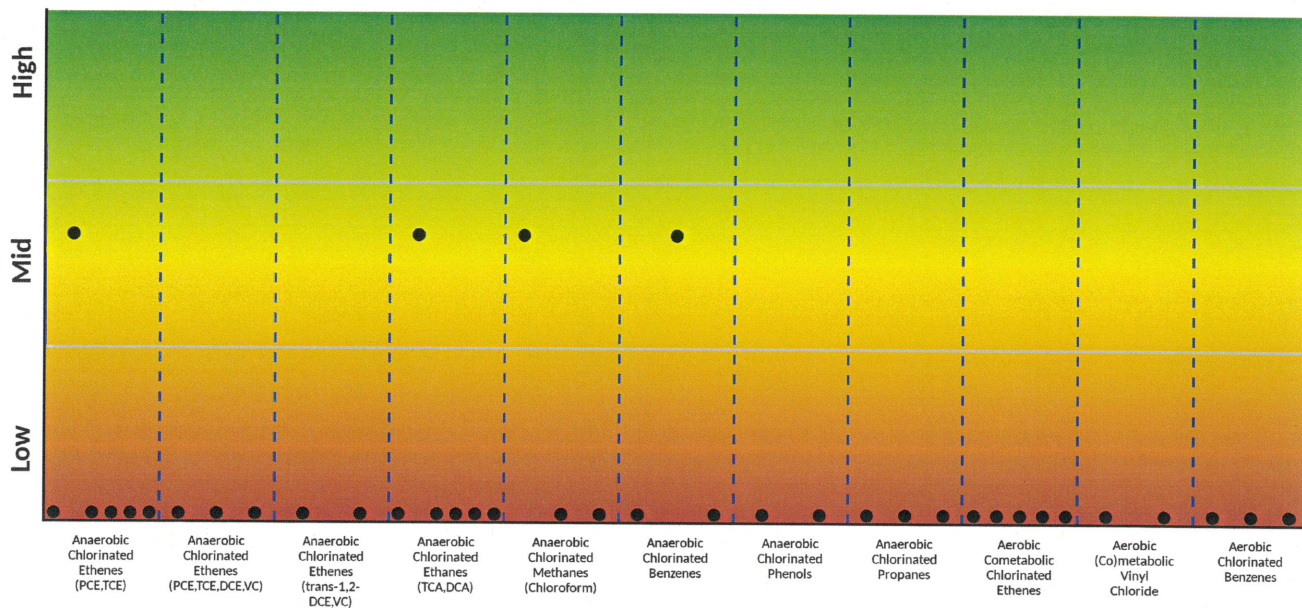


Figure 2: Microbial population summary to aid in evaluating potential pathways and biodegradation of specific contaminants.

Anaerobic - Reductive Dechlorination or Dichloroelimination			Aerobic - (Co)metabolism	
Chlorinated Ethenes (PCE, TCE)	DHC, DHBt, DSB, DSM, PCE-1, PCE-2	Chlorinated Ethenes (TCE,DCE,VC)	sMMO, TOD, PHE, RDEG, RMO	
Chlorinated Ethenes (PCE, TCE, DCE, VC)	DHC, BVC, VCR	(Co)metabolic Vinyl Chloride	etnC, etnE	
Chlorinated Ethenes (trans-1,2-DCE, VC)	TDR, CER	Chlorinated Benzenes	TOD, TCBO, PHE	
Chlorinated Ethanes (TCA and 1,2-DCA)	DHC, DHBt, DHG, DSB ¹ , DCA, DCAR			
Chlorinated Methanes (Chloroform)	DHBt, DCM, CFR			
Chlorinated Benzenes	DHC, DHBt ² , DECO			
Chlorinated Phenols	DHC, DSB			
Chlorinated Propanes	DHC, DHG, DSB ¹			

¹*Desulfotobacterium dichloroeliminans* DCA1. ²Implicated in reductive dechlorination of dichlorobenzene and potentially chlorobenzene.

Microbial Populations BM-V2-131218-144

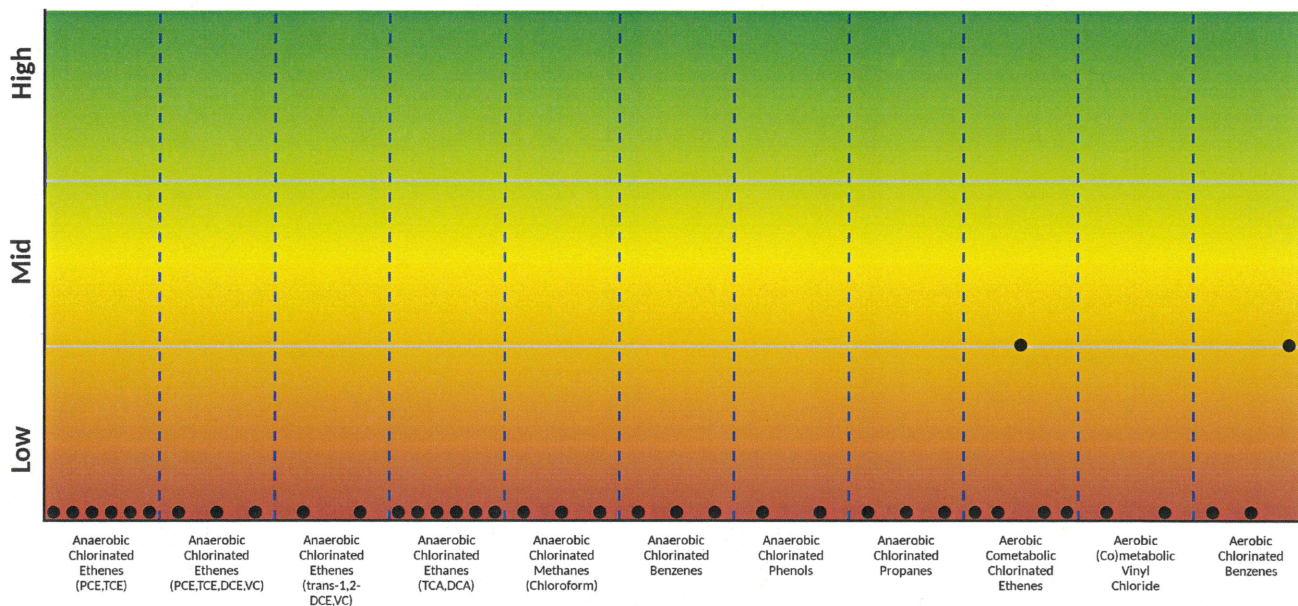


Figure 3: Microbial population summary to aid in evaluating potential pathways and biodegradation of specific contaminants.

Anaerobic - Reductive Dechlorination or Dichloroelimination

Chlorinated Ethenes (PCE, TCE)	DHC, DHBt, DSB, DSM, PCE-1, PCE-2
Chlorinated Ethenes (PCE, TCE, DCE, VC)	DHC, BVC, VCR
Chlorinated Ethenes (trans-1,2-DCE, VC)	TDR, CER
Chlorinated Ethanes (TCA and 1,2-DCA)	DHC, DHBt, DHG, DSB ¹ , DCA, DCAR
Chlorinated Methanes (Chloroform)	DHBt, DCM, CFR
Chlorinated Benzenes	DHC, DHBt ² , DECO
Chlorinated Phenols	DHC, DSB
Chlorinated Propanes	DHC, DHG, DSB ¹

Aerobic - (Co)metabolism

Chlorinated Ethenes (TCE,DCE,VC)	sMMO, TOD, PHE, RDEG, RMO
(Co)metabolic Vinyl Chloride	etnC, etnE
Chlorinated Benzenes	TOD, TCBO, PHE

¹Desulfotobacterium dichloroeliminans DCA1. ²Implicated in reductive dechlorination of dichlorobenzene and potentially chlorobenzene.

Microbial Populations BM-V2-141218-215

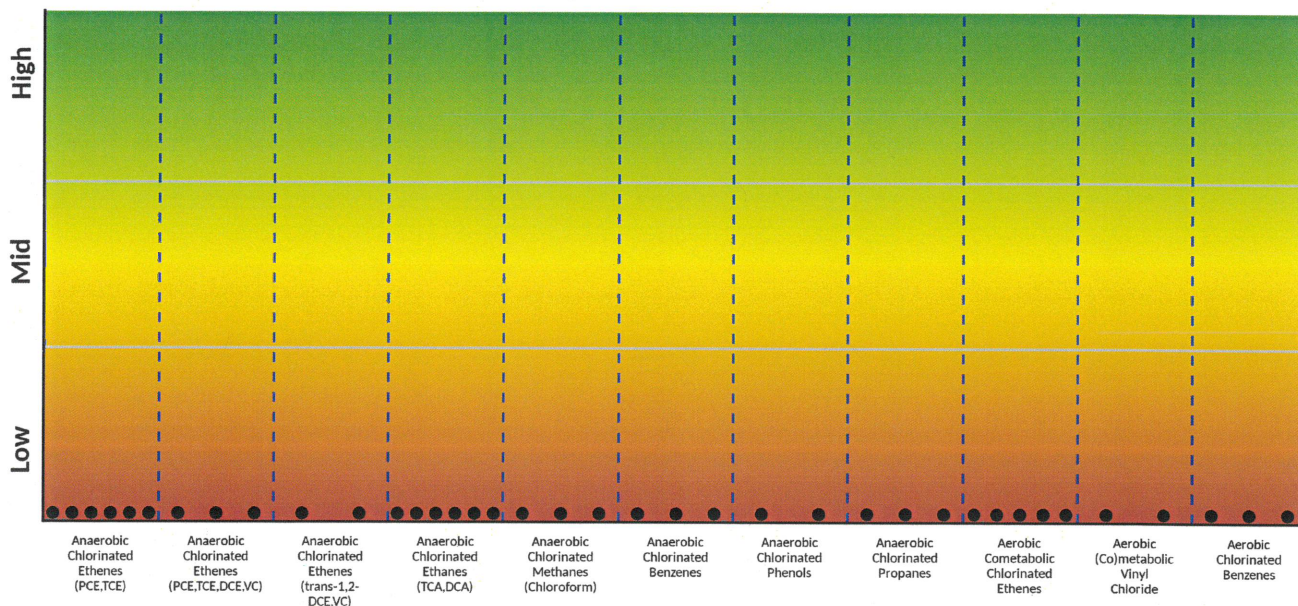


Figure 4: Microbial population summary to aid in evaluating potential pathways and biodegradation of specific contaminants.

Anaerobic - Reductive Dechlorination or Dichloroelimination		Aerobic - (Co)metabolism	
Chlorinated Ethenes (PCE, TCE)	DHC, DHBt, DSB, DSM, PCE-1, PCE-2	Chlorinated Ethenes (TCE,DCE,VC)	sMMO, TOD, PHE, RDEG, RMO
Chlorinated Ethenes (PCE, TCE, DCE, VC)	DHC, BVC, VCR	(Co)metabolic Vinyl Chloride	etnC, etnE
Chlorinated Ethenes (trans-1,2-DCE, VC)	TDR, CER	Chlorinated Benzenes	TOD, TCBO, PHE
Chlorinated Ethanes (TCA and 1,2-DCA)	DHC, DHBt, DHG, DSB ¹ , DCA, DCAR		
Chlorinated Methanes (Chloroform)	DHBt, DCM, CFR		
Chlorinated Benzenes	DHC, DHBt ² , DECO		
Chlorinated Phenols	DHC, DSB		
Chlorinated Propanes	DHC, DHG, DSB ¹		

¹*Desulfotobacterium dichloroeliminans* DCA1. ²Implicated in reductive dechlorination of dichlorobenzene and potentially chlorobenzene.

Microbial Populations BM-V1-171218-122

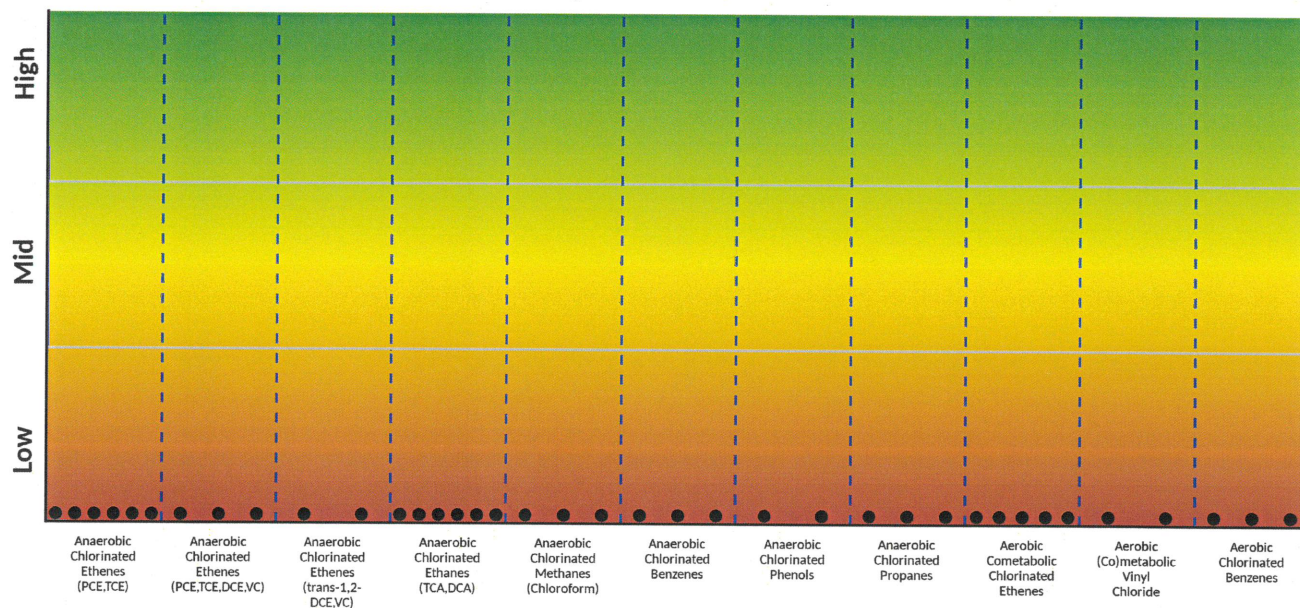


Figure 5: Microbial population summary to aid in evaluating potential pathways and biodegradation of specific contaminants.

Anaerobic - Reductive Dechlorination or Dichloroelimination		Aerobic - (Co)metabolism	
Chlorinated Ethenes (PCE, TCE)	DHC, DHBt, DSB, DSM, PCE-1, PCE-2	Chlorinated Ethenes (TCE,DCE,VC)	sMMO, TOD, PHE, RDEG, RMO
Chlorinated Ethenes (PCE, TCE, DCE, VC)	DHC, BVC, VCR	(Co)metabolic Vinyl Chloride	etnC, etnE
Chlorinated Ethenes (trans-1,2-DCE, VC)	TDR, CER	Chlorinated Benzenes	TOD, TCBO, PHE
Chlorinated Ethanes (TCA and 1,2-DCA)	DHC, DHBt, DHG, DSB ¹ , DCA, DCAR		
Chlorinated Methanes (Chloroform)	DHBt, DCM, CFR		
Chlorinated Benzenes	DHC, DHBt ² , DECO		
Chlorinated Phenols	DHC, DSB		
Chlorinated Propanes	DHC, DHG, DSB ¹		

¹*Desulfotobacterium dichloroelimians* DCA1. ²Implicated in reductive dechlorination of dichlorobenzene and potentially chlorobenzene.

Microbial Populations BM-V1-171218-161

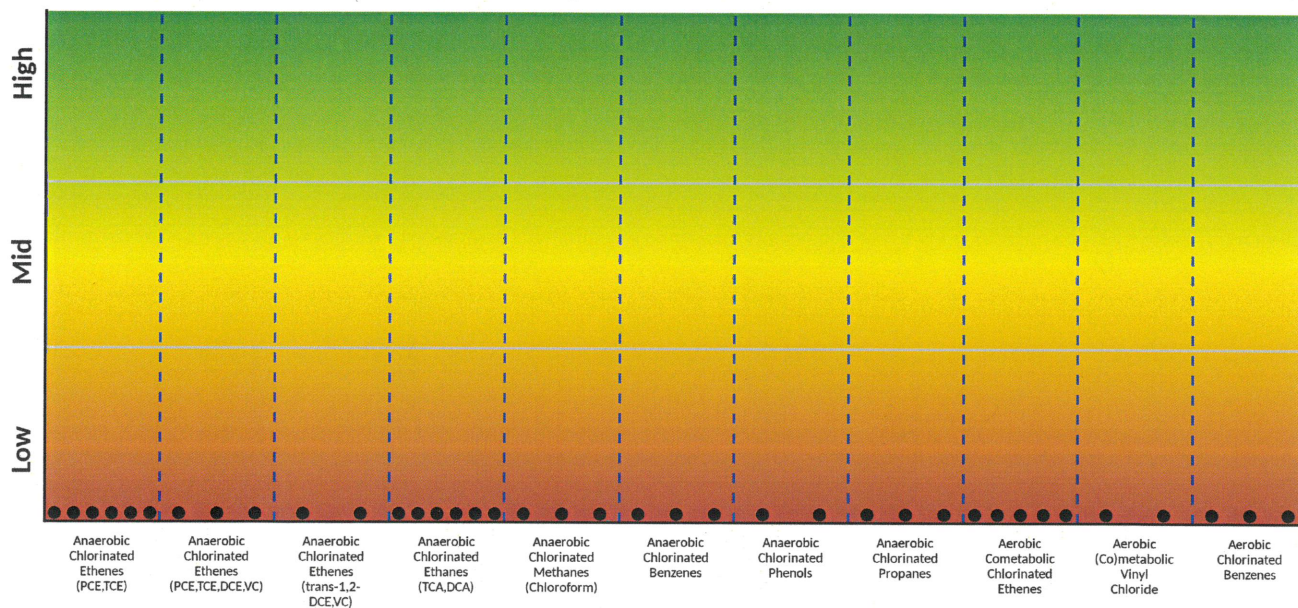


Figure 6: Microbial population summary to aid in evaluating potential pathways and biodegradation of specific contaminants.

Anaerobic - Reductive Dechlorination or Dichloroelimination		Aerobic - (Co)metabolism	
Chlorinated Ethenes (PCE, TCE)	DHC, DHBt, DSB, DSM, PCE-1, PCE-2	Chlorinated Ethenes (TCE,DCE,VC)	sMMO, TOD, PHE, RDEG, RMO
Chlorinated Ethenes (PCE, TCE, DCE, VC)	DHC, BVC, VCR	(Co)metabolic Vinyl Chloride	etnC, etnE
Chlorinated Ethenes (trans-1,2-DCE, VC)	TDR, CER	Chlorinated Benzenes	TOD, TCBO, PHE
Chlorinated Ethanes (TCA and 1,2-DCA)	DHC, DHBt, DHG, DSB ¹ , DCA, DCAR		
Chlorinated Methanes (Chloroform)	DHBt, DCM, CFR		
Chlorinated Benzenes	DHC, DHBt ² , DECO		
Chlorinated Phenols	DHC, DSB		
Chlorinated Propanes	DHC, DHG, DSB ¹		

¹ *Desulfotobacterium dichloroeliminans* DCA1. ² Implicated in reductive dechlorination of dichlorobenzene and potentially chlorobenzene.

Table 3: Summary of the QuantArray®-Chlor results for microorganisms responsible for reductive dechlorination for samples BM-S1-051218-480, BM-S1-051218-489, and BM-V2-131218-144.

Sample Name	BM-S1-051218-480	BM-S1-051218-489	BM-V2-131218-144
Sample Date	12/05/2018	12/05/2018	12/13/2018
Reductive Dechlorination	cells/g	cells/g	cells/g
<i>Dehalococcoides</i> (DHC)	<1.00E+03	<1.00E+03	<1.00E+03
tceA Reductase (TCE)	<1.00E+03	<1.00E+03	<1.00E+03
BAV1 Vinyl Chloride Reductase (BVC)	<1.00E+03	<1.00E+03	<1.00E+03
Vinyl Chloride Reductase (VCR)	<1.00E+03	<1.00E+03	<1.00E+03
<i>Dehalobacter</i> spp. (DHBt)	<1.00E+04	3.35E+04	<1.00E+04
<i>Dehalobacter</i> DCM (DCM)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalogenimonas</i> spp. (DHG)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Desulfotobacterium</i> spp. (DSB)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalobium chlorocoercia</i> (DECO)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Desulfuromonas</i> spp. (DSM)	5.31E+03 (J)	<1.00E+04	<1.00E+04

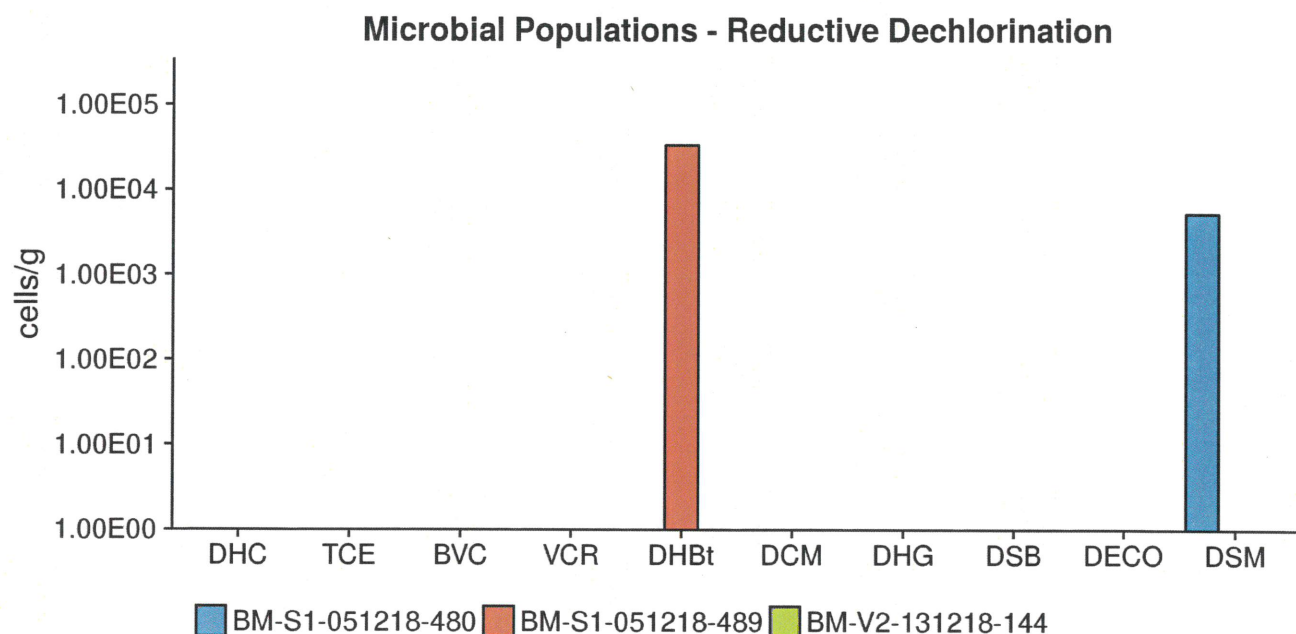


Figure 7: Comparison - microbial populations involved in reductive dechlorination.

Table 4: Summary of the QuantArray®-Chlor results for microorganisms responsible for reductive dechlorination for samples BM-S1-051218-480, BM-S1-051218-489, and BM-V2-131218-144.

Sample Name	BM-S1-051218-480	BM-S1-051218-489	BM-V2-131218-144
Sample Date	12/05/2018	12/05/2018	12/13/2018
Reductive Dechlorination	cells/g	cells/g	cells/g
Chloroform Reductase (CFR)	<1.00E+04	<1.00E+04	<1.00E+04
1,1 DCA Reductase (DCA)	<1.00E+04	<1.00E+04	<1.00E+04
1,2 DCA Reductase (DCAR)	<1.00E+04	<1.00E+04	<1.00E+04
PCE Reductase (PCE-1)	<1.00E+04	<1.00E+04	<1.00E+04
PCE Reductase (PCE-2)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalogenimonas trans</i> -1,2-DCE Reductase (TDR)	<1.00E+04	<1.00E+04	<1.00E+04
<i>Dehalogenimonas cerA</i> Reductase (CER)	<1.00E+04	<1.00E+04	<1.00E+04

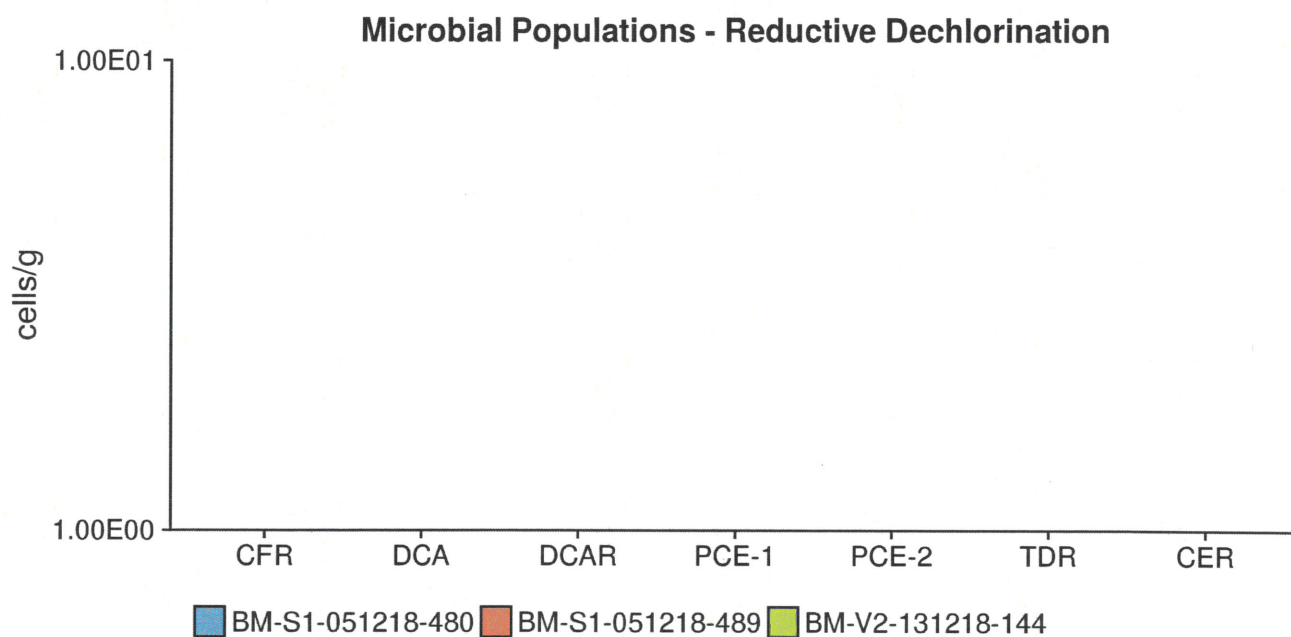


Figure 8: Comparison - microbial populations involved in reductive dechlorination.

Table 5: Summary of the QuantArray®-Chlor results for microorganisms responsible for reductive dechlorination for samples BM-V2-141218-215, BM-V1-171218-122, and BM-V1-171218-161.

Sample Name	BM-V2-141218-215	BM-V1-171218-122	BM-V1-171218-161
Sample Date	12/14/2018	12/17/2018	12/17/2018
Reductive Dechlorination	cells/g	cells/g	cells/g
<i>Dehalococcoides</i> (DHC)	<1.00E+03	<2.00E+03	<2.00E+03
tceA Reductase (TCE)	<1.00E+03	<2.00E+03	<2.00E+03
BAV1 Vinyl Chloride Reductase (BVC)	<1.00E+03	<2.00E+03	<2.00E+03
Vinyl Chloride Reductase (VCR)	<1.00E+03	<2.00E+03	<2.00E+03
<i>Dehalobacter</i> spp. (DHBt)	<1.00E+04	<2.00E+04	<2.00E+04
<i>Dehalobacter</i> DCM (DCM)	<1.00E+04	<2.00E+04	<2.00E+04
<i>Dehalogenimonas</i> spp. (DHG)	<1.00E+04	<2.00E+04	<2.00E+04
<i>Desulfotobacterium</i> spp. (DSB)	<1.00E+04	<2.00E+04	<2.00E+04
<i>Dehalobium chlorocoercia</i> (DECO)	<1.00E+04	<2.00E+04	<2.00E+04
<i>Desulfuromonas</i> spp. (DSM)	<1.00E+04	<2.00E+04	<2.00E+04

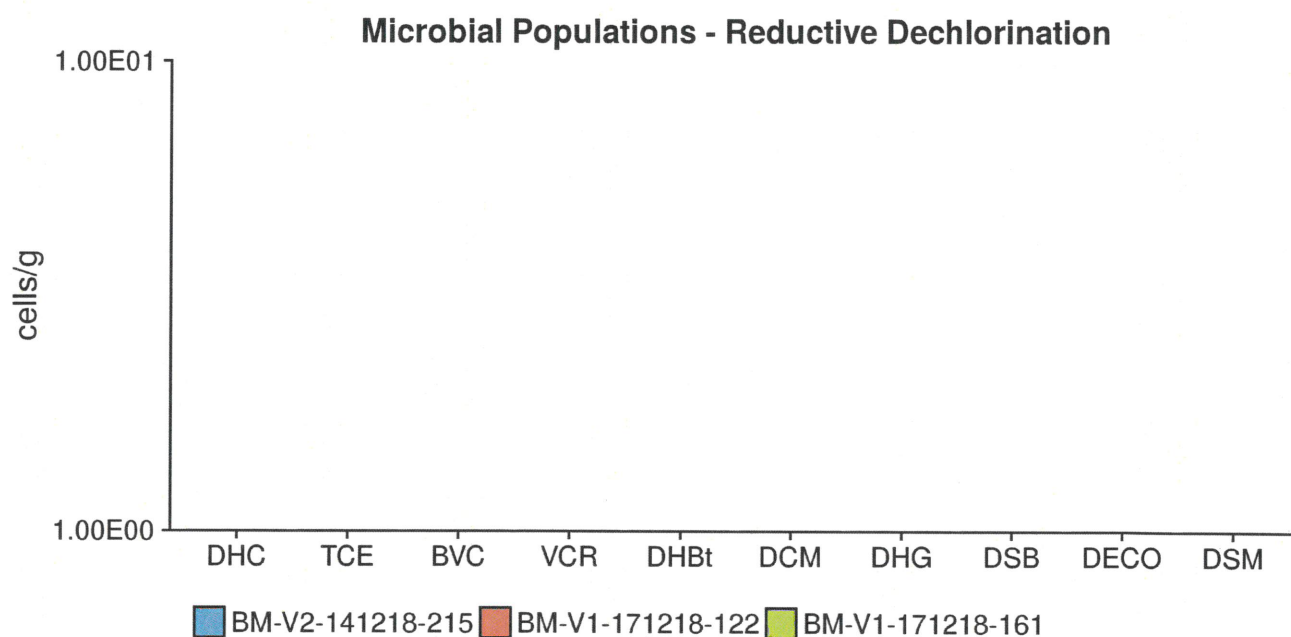


Figure 9: Comparison - microbial populations involved in reductive dechlorination.

Table 6: Summary of the QuantArray®-Chlor results for microorganisms responsible for reductive dechlorination for samples BM-V2-141218-215, BM-V1-171218-122, and BM-V1-171218-161.

Sample Name	BM-V2-141218-215	BM-V1-171218-122	BM-V1-171218-161
Sample Date	12/14/2018	12/17/2018	12/17/2018
Reductive Dechlorination	cells/g	cells/g	cells/g
Chloroform Reductase (CFR)	<1.00E+04	<2.00E+04	<2.00E+04
1,1 DCA Reductase (DCA)	<1.00E+04	<2.00E+04	<2.00E+04
1,2 DCA Reductase (DCAR)	<1.00E+04	<2.00E+04	<2.00E+04
PCE Reductase (PCE-1)	<1.00E+04	<2.00E+04	<2.00E+04
PCE Reductase (PCE-2)	<1.00E+04	<2.00E+04	<2.00E+04
Dehalogenimonas trans-1,2-DCE Reductase (TDR)	<1.00E+04	<2.00E+04	<2.00E+04
Dehalogenimonas cerA Reductase (CER)	<1.00E+04	<2.00E+04	<2.00E+04

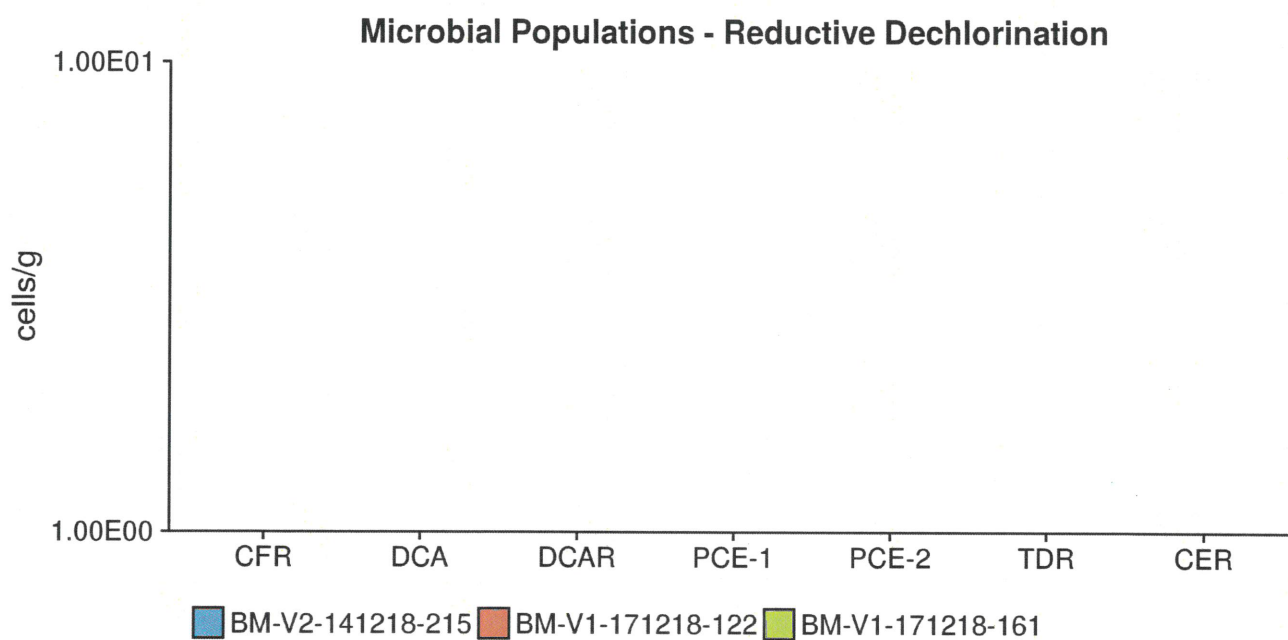


Figure 10: Comparison - microbial populations involved in reductive dechlorination.

Table 7: Summary of the QuantArray®-Chlor results for microorganisms responsible for aerobic (co)metabolism for samples BM-S1-051218-480, BM-S1-051218-489, and BM-V2-131218-144.

Sample Name	BM-S1-051218-480	BM-S1-051218-489	BM-V2-131218-144
Sample Date	12/05/2018	12/05/2018	12/13/2018
<i>Aerobic (Co)Metabolic</i>	cells/g	cells/g	cells/g
Soluble Methane Monooxygenase (SMMO)	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Dioxygenase (TOD)	<1.00E+04	<1.00E+04	<1.00E+04
Phenol Hydroxylase (PHE)	1.04E+07	<1.00E+04	5.18E+03 (J)
Trichlorobenzene Dioxygenase (TCBO)	<1.00E+04	<1.00E+04	<1.00E+04
Toluene Monooxygenase 2 (RDEG)	2.86E+06	<1.00E+04	<1.00E+04
Toluene Monooxygenase (RMO)	4.53E+06	<1.00E+04	<1.00E+04
Ethene Monooxygenase (EtnC)	<1.00E+04	<1.00E+04	<1.00E+04
Epoxyalkane Transferase (EtnE)	<1.00E+04	<1.00E+04	<1.00E+04
Dichloromethane Dehalogenase (DCMA)	<1.00E+04	<1.00E+04	<1.00E+04

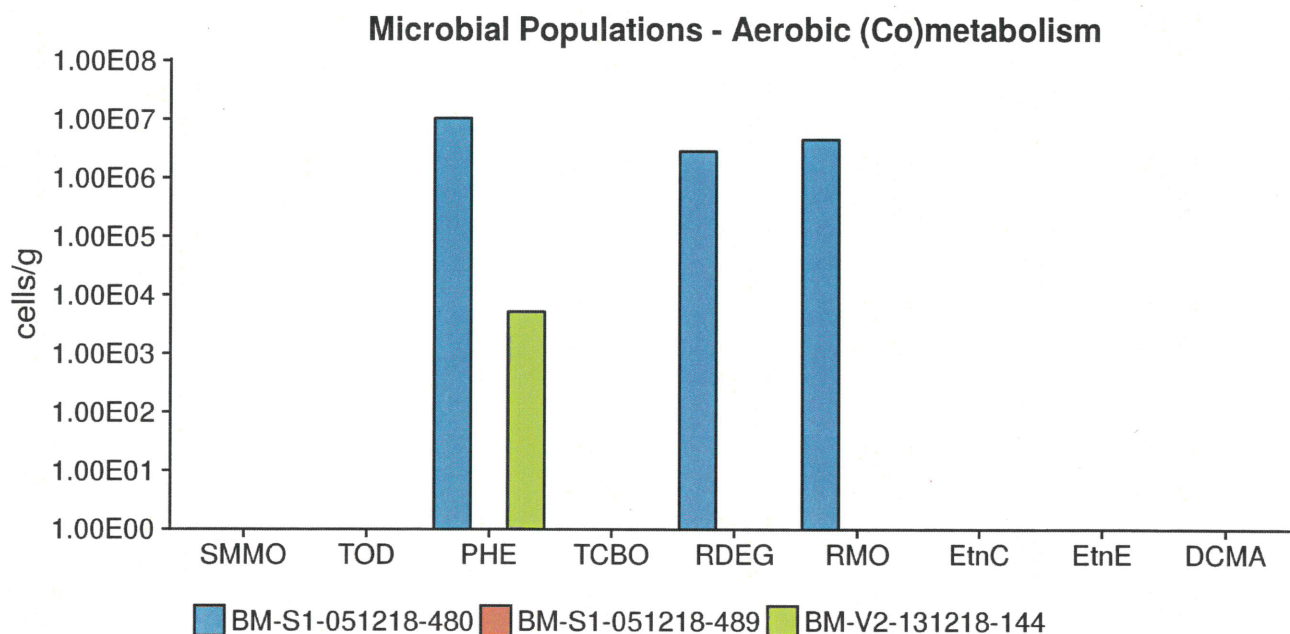


Figure 11: Comparison - microbial populations involved in aerobic (co)metabolism.

Table 8: Summary of the QuantArray®-Chlor results for microorganisms responsible for aerobic (co)metabolism for samples BM-V2-141218-215, BM-V1-171218-122, and BM-V1-171218-161.

Sample Name	BM-V2-141218-215	BM-V1-171218-122	BM-V1-171218-161
Sample Date	12/14/2018	12/17/2018	12/17/2018
<i>Aerobic (Co)Metabolic</i>	cells/g	cells/g	cells/g
Soluble Methane Monooxygenase (SMMO)	<1.00E+04	<2.00E+04	<2.00E+04
Toluene Dioxygenase (TOD)	<1.00E+04	<2.00E+04	<2.00E+04
Phenol Hydroxylase (PHE)	<1.00E+04	<2.00E+04	<2.00E+04
Trichlorobenzene Dioxygenase (TCBO)	<1.00E+04	<2.00E+04	<2.00E+04
Toluene Monooxygenase 2 (RDEG)	<1.00E+04	<2.00E+04	<2.00E+04
Toluene Monooxygenase (RMO)	<1.00E+04	<2.00E+04	<2.00E+04
Ethene Monooxygenase (EtnC)	<1.00E+04	<2.00E+04	<2.00E+04
Epoxyalkane Transferase (EtnE)	<1.00E+04	<2.00E+04	<2.00E+04
Dichloromethane Dehalogenase (DCMA)	<1.00E+04	<2.00E+04	<2.00E+04

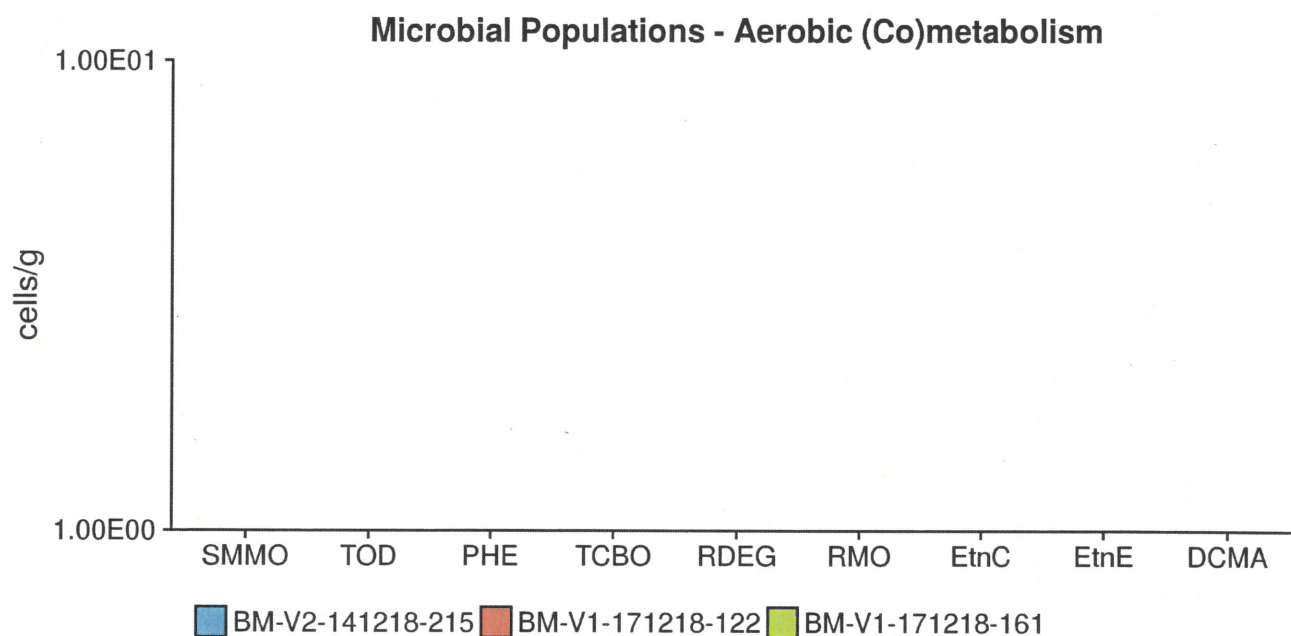


Figure 12: Comparison - microbial populations involved in aerobic (co)metabolism.

Table 9: Summary of the QuantArray® results for total bacteria and other populations for samples BM-S1-051218-480, BM-S1-051218-489, and BM-V2-131218-144.

Sample Name	BM-S1-051218-480	BM-S1-051218-489	BM-V2-131218-144
Sample Date	12/05/2018	12/05/2018	12/13/2018
Other	cells/g	cells/g	cells/g
Total Eubacteria (EBAC)	1.96E+07	3.45E+05	5.30E+06
Sulfate Reducing Bacteria (APS)	<1.00E+04	5.66E+05	<1.00E+04
Methanogens (MGN)	<1.00E+04	<1.00E+04	<1.00E+04

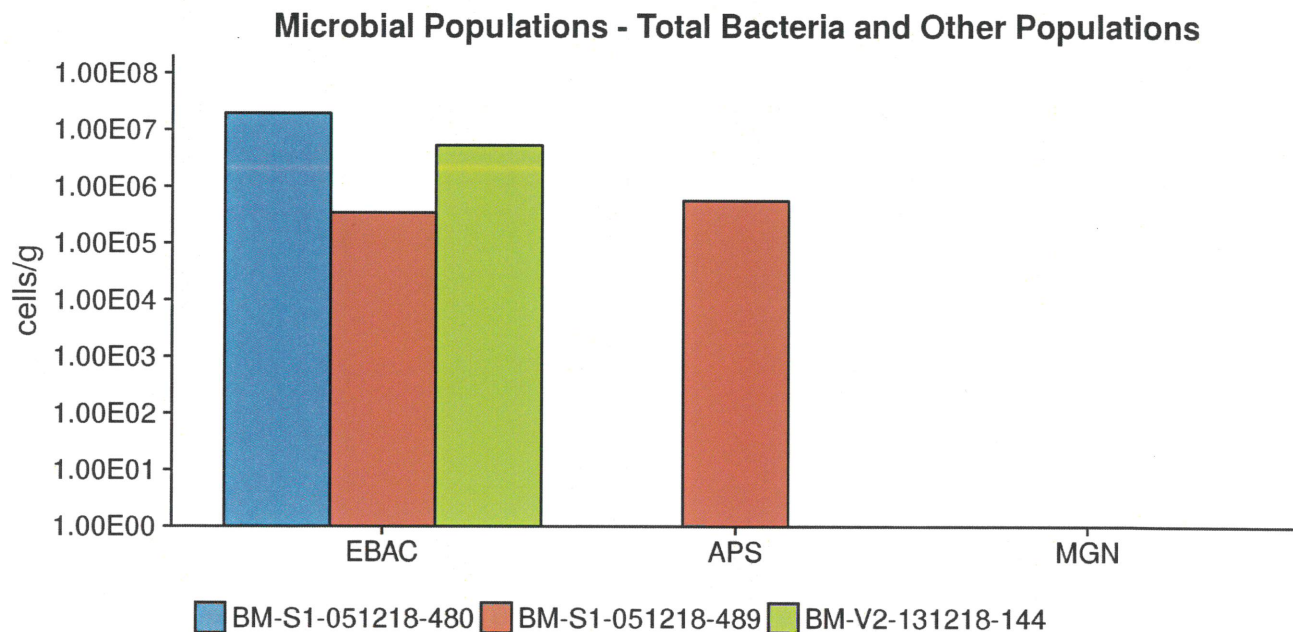


Figure 13: Comparison - microbial populations.

Table 10: Summary of the QuantArray® results for total bacteria and other populations for samples BM-V2-141218-215, BM-V1-171218-122, and BM-V1-171218-161.

Sample Name	BM-V2-141218-215	BM-V1-171218-122	BM-V1-171218-161
Sample Date	12/14/2018	12/17/2018	12/17/2018
Other	cells/g	cells/g	cells/g
Total Eubacteria (EBAC)	2.31E+04 (I)	9.96E+04 (I)	2.94E+04 (I)
Sulfate Reducing Bacteria (APS)	<1.00E+04	<2.00E+04	<2.00E+04
Methanogens (MGN)	<1.00E+04	<2.00E+04	<2.00E+04

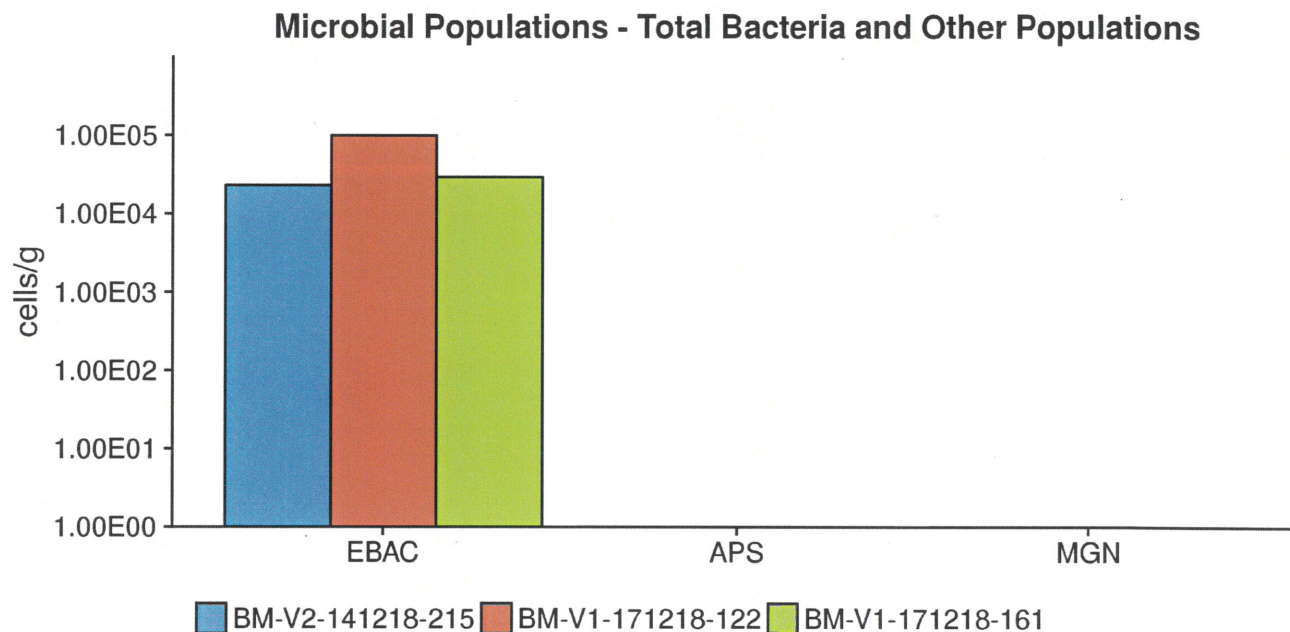


Figure 14: Comparison - microbial populations.

Interpretation

The overall purpose of the QuantArray®-Chlor is to give site managers the ability to simultaneously yet economically evaluate the potential for biodegradation of a spectrum of common chlorinated contaminants through a multitude of anaerobic and aerobic (co)metabolic pathways in order to provide a clearer and more comprehensive view of contaminant biodegradation. The following discussion describes the interpretation of results in general terms and is meant to serve as a guide.

Reductive Dechlorination - Chlorinated Ethenes: While a number of bacterial cultures including *Dehalococcoides*, *Dehalobacter*, *Desulfotobacterium*, and *Desulfuromonas* spp. capable of utilizing PCE and TCE as growth-supporting electron acceptors have been isolated [1–5], *Dehalococcoides* may be the most important because they are the only bacterial group that has been isolated to date which is capable of complete reductive dechlorination of PCE to ethene [6]. In fact, the presence of *Dehalococcoides* has been associated with complete reductive dechlorination to ethene at sites across North America and Europe [7], and Lu et al. [8] have proposed using a *Dehalococcoides* concentration of 1×10^4 cells/mL as a screening criterion to identify sites where biological reductive dechlorination is predicted to proceed at “generally useful” rates.

At chlorinated ethene sites, any “stall” leading to the accumulation of daughter products, especially vinyl chloride, would be a substantial concern. While *Dehalococcoides* concentrations greater than 1×10^4 cells/mL correspond to ethene production and useful rates of dechlorination, the range of chlorinated ethenes degraded varies by strain within the *Dehalococcoides* genus [6, 9], and the presence of co-contaminants and competitors can have complex impacts on the halo-respiring microbial community [10–15]. Therefore, QuantArray®-Chlor also provides quantification of a suite of reductive dehalogenase genes (PCE, TCE, BVC, VCR, CER, and TDR) to more definitively confirm the potential for reductive dechlorination of all chlorinated ethene compounds including vinyl chloride.

Perhaps most importantly, QuantArray®-Chlor quantifies TCE reductase (TCE) and both known vinyl chloride reductase genes (BVC, VCR) from *Dehalococcoides* to conclusively evaluate the potential for complete reductive dechlorination of chlorinated ethenes to non-toxic ethene [16–18]. In addition, the analysis also includes quantification of reductive dehalogenase genes from *Dehalogenimonas* spp. capable of reductive dechlorination of chlorinated ethenes. More specifically, these are the trans-1,2-DCE dehalogenase gene (TDR) from strain WBC-2 [19] and the vinyl chloride reductase gene (CER) from GP, the only known organisms other than *Dehalococcoides* capable of vinyl chloride reduction [20]. Finally, PCE reductase genes responsible for sequential reductive dechlorination of PCE to cis-DCE by *Sulfurospirillum* and *Geobacter* spp. are also quantified. In mixed cultures, evidence increasingly suggests that partial dechlorinators like *Sulfurospirillum* and *Geobacter* may be responsible for the majority of reductive dechlorination of PCE to TCE and cis-DCE while *Dehalococcoides* functions more as cis-DCE and vinyl chloride reducing specialists [10, 21].

Reductive Dechlorination - Chlorinated Ethanes: Under anaerobic conditions, chlorinated ethanes are susceptible to reductive dechlorination by several groups of halo-respiring bacteria including *Dehalobacter*, *Dehalogenimonas*, and *Dehalococcoides*. While the reported range of chlorinated ethanes utilized varies by genus, species, and sometimes at the strain level, several general observations can be made regarding biodegradation pathways and daughter product formation. *Dehalobacter* spp. have been isolated that are capable of sequential reductive dechlorination of 1,1,1-TCA through 1,1-DCA to chloroethane [13]. Biodegradation of 1,1,2-TCA by several halo-respiring bacteria including *Dehalobacter* and *Dehalogenimonas* spp. proceeds via dichloroelimination producing vinyl chloride [22–24]. Similarly, 1,2-DCA biodegradation by *Dehalobacter*, *Dehalogenimonas*, and *Dehalococcoides* occurs via dichloroelimination producing ethene. While not utilized by many *Desulfotobacterium* isolates, at least one strain, *Desulfotobacterium dichloroeliminans* strain DCA1, is also capable of dichloroelimination of 1,2-DCA [25]. The 1,2-dichloroethane reductive dehalogenase gene (DCAR) from members of *Desulfotobacterium* and *Dehalobacter* is known to dechlorinate 1,2-DCA to ethene, while the 1,1-dichloroethane reductive dehalogenase (DCA) targets the gene responsible for 1,1-DCA dechlorination in some strains of *Dehalobacter*. In addition to chloroform, chloroform reductase (CFR) has also been shown to be responsible for reductive dechlorination of 1,1,1-TCA [26].

Reductive Dechlorination - Chlorinated Methanes: Chloroform is a common co-contaminant at chlorinated solvent sites and can inhibit reductive dechlorination of chlorinated ethenes. Grostern et al. demonstrated that a *Dehalobacter* population was capable of reductive dechlorination of chloroform to produce dichloromethane [27]. The *cfrA* gene encodes the reductase which catalyzes this initial step in chloroform biodegradation [26]. Justicia-Leon et al. have since shown that dichloromethane can support growth of a distinct group of *Dehalobacter* strains via fermentation [28]. The *Dehalobacter* DCM assay targets the 16S rRNA gene of these strains.

Reductive Dechlorination - Chlorinated Benzenes: Chlorinated benzenes are an important class of industrial solvents and chemical intermediates in the production of drugs, dyes, herbicides, and insecticides. The physical-chemical properties of chlorinated benzenes as well as susceptibility to biodegradation are functions of their degree of chlorination and the positions of chlorine substituents. Under anaerobic conditions, reductive dechlorination of higher chlorinated benzenes including hexachlorobenzene (HCB),

pentachlorobenzene (PeCB), tetrachlorobenzene (TeCB) isomers, and trichlorobenzene (TCB) isomers has been well documented [29], although biodegradation of individual compounds and isomers varies between isolates. For example, *Dehalococcoides* strain CBDB1 reductively dechlorinates HCB, PeCB, all three TeCB isomers, 1,2,3-TCB, and 1,2,4-TCB [9, 30]. *Dehalobium chlorocoercia* DF-1 has been shown to be capable of reductive dechlorination of HCB, PeCB, and 1,2,3,5-TeCB [31]. The dichlorobenzene (DCB) isomers and chlorobenzene (CB) were considered relatively recalcitrant under anaerobic conditions. However, new evidence has demonstrated reductive dechlorination of DCBs to CB and CB to benzene [32] with corresponding increases in concentrations of *Dehalobacter* spp. [33].

Reductive Dechlorination - Chlorinated Phenols: Pentachlorophenol (PCP) was one of the most widely used biocides in the U.S. and despite residential use restrictions, is still extensively used industrially as a wood preservative. Along with PCP, the tetrachlorophenol and trichlorophenol isomers were also used as fungicides in wood preserving formulations. 2,4-Dichlorophenol and 2,4,5-TCP were used as chemical intermediates in herbicide production (e.g. 2,4-D) and chlorophenols are known byproducts of chlorine bleaching in the pulp and paper industry. While the range of compounds utilized varies by strain, some *Dehalococcoides* isolates are capable of reductive dechlorination of PCP and other chlorinated phenols. For example, *Dehalococcoides* strain CBDB1 is capable of utilizing PCP, all three tetrachlorophenol (TeCP) congeners, all six trichlorophenol (TCP) congeners, and 2,3-dichlorophenol (2,3-DCP). PCP dechlorination by strain CBDB1 produces a mixture of 3,5-DCP, 3,4-DCP, 2,4-DCP, 3-CP, and 4-CP [34]. In the same study, however, *Dehalococcoides* strain 195 dechlorinated a more narrow spectrum of chlorophenols which included 2,3-DCP, 2,3,4-TCP, and 2,3,6-TCP, but no other TCPs or PCP. Similar to *Dehalococcoides*, some species and strains of *Desulfotobacterium* are capable of utilizing PCP and other chlorinated phenols. *Desulfotobacterium hafniense* PCP-1 is capable of reductive dechlorination of PCP to 3-CP [35]. However, the ability to biodegrade PCP is not universal among *Desulfotobacterium* isolates. *Desulfotobacterium* sp. strain PCE1 and *D. chlororespirans* strain Co23, for example, can utilize some TCP and DCP isomers, but not PCP for growth [2, 36].

Reductive Dechlorination - Chlorinated Propanes: *Dehalogenimonas* is a recently described bacterial genus of the phylum Chloroflexi which also includes the well-known chloroethene-respiring *Dehalococcoides* [23]. The *Dehalogenimonas* isolates characterized to date are also halo-respiring bacteria, but utilize a rather unique range of chlorinated compounds as electron acceptors including chlorinated propanes (1,2,3-TCP and 1,2-DCP) and a variety of other vicinally chlorinated alkanes including 1,1,2,2-tetrachloroethane, 1,1,2-trichloroethane, and 1,2-dichloroethane [23].

Aerobic - Chlorinated Ethene Cometabolism: Under aerobic conditions, several different types of bacteria including methane-oxidizing bacteria (methanotrophs), and many benzene, toluene, ethylbenzene, xylene, and (BTEX)-utilizing bacteria can cometabolize or co-oxidize TCE, DCE, and vinyl chloride [37]. In general, cometabolism of chlorinated ethenes is mediated by monooxygenase enzymes with "relaxed" specificity that oxidize a primary (growth supporting) substrate (e.g. methane) and co-oxidize the chlorinated compound (e.g. TCE). QuantArray[®]-Chlor provides quantification of a suite of genes encoding oxygenase enzymes capable of co-oxidation of chlorinated ethenes including soluble methane monooxygenase (sMMO). Soluble methane monooxygenases co-oxidize a broad range of chlorinated compounds [38–41] including TCE, *cis*-DCE, and vinyl chloride. Furthermore, soluble methane monooxygenases are generally believed to support greater rates of aerobic cometabolism [40]. QuantArray[®]-Chlor also quantifies aromatic oxygenase genes encoding ring hydroxylating toluene monooxygenase genes (RMO, RDEG), toluene dioxygenase (TOD) and phenol hydroxylases (PHE) capable of TCE co-oxidation [42–46]. TCE or a degradation product has been shown to induce expression of toluene monooxygenases in some laboratory studies [43, 47] raising the possibility of TCE cometabolism with an alternative (non-aromatic) growth substrate. Moreover, while a number of additional factors must be considered, recent research under ESTCP Project 201584 has shown positive correlations between concentrations of monooxygenase genes (soluble methane monooxygenase, ring hydroxylating monooxygenases, and phenol hydroxylase) and the rate of TCE degradation [48].

Aerobic - Chlorinated Ethane Cometabolism: While less widely studied than cometabolism of chlorinated ethenes, some chlorinated ethanes are also susceptible to co-oxidation. As mentioned previously, soluble methane monooxygenases (sMMO) exhibit very relaxed specificity. In laboratory studies, sMMO has been shown to co-oxidize a number of chlorinated ethanes including 1,1,1-TCA and 1,2-DCA [38, 40].

Aerobic - Vinyl Chloride Cometabolism: Beginning in the early 1990s, numerous microcosm studies demonstrated aerobic oxidation of vinyl chloride under MNA conditions without the addition of exogenous primary substrates. Since then, strains of

Mycobacterium, *Nocardioides*, *Pseudomonas*, *Ochrobactrum*, and *Ralstonia* species have been isolated which are capable of aerobic growth on both ethene and vinyl chloride (see Mattes et al. [49] for a review). The initial steps in the pathway are the monooxygenase (*etnABCD*) catalyzed conversion of ethene and vinyl chloride to their respective epoxyalkanes (epoxyethane and chlorooxirane), followed by epoxyalkane:CoM transferase (*etnE*) mediated conjugation and breaking of the epoxide [50].

Aerobic - Chlorinated Benzenes: In general, chlorobenzenes with four or less chlorine groups are susceptible to aerobic biodegradation and can serve as growth-supporting substrates. Toluene dioxygenase (TOD) has a relatively relaxed substrate specificity and mediates the incorporation of both atoms of oxygen into the aromatic ring of benzene and substituted benzenes (toluene and chlorobenzene). Comparison of TOD levels in background and source zone samples from a CB-impacted site suggested that CBs promoted growth of TOD-containing bacteria [51]. In addition, aerobic biodegradation of some trichlorobenzene and even tetrachlorobenzene isomers is initiated by a group of related trichlorobenzene dioxygenase genes (TCBO). Finally, phenol hydroxylases catalyze the continued oxidation and in some cases, the initial oxidation of a variety of monoaromatic compounds. In an independent study, significant increases in numbers of bacteria containing PHE genes corresponded to increases in biodegradation of DCB isomers [51].

Aerobic - Chlorinated Methanes: Many aerobic methylotrophic bacteria, belonging to diverse genera (*Hyphomicrobium*, *Methylobacterium*, *Methylophilus*, *Pseudomonas*, *Paracoccus*, and *Alibacter*) have been isolated which are capable of utilizing dichloromethane (DCM) as a growth substrate. The DCM metabolic pathway in methylotrophic bacteria is initiated by a dichloromethane dehalogenase (DCMA) gene. DCMA is responsible for aerobic biodegradation of dichloromethane by methylotrophs by first producing formaldehyde which is then further oxidized [52]. As discussed in previous sections, soluble methane monooxygenase (sMMO) exhibits relaxed specificity and co-oxidizes a broad spectrum of chlorinated hydrocarbons. In addition to chlorinated ethenes, sMMO has been shown to co-oxidize chloroform in laboratory studies [38, 41].

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SITE LOGIC Report

Abiotic Potential

Contact: Pamela J. Moss

Phone: 303-590-9143

Address: EA Engineering- Greenwood Village, CO
7995 E. Prentice Ave, Suite 206E
Greenwood Village, CO 80111

Email: pmoss@eaest.com

MI Identifier: 028PL

Report Date: 1-3-2019

Project Name: KAFB-Bulk Fuels Fac Vadose Zone

Comments:

Overview

Although not always fully considered, abiotic degradation can be a substantial or even the primary process for chlorinated hydrocarbon destruction at sites undergoing or transitioning to monitored natural attenuation (MNA). A variety of iron-bearing minerals including iron sulfides (mackinawite and pyrite), iron oxides (magnetite), green rust, and iron-bearing clays are capable of complete or nearly complete degradation of PCE, TCE, and carbon tetrachloride (He et al. 2009). Some iron-bearing minerals also catalyze the degradation of chlorinated ethanes and the lesser chlorinated ethenes cis-dichloroethene (DCE) and vinyl chloride. While the quantities and types will vary, these reactive iron minerals are frequently identified in subsurface environments under iron reducing and sulfate reducing conditions.

Brown et al. (2007) recommend four avenues for evaluating the role of abiotic processes in contaminant attenuation. First, examining contaminant concentrations along the flow path - decreasing parent compound concentrations with no evidence of accumulation of chlorinated transformation products like cis-DCE and vinyl chloride suggest abiotic degradation. Performing compound specific isotope analysis (CSIA) or monitoring for products unique to abiotic reactions such as acetylene can also provide a strong line of evidence. Microcosm studies with native sediment and killed controls can also be performed. Finally, Brown et al. (2007) suggest performing mineralogical analyses on aquifer sediment to characterize reactive minerals such as magnetite or iron monosulfides

Magnetic Susceptibility - Magnetite

Magnetite (Fe_3O_4) is a mixed valence iron mineral shown to react with PCE, TCE, and carbon tetrachloride. Furthermore, Ferrey et al. (2004) conclusively linked the observed degradation of cis-DCE at a former ammunition plant to magnetite in the subsurface. No direct chemical test is available for quantification of magnetite. However, magnetite is the most abundant mineral in natural sediments that exhibits magnetic behavior. Therefore, magnetic susceptibility provides an inexpensive and valuable estimate of the quantity of magnetite in environmental samples.

X-ray Diffraction (XRD) - Mackinawite, Pyrite, Magnetite and Green Rust

XRD is one of the primary techniques used to identify unknown crystalline materials. Most minerals are crystalline and will scatter X-rays in a regular, characteristic manner dependent on their crystal structure.

- Mackinawite is the most reactive of the iron-bearing minerals and a crystalline form (tetragonal FeS) can be detected by XRD. Mackinawite will transform PCE and TCE primarily by elimination to acetylene. Carbon tetrachloride is transformed mainly to chloroform but carbon dioxide, formate, and carbon disulfide have also been detected. Finally, the more heavily chlorinated ethanes including hexachloroethane, pentachloroethane, and tetrachloroethanes react to form chlorinated ethenes which can be further degraded.
- Pyrite (FeS_2) catalyzes beta elimination transforming PCE, TCE, and cis-DCE to acetylene and ethene. Vinyl chloride is transformed to ethene and ethane. Pyrite is also capable of degradation of carbon tetrachloride potentially forming a number of products including chloroform, carbon dioxide, carbon disulfide, and formate depending on reaction conditions.
- While not quantitative like the magnetic susceptibility test, XRD can also detect magnetite when present at between 2% and 5% on a weight basis.
- Green rusts have been reported to transform a number of common chlorinated contaminants including cis-DCE, vinyl chloride, trichloroethanes, and tetrachloroethanes. While special sample care to prevent oxidation would be needed, XRD can be used to detect green rust.

Percent Clay

Clays have large surface areas, balanced by exchangeable cations, which can bind a large number of both organic and inorganic molecules impacting their availability and reactivity in the subsurface. While less well studied than the other iron-bearing minerals, various phyllosilicate clays have been shown to be capable of degradation of PCE, TCE, cis-DCE, vinyl chloride, and carbon tetrachloride

Sample Location	BM-S1-051218-480	BM-S1-051218-489	BM-V2-131218-144
Date Sampled	12/05/2018	12/05/2018	12/13/2018

Magnetic Susceptibility Analysis

Magnetic Susceptibility (m3/kg)	2.5E-6 ± 4.4E-8	1.9E-6 ± 4.0E-8	5.5E-7 ± 1.1E-8
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Sample Location	BM-V2-141218-215	BM-V1-171218-122	BM-V1-171218-161
Date Sampled	12/14/2018	12/17/2018	12/17/2018

Magnetic Susceptibility Analysis

Magnetic Susceptibility (m3/kg)	3.1E-6 ± 7.2E-8	1.0E-6 ± 1.2E-8	1.5E-7 ± 3.2E-9
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* Analysis performed in triplicate and results reported as the mean followed by +/- standard deviation.

References

- Brown, R. A., J. T. Wilson and M. Ferrey (2007). "Monitored natural attenuation forum: The case for abiotic MNA." Remediation Journal **17**(2): 127-137.
- Ferrey, M. L., R. T. Wilkin, R. G. Ford and J. T. Wilson (2004). "Nonbiological Removal of cis-Dichloroethylene and 1,1-Dichloroethylene in Aquifer Sediment Containing Magnetite." Environmental Science & Technology **38**(6): 1746-1752.
- He, Y., C. Su, J. T. Wilson, R. T. Wilkin, C. Adair, T. Lee, P. Bradley and M. Ferrey (2009). Identification and characterization of methods for reactive minerals responsible for natural attenuation of chlorinated organic compounds in ground water, US EPA.
- Liu, Y., S. A. Majetich, R. D. Tilton, D. S. Sholl and G. V. Lowry (2005). "TCE Dechlorination Rates, Pathways, and Efficiency of Nanoscale Iron Particles with Different Properties." Environmental Science & Technology **39**(5): 1338-1345.
- Song, H. and E. R. Carraway (2005). "Reduction of Chlorinated Ethanes by Nanosized Zero-Valent Iron: Kinetics, Pathways, and Effects of Reaction Conditions." Environmental Science & Technology **39**(16): 6237-6245.

Abiotic Reactions of Chlorinated Compounds with Iron Bearing Minerals and Zero Valent Iron (ZVI). Summaries for iron bearing minerals are based on He et al. (2009) and references therein. He et al. available at <http://nepis.epa.gov/>. Summary of ZVI based on Liu et al. (2005) and Song et al. (2005).

Contaminant	Mineral	Degradation	Reported Degradation Intermediates and Products ¹
PCE	FeS	Yes	Acetylene, TCE, cis-DCE, 1,1-DCE, ethene
	Pyrite	Yes	TCE, acetylene, ethene
	Magnetite	Yes	Unknown ²
	³ GR(SO ₄) phyllosilicate clays	Reports Differ	
TCE	ZVI	Yes	TCE, 1,1-DCE, vinyl chloride, 1,1,2-TCA, 1,1-DCA, chloroacetylene, acetylene, ethene, ethane
	FeS	Yes	Ethene and ethane
	Pyrite	Yes	Acetylene, cis-DCE, vinyl chloride, 1,1-DCE
	Magnetite	Yes	Acetylene, ethene, cis-DCE, (organic acids with DO present)
cis-DCE	GR(SO ₄), GR(CO ₃)	No	Unknown ¹
	phyllosilicate clays	Yes	Only observed degradation when Cu(II) added
	ZVI	Yes	cDCE, vinyl chloride, acetylene, ethene, ethane
	FeS	No	Ethane, ethene, acetylene with minor amounts of DCE, VC depending on conditions
	Pyrite	Yes	None detected
	Magnetite	Yes	Acetylene, ethene
Vinyl chloride	GR(SO ₄)	Yes	Unknown ²
	phyllosilicate clays	Yes	
	ZVI	Yes	Primarily acetylene and ethene but also much lesser amounts of ethane and VC and traces of methane, propane, propene, butane and butene
	FeS	Unknown	
	Pyrite	Yes	Ethene, ethane
	Magnetite	Yes	Unknown ²
	GR(SO ₄)	Yes	
	phyllosilicate clays	Yes	
	ZVI	Yes	Ethene, ethane, (no evidence of acetylene)

Contaminant	Mineral	Degradation	Reported Degradation Intermediates and Products ¹
1,1-DCA	FeS	Not Significant	None detected
1,1-DCA	GR(SO ₄)	Low conversion	Ethane and ethane (w/ Cu or Ag)
1,1-DCA	ZVI	Yes (low)	Ethane
1,2-DCA	FeS	Not Significant	None detected
1,2-DCA	FeS (Biogenic)	Yes	Not monitored
1,2-DCA	GR(SO ₄)	No	
1,2-DCA	ZVI	No	
1,1,1-TCA	FeS	Yes	1,1-DCA, ethene, 2-butyne
1,1,1-TCA	GR(SO ₄)	Yes	1,1-DCA, CA, ethene ethane
1,1,1-TCA	ZVI	Yes	1,1-DCA, ethane
1,1,2-TCA	FeS	Rate not significant	Small amounts of 1,1-DCE and vinyl chloride but rate not significant
1,1,2-TCA	GR(SO ₄)	Yes	Vinyl chloride, 1,1-DCE, ethene, ethane
1,1,2-TCA	ZVI	Yes	Ethane
1,1,1,2-TeCA	FeS	Yes	1,1-DCE
1,1,1,2-TeCA	GR(SO ₄)	Yes	1,1-DCE and minor (<1%) vinyl chloride, ethene, ethane
1,1,1,2-TeCA	phyllosilicate clays	Yes	1,1-DCE
1,1,1,2-TeCA	ZVI	Yes	TCE, 1,1-DCE
1,1,2,2-TeCA	FeS	Yes	TCE, cis-DCE, trans-DCE, acetylene
1,1,2,2-TeCA	GR(SO ₄)	Yes	TCE (major), cis-DCE, trans-DCE
1,1,2,2-TeCA	phyllosilicate clays	Yes	TCE
1,1,2,2-TeCA	ZVI	Yes	TCE, trans-DCE, cis-DCE
Carbon Tetrachloride	FeS	Yes	Chloroform, carbon disulfide, possibly methane, ethene, ethane
CT	Pyrite	Yes	Chloroform, CO ₂ , carbon disulfide, formate (highly dependent on conditions)
CT	Magnetite	Yes	Chloroform, carbon monoxide, methane, formate (highly dependent on conditions)
CT	GR(SO ₄)	Yes	Chloroform and hexachloroethane; Chloroform, DCM, methane, ethene
CT	phyllosilicate clays	Yes	Chloroform
CT	ZVI	Yes	Chloroform, dichloromethane, methane (depending on conditions)

Notes: GR(SO₄) sulfate green rust. GR(CO₃) carbonate green rust. ZVI zero valent iron

¹Compilation of reported degradation products. Mass recovery of products typically low --- additional undetected and unreported products are likely. Reported reaction products or proportions of reaction products were often a function of environmental conditions.

²No published studies that identify the transformation products of PCE, TCE, cis-DCE or vinyl chloride with magnetite. Ferrey et al (2004) analyzed for products of cis-DCE dechlorination including vinyl chloride, ethene, and ethane and did not find them. If Fe²⁺ sorbed to magnetite stabilizes carbene ions, the ultimate degradation product of cis-DCE on magnetite would be CO₂.

SITE LOGIC Report

Abiotic Potential

Contact: Pamela J. Moss
Address: EA Engineering
7995 E. Prentice Ave
Suite 206E
Greenwood Village, CO 80111

Phone: 303-590-9143

Email: pmoss@eaest.com

MI Identifier: 028PL

Report Date: January 15, 2019

Comments: KAFB-Bulk Fuels Fac Vadose Zone, 62735DM02

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Overview

Although not always fully considered, abiotic degradation can be a substantial or even the primary process for chlorinated hydrocarbon destruction at sites undergoing or transitioning to monitored natural attenuation (MNA). A variety of iron-bearing minerals including iron sulfides (mackinawite and pyrite), iron oxides (magnetite), green rust, and iron-bearing clays are capable of complete or nearly complete degradation of PCE, TCE, and carbon tetrachloride (He et al. 2009). Some iron-bearing minerals also catalyze the degradation of chlorinated ethanes and the lesser chlorinated ethenes cis-dichloroethene (DCE) and vinyl chloride. While the quantities and types will vary, these reactive iron minerals are frequently identified in subsurface environments under iron reducing and sulfate reducing conditions.

Brown et al. (2007) recommend four avenues for evaluating the role of abiotic processes in contaminant attenuation. First, examining contaminant concentrations along the flow path - decreasing parent compound concentrations with no evidence of accumulation of chlorinated transformation products like cis-DCE and vinyl chloride suggest abiotic degradation. Performing compound specific isotope analysis (CSIA) or monitoring for products unique to abiotic reactions such as acetylene can also provide a strong line of evidence. Microcosm studies with native sediment and killed controls can also be performed. Finally, Brown et al. (2007) suggest performing mineralogical analyses on aquifer sediment to characterize reactive minerals such as magnetite or iron monosulfides.

Magnetic Susceptibility - Magnetite

Magnetite (Fe_3O_4) is a mixed valence iron mineral shown to react with PCE, TCE, and carbon tetrachloride. Furthermore, Ferrey et al. (2004) conclusively linked the observed degradation of cis-DCE at a former ammunition plant to magnetite in the subsurface. No direct chemical test is available for quantification of magnetite. However, magnetite is the most abundant mineral in natural sediments that exhibits magnetic behavior. Therefore, magnetic susceptibility provides an inexpensive and valuable estimate of the quantity of magnetite in environmental samples.

X-ray Diffraction (XRD) - Mackinawite, Pyrite, Magnetite and Green Rust

XRD is one of the primary techniques used to identify unknown crystalline materials. Most minerals are crystalline and will scatter X-rays in a regular, characteristic manner dependent on their crystal structure.

- Mackinawite is the most reactive of the iron-bearing minerals and a crystalline form (tetragonal FeS) can be detected by XRD. Mackinawite will transform PCE and TCE primarily by elimination to acetylene. Carbon tetrachloride is transformed mainly to chloroform but carbon dioxide, formate, and carbon disulfide have also been detected. Finally, the more heavily chlorinated ethanes including hexachloroethane, pentachloroethane, and tetrachloroethanes react to form chlorinated ethenes which can be further degraded.
- Pyrite (FeS_2) catalyzes beta elimination transforming PCE, TCE, and cis-DCE to acetylene and ethene. Vinyl chloride is transformed to ethene and ethane. Pyrite is also capable of degradation of carbon tetrachloride potentially forming a number of products including chloroform, carbon dioxide, carbon disulfide, and formate depending on reaction conditions.
- While not quantitative like the magnetic susceptibility test, XRD can also detect magnetite when present at between 2% and 5% on a weight basis.
- Green rusts have been reported to transform a number of common chlorinated contaminants including cis-DCE, vinyl chloride, trichloroethanes, and tetrachloroethanes. While special sample care to prevent oxidation would be needed, XRD can be used to detect green rust.

Results

Table 1. Summary of X-Ray Diffraction and X-ray fluorescence results.

Sample Location	BM-S1-051218-480	BM-S1-051218-489	BM-V2-131218-144
Sample Type	Solid	Solid	Solid
Sample Date	12/05/2018	12/05/2018	12/13/2018
MI ID	028PL1	028PL2	028PL3
X-Ray Diffraction Results			
Mineral Constituent	Relative Abundance (%)		
Quartz	47	40	26.5
Plagioclase Feldspar	30	34.5	13
Microcline	10	14	6
Calcite	1	ND	34
Magnetite	4	2.5	2.5
Kaolinite	0.5	0.5	1
Chlorite	ND	ND	<0.5
Illite/Mica	0.5	0.5	3
Montmorillonite	7	8	ND
Mixed-layered Illite/ Smectite	ND	ND	14
Total	100	100	100
% Illite Layers in ML Illite/Smectite	ND	ND	60.0
X-Ray Fluorescence Results			
Compound	Results (mass %)		
MgO	0.2096	0.2686	0.9677
Al ₂ O ₃	8.6217	8.0828	6.3591
SiO ₂	74.6048	76.7199	49.7724
P ₂ O ₅	0.9668	0.9646	0.8553
SO ₃	ND	ND	0.1172
K ₂ O	6.3338	5.9193	3.9246
CaO	3.6131	3.9177	32.5081
TiO ₂	0.4273	0.4601	0.5222
MnO	0.0944	0.0661	0.2670
Fe ₂ O ₃	4.5467	3.1964	4.4103
ZnO	ND	ND	ND
Rb ₂ O	0.0227	0.0195	0.0287
SrO	0.0742	0.0842	0.1562
Y ₂ O ₃	ND	ND	ND
ZrO ₂	0.0352	0.0191	0.1111
Nb ₂ O ₅	ND	ND	ND
Ag ₂ O	ND	ND	ND
BaO	ND	0.2817	ND

Legend: J = Estimated result below PQL but above LQL ND = Result not detected < = Below detection limit

Table 2. Summary of X-Ray Diffraction and X-ray fluorescence results continued.

Sample Location	BM-V2-141218-215	BM-V1-171218-122	BM-V1-171218-161
Sample Type	Solid	Solid	Solid
Sample Date	12/14/2018	12/17/2018	12/17/2018
MI ID	028PL4	028PL5	028PL6
X-Ray Diffraction Results			
Mineral Constituent	Relative Abundance (%)		
Quartz	39	32	19
Plagioclase Feldspar	34	10	3
Microcline	13	6	3
Calcite	ND	4	12
Magnetite	5	4	6
Kaolinite	<0.5	ND	10
Chlorite	ND	ND	ND
Illite/Mica	<0.5	3	5
Montmorillonite	ND	ND	ND
Mixed-layered Illite/ Smectite	9	41	42
Total	100	100	100
% Illite Layers in ML Illite/Smectite	30.0%	30.0%	30.0%
X-Ray Fluorescence Results			
Compound	Results (mass %)		
MgO	0.4076	1.2093	1.0533
Al ₂ O ₃	8.3962	11.9004	14.7724
SiO ₂	75.4545	68.4083	57.4354
P ₂ O ₅	1.1154	1.1017	1.0820
SO ₃	0.0565	0.0834	ND
K ₂ O	5.2092	5.2766	5.1542
CaO	3.1812	5.6122	10.0372
TiO ₂	0.4674	0.8061	1.3946
MnO	0.0986	0.1384	0.2202
Fe ₂ O ₃	5.2604	5.1679	8.6051
ZnO	ND	0.0242	ND
Rb ₂ O	0.0164	0.0257	0.0371
SrO	0.0802	0.0870	0.0608
Y ₂ O ₃	ND	ND	ND
ZrO ₂	0.0281	0.1460	0.0398
Nb ₂ O ₅	ND	0.0127	ND
Ag ₂ O	ND	ND	0.1080
BaO	0.2283	ND	ND

Legend: J = Estimated result below PQL but above LQL ND = Result not detected < = Below detection limit

References

- Brown, R. A., J. T. Wilson and M. Ferrey (2007). "Monitored natural attenuation forum: The case for abiotic MNA." Remediation Journal **17**(2): 127-137.
- Ferrey, M. L., R. T. Wilkin, R. G. Ford and J. T. Wilson (2004). "Nonbiological Removal of cis-Dichloroethylene and 1,1-Dichloroethylene in Aquifer Sediment Containing Magnetite." Environmental Science & Technology **38**(6): 1746-1752.
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- Liu, Y., S. A. Majetich, R. D. Tilton, D. S. Sholl and G. V. Lowry (2005). "TCE Dechlorination Rates, Pathways, and Efficiency of Nanoscale Iron Particles with Different Properties." Environmental Science & Technology **39**(5): 1338-1345.
- Song, H. and E. R. Carraway (2005). "Reduction of Chlorinated Ethanes by Nanosized Zero-Valent Iron: Kinetics, Pathways, and Effects of Reaction Conditions." Environmental Science & Technology **39**(16): 6237-6245.

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Contaminant	Mineral	Degradation	Reported Degradation Intermediates and Products ¹
PCE	FeS	Yes	Acetylene, TCE, cis-DCE, 1,1-DCE, ethene
	Pyrite	Yes	TCE, acetylene, ethene
	Magnetite	Yes	Unknown ²
	³ GR(SO ₄)	Reports Differ	
	phyllosilicate clays	Yes	TCE, 1,1-DCE, vinyl chloride, 1,1,2-TCA, 1,1-DCA, chloroacetylene, acetylene, ethene, ethane
	ZVI	Yes	Ethene and ethane
TCE	FeS	Yes	Acetylene, cis-DCE, vinyl chloride, 1,1-DCE
	Pyrite	Yes	Acetylene, ethene, cis-DCE, (organic acids with DO present)
	Magnetite	Yes	Unknown ¹
	GR(SO ₄), GR(CO ₃)	No	Only observed degradation when Cu(II) added
	phyllosilicate clays	Yes	cDCE, vinyl chloride, acetylene, ethene, ethane
	ZVI	Yes	Ethane, ethene, acetylene with minor amounts of DCE, VC depending on conditions
cis-DCE	FeS	No	None detected
	Pyrite	Yes	Acetylene, ethene
	Magnetite	Yes	Unknown ²
	GR(SO ₄)	Yes	
	phyllosilicate clays	Yes	
	ZVI	Yes	Primarily acetylene and ethene but also much lesser amounts of ethane and VC and traces of methane, propane, propene, butane and butene
Vinyl chloride	FeS	Unknown	
	Pyrite	Yes	Ethene, ethane
	Magnetite	Yes	Unknown ²
	GR(SO ₄)	Yes	
	phyllosilicate clays	Yes	
	ZVI	Yes	Ethene, ethane, (no evidence of acetylene)

Contaminant	Mineral	Degradation	Reported Degradation Intermediates and Products ¹
1,1-DCA	FeS	Not Significant	None detected
1,1-DCA	GR(SO ₄)	Low conversion	Ethene and ethane (w/ Cu or Ag)
1,1-DCA	ZVI	Yes (low)	Ethane
1,2-DCA	FeS	Not Significant	None detected
1,2-DCA	FeS (Biogenic)	Yes	Not monitored
1,2-DCA	GR(SO ₄)	No	
1,2-DCA	ZVI	No	
1,1,1-TCA	FeS	Yes	1,1-DCA, ethene, 2-butyne
1,1,1-TCA	GR(SO ₄)	Yes	1,1-DCA, CA, ethene ethane
1,1,1-TCA	ZVI	Yes	1,1-DCA, ethane
1,1,2-TCA	FeS	Rate not significant	Small amounts of 1,1-DCE and vinyl chloride but rate not significant
1,1,2-TCA	GR(SO ₄)	Yes	Vinyl chloride, 1,1-DCE, ethene, ethane
1,1,2-TCA	ZVI	Yes	Ethane
1,1,1,2-TeCA	FeS	Yes	1,1-DCE
1,1,1,2-TeCA	GR(SO ₄)	Yes	1,1-DCE and minor (<1%) vinyl chloride, ethene, ethane
1,1,1,2-TeCA	phyllosilicate clays	Yes	1,1-DCE
1,1,1,2-TeCA	ZVI	Yes	TCE, 1,1-DCE
1,1,2,2-TeCA	FeS	Yes	TCE, cis-DCE, trans-DCE, acetylene
1,1,2,2-TeCA	GR(SO ₄)	Yes	TCE (major), cis-DCE, trans-DCE
1,1,2,2-TeCA	phyllosilicate clays	Yes	TCE
1,1,2,2-TeCA	ZVI	Yes	TCE, trans-DCE, cis-DCE
Carbon Tetrachloride	FeS	Yes	Chloroform, carbon disulfide, possibly methane, ethene, ethane
CT	Pyrite	Yes	Chloroform, CO ₂ , carbon disulfide, formate (highly dependent on conditions)
CT	Magnetite	Yes	Chloroform, carbon monoxide, methane, formate (highly dependent on conditions)
CT	GR(SO ₄)	Yes	Chloroform and hexachloroethane; Chloroform, DCM, methane, ethene
CT	phyllosilicate clays	Yes	Chloroform
CT	ZVI	Yes	Chloroform, dichloromethane, methane (depending on conditions)

Notes: GR(SO₄) sulfate green rust. GR(CO₃) carbonate green rust. ZVI zero valent iron

¹Compilation of reported degradation products. Mass recovery of products typically low - additional undetected and unreported products are likely. Reported reaction products or proportions of reaction products were often a function of environmental conditions.

²No published studies that identify the transformation products of PCE, TCE, cis-DCE or vinyl chloride with magnetite. Ferrey et al (2004) analyzed for products of cis-DCE dechlorination including vinyl chloride, ethene, and ethane and did not find them. If Fe²⁺ sorbed to magnetite stabilizes carbene ions, the ultimate degradation product of cis-DCE on magnetite would be CO₂.

2021-22-III-20181213



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REPORT TO:

Name: P. Moss
Company: EA Engineering, Science & Technology, Inc.
Address: 7995 E Prentic Ave.
Greenwood Village CO 80111
email: p.moss@eaest.com
Phone: 303-590-9143
Fax:

Project Manager:	Bernie Bockisch
Project Name:	Kirtland Vadose Zone Coring
Project No.:	62735DM02.1017

Report Type: ☒ Standard (default) ☐ Microbial Insights Level III raw data

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Dallas Accounts Payable via Email below

Company: EA Engineering, Science & Technology, Inc.

Address:

EA Engineering, Science & Technology, Inc.
7995 E Prentice Ave.
Greenwood Village CO 80111

p.moss@eaast.com

303-590-9143

Bernie Bockisch

Kirtland Vadose Zone Coring

62735DM02.1017

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MI Quote No. Q2018628.0001

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10515 Research Dr
Knoxville, TN 37932
865-573-8188

www.microbe.com

dallasan@east.com

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REPORT TO:

Name: P. Moss
Company: EA Engineering, Science & Technology, Inc.
Address: 7995 E Prentic Ave.
Greenwood Village CO 80111
p.moss@eaest.com
Phone: 303-590-9143
Fax:

Project Manager:	<u>Bernie Bockisch</u>
Project Name:	<u>Kirtland Vadose Zone Coring</u>
Project No.:	<u>62735DM02.1017</u>

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EA Engineering, Science & Technology, Inc.

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Laboratory Report for EA Engineering, Science and Technology, Inc.

KAFB-Bulk Fuels Fac Vadose Zone

Project #: 62735DM02, PO: 18189

January 11, 2019



Daniel B. Stephens & Associates, Inc.

4400 Alameda Blvd. NE, Suite C • Albuquerque, New Mexico 87113



January 11, 2019

Bernie Bockisch
EA Engineering, Science and Technology, Inc.
320 Gold Ave SW, Ste. 1300
Albuquerque, NM 87102
(505) 234-1105

Re: DBS&A Laboratory Report for the EA Engineering, Science, and Technology, Inc.
62735DM02 KAFB-Bulk Fuels Fac Vadose Zone Project

Dear Mr. Bockisch:

Enclosed is the report for the EA Engineering, Science, and Technology, Inc. 62735DM02 KAFB-Bulk Fuels Fac Vadose Zone project samples. Please review this report and provide any comments as samples will be held for a maximum of 30 days. After 30 days samples will be returned or disposed of in an appropriate manner.

All testing results were evaluated subjectively for consistency and reasonableness, and the results appear to be reasonably representative of the material tested. However, DBS&A does not assume any responsibility for interpretations or analyses based on the data enclosed, nor can we guarantee that these data are fully representative of the undisturbed materials at the field site. We recommend that careful evaluation of these laboratory results be made for your particular application.

The testing utilized to generate the enclosed report employs methods that are standard for the industry. The results do not constitute a professional opinion by DBS&A, nor can the results affect any professional or expert opinions rendered with respect thereto by DBS&A. You have acknowledged that all the testing undertaken by us, and the report provided, constitutes mere test results using standardized methods, and cannot be used to disqualify DBS&A from rendering any professional or expert opinion, having waived any claim of conflict of interest by DBS&A.

We are pleased to provide this service to EA Engineering and look forward to future laboratory testing on other projects. If you have any questions about the enclosed data, please do not hesitate to call.

Sincerely,

DANIEL B. STEPHENS & ASSOCIATES, INC.
SOIL TESTING & RESEARCH LABORATORY

Joleen Hines
Laboratory Manager

Enclosure

Daniel B. Stephens & Associates, Inc.
Soil Testing & Research Laboratory

4400 Alameda Blvd. NE, Suite C
Albuquerque, NM 87113

505-889-7752
FAX 505-889-0258

Summaries



Summary of Tests Performed

Laboratory Sample Number	Initial Soil Properties ¹			Saturated Hydraulic Conductivity ²			Moisture Characteristics ³								Particle Size ⁴			Specific Gravity ⁵		Air Perm- eability	Atterberg Limits	Thermal Properties
	G	VM	VD	CH	FH	FW	HC	PP	FP	DPP	RH	EP	WHC	K _{unsat}	DS	WS	H	F	C			
T-V2-131218-144 Clay	X	X																				X
T-V2-131218-144 Sand	X	X																				X

¹ G = Gravimetric Moisture Content, VM = Volume Measurement Method, VD = Volume Displacement Method

² CH = Constant Head Rigid Wall, FH = Falling Head Rigid Wall, FW = Falling Head Rising Tail Flexible Wall

³ HC = Hanging Column, PP = Pressure Plate, FP = Filter Paper, DPP = Dew Point Potentiometer, RH = Relative Humidity Box,
EP = Effective Porosity, WHC = Water Holding Capacity, K_{unsat} = Calculated Unsaturated Hydraulic Conductivity

⁴ DS = Dry Sieve, WS = Wet Sieve, H = Hydrometer

⁵ F = Fine (<4.75mm), C = Coarse (>4.75mm)



Notes

Sample Receipt:

One sample, in a full 6" x 24" plastic bag sealed with duct tape, was hand-delivered on December 13, 2018. The sample was received inside a cardboard box with a lid and was received in good order.

Sample Preparation and Testing Notes:

Two test samples were prepared, one from the more clay portion and one from the more sand portion. The clay sub-sample was prepared by remolding the material to target the density as measured from a small clay clod. The sand sub-sample was prepared by remolding the material to target a moderate compactive state, approximating 85% of standard proctor compaction maximum dry bulk density, based on technician experience and judgement. Both sub-samples were remolded using material near the initial as received moisture content. The remolded sub-samples were subjected to thermal properties testing at the initial moisture content, the saturated moisture content, at several air drying moisture contents, and at the oven dry state. Each thermal properties reading was obtained in the same location.

Porosity calculations are based on the use of an assumed specific gravity value of 2.65.

Volumetric water contents were adjusted for changes in volume, where applicable. Due to the irregularities formed on the sample surfaces during settling or swelling, volume measurements obtained after the initial reading should be considered estimates.



Summary of Sample Preparation/Volume Changes

Sample Number	Target Remold Parameters ¹		Actual Remold Data			Volume Change Post Saturation ²		
	Moisture Content (%, g/g)	Dry Bulk Density (g/cm ³)	Moisture Content (%, g/g)	Dry Bulk Density (g/cm ³)	% of Target Density (%)	Dry Bulk Density (g/cm ³)	% Volume Change (%)	% of Initial Density (%)
T-V2-131218-144 Clay	20.3	1.64	20.0	1.64	100.3%	1.64	---	100%
T-V2-131218-144 Sand	As Received	'Moderate'	6.5	1.72	NA	1.72	---	100%

¹Target Remold Parameters: Provided by the client. See "Notes" page.

²Volume Change Post Saturation: Volume change measurements were obtained after saturated hydraulic conductivity testing.

Notes:

"+" indicates sample swelling, "-" indicates sample settling, and "---" indicates no volume change occurred.



**Summary of Initial Moisture Content, Dry Bulk Density
Wet Bulk Density and Calculated Porosity**

Sample Number	Moisture Content				Dry Bulk Density (g/cm ³)	Wet Bulk Density (g/cm ³)	Calculated Porosity (%)
	As Received		Remolded				
	Gravimetric (%, g/g)	Volumetric (%, cm ³ /cm ³)	Gravimetric (%, g/g)	Volumetric (%, cm ³ /cm ³)			
T-V2-131218-144 Clay	20.3	33.3	20.0	32.8	1.64	1.97	37.9
T-V2-131218-144 Sand	NA	NA	6.5	11.2	1.72	1.83	35.0

NA = Not analyzed

--- = This sample was not remolded



Summary of Thermal Properties

Sample	Reading	Gravimetric Moisture Content (g/g, %)	Volumetric Moisture Content ¹ (vol/vol, %)	Dry Bulk Density ¹ (g/cm ³)	Temp °C	K W/(m·K)	ρ °C·cm/W	C MJ/(m ³ ·K)	D mm ² /s
T-V2-131218-144 Clay	Initial	19.96	32.83	1.64	20.08	1.552	64.4	1.939	0.800
T-V2-131218-144 Clay	0	23.41	38.50	1.64	18.31	1.686	59.3	3.865	0.436
T-V2-131218-144 Clay	Air Dry #1	18.89	31.07	1.64	15.66	1.583	63.2	1.947	0.813
T-V2-131218-144 Clay	Air Dry #2	15.48	27.01	1.74	17.68	1.411	70.9	1.766	0.799
T-V2-131218-144 Clay	Air Dry #3	11.40	20.60	1.81	17.57	1.252	79.9	1.568	0.798
T-V2-131218-144 Clay	Air Dry #4	8.94	16.20	1.81	19.71	1.110	90.1	1.418	0.783
T-V2-131218-144 Clay	Air Dry #5	5.79	10.48	1.81	18.69	0.923	108.3	1.313	0.703
T-V2-131218-144 Clay	Air Dry #6	4.33	7.85	1.81	19.01	0.804	124.3	1.187	0.678
T-V2-131218-144 Clay	Oven Dry	0	0	1.81	21.76	0.524	190.9	1.118	0.468
T-V2-131218-144 Sand	Initial	6.49	11.18	1.72	20.47	0.862	116.0	1.579	0.546
T-V2-131218-144 Sand	0	20.43	35.17	1.72	18.21	1.838	54.4	2.756	0.667
T-V2-131218-144 Sand	Air Dry #1	15.23	26.23	1.72	16.02	1.462	68.4	2.140	0.683
T-V2-131218-144 Sand	Air Dry #2	8.51	14.66	1.72	17.91	1.068	93.6	1.738	0.615
T-V2-131218-144 Sand	Air Dry #3	5.63	9.70	1.72	18.74	0.753	132.8	1.307	0.576
T-V2-131218-144 Sand	Air Dry #4	4.51	7.77	1.72	20.41	0.737	135.7	1.516	0.486
T-V2-131218-144 Sand	Air Dry #5	3.03	5.21	1.72	19.19	0.480	208.3	1.627	0.295
T-V2-131218-144 Sand	Air Dry #6	2.26	3.90	1.72	19.17	0.434	230.5	1.470	0.295
T-V2-131218-144 Sand	Oven Dry	0	0	1.72	21.86	0.298	335.3	1.299	0.230

¹ Adjusted for volume changes during testing, if applicable.

Initial Properties



**Summary of Initial Moisture Content, Dry Bulk Density
Wet Bulk Density and Calculated Porosity**

Sample Number	Moisture Content				Dry Bulk Density (g/cm ³)	Wet Bulk Density (g/cm ³)	Calculated Porosity (%)
	As Received		Remolded				
	Gravimetric (%, g/g)	Volumetric (%, cm ³ /cm ³)	Gravimetric (%, g/g)	Volumetric (%, cm ³ /cm ³)			
T-V2-131218-144 Clay	20.3	33.3	20.0	32.8	1.64	1.97	37.9
T-V2-131218-144 Sand	NA	NA	6.5	11.2	1.72	1.83	35.0

NA = Not analyzed

--- = This sample was not remolded



**Data for Initial Moisture Content,
Bulk Density, Porosity, and Percent Saturation**

Job Name: EA Engineering, Science, and Technology, Inc.
Job Number: DB18.1333.00
Sample Number: T-V2-131218-144 Clay
Project Name: 62735DM02 KAFB-Bulk Fuels Fac Vadose Zone
PO Number: 18189

	<u>As Received</u>	<u>Remolded</u>
Test Date:	17-Dec-19	18-Dec-19
Field weight* of sample (g):	132.67	571.60
Tare weight, ring (g):	0.00	136.65
Tare weight, pan/plate (g):	0.00	0.00
Tare weight, other (g):	0.00	0.00
Dry weight of sample (g):	110.26	362.58
Sample volume (cm ³):	67.24	220.41
Assumed particle density (g/cm ³):	2.65	2.65
Gravimetric Moisture Content (% g/g):	20.3	20.0
Volumetric Moisture Content (% vol):	33.3	32.8
Dry bulk density (g/cm ³):	1.64	1.64
Wet bulk density (g/cm ³):	1.97	1.97
Calculated Porosity (% vol):	38.1	37.9
Percent Saturation:	87.4	86.6

Laboratory analysis by:	D. O'Dowd	D. O'Dowd
Data entered by:	C. Krous	C. Krous
Checked by:	J. Hines	J. Hines

Comments:

* Weight including tares
NA = Not analyzed
--- = This sample was not remolded



**Data for Initial Moisture Content,
Bulk Density, Porosity, and Percent Saturation**

Job Name: EA Engineering, Science, and Technology, Inc.
Job Number: DB18.1333.00
Sample Number: T-V2-131218-144 Sand
Project Name: 62735DM02 KAFB-Bulk Fuels Fac Vadose Zone
PO Number: 18189

	<u>As Received</u>	<u>Remolded</u>
Test Date:	NA	18-Dec-19
Field weight* of sample (g):		540.04
Tare weight, ring (g):		126.69
Tare weight, pan/plate (g):		0.00
Tare weight, other (g):		0.00
Dry weight of sample (g):		388.15
Sample volume (cm ³):		225.43
Assumed particle density (g/cm ³):		2.65
Gravimetric Moisture Content (% g/g):		6.5
Volumetric Moisture Content (% vol):		11.2
Dry bulk density (g/cm ³):		1.72
Wet bulk density (g/cm ³):		1.83
Calculated Porosity (% vol):		35.0
Percent Saturation:		31.9
Laboratory analysis by:		D. O'Dowd
Data entered by:		C. Krous
Checked by:		J. Hines

Comments:

* Weight including tares
NA = Not analyzed
--- = This sample was not remolded

Thermal Properties



Summary of Thermal Properties

Sample	Reading	Gravimetric Moisture Content (g/g, %)	Volumetric Moisture Content ¹ (vol/vol, %)	Dry Bulk Density ¹ (g/cm ³)	Temp °C	K W/(m·K)	ρ °C·cm/W	C MJ/(m ³ ·K)	D mm ² /s
T-V2-131218-144 Clay	Initial	19.96	32.83	1.64	20.08	1.552	64.4	1.939	0.800
T-V2-131218-144 Clay	0	23.41	38.50	1.64	18.31	1.686	59.3	3.865	0.436
T-V2-131218-144 Clay	Air Dry #1	18.89	31.07	1.64	15.66	1.583	63.2	1.947	0.813
T-V2-131218-144 Clay	Air Dry #2	15.48	27.01	1.74	17.68	1.411	70.9	1.766	0.799
T-V2-131218-144 Clay	Air Dry #3	11.40	20.60	1.81	17.57	1.252	79.9	1.568	0.798
T-V2-131218-144 Clay	Air Dry #4	8.94	16.20	1.81	19.71	1.110	90.1	1.418	0.783
T-V2-131218-144 Clay	Air Dry #5	5.79	10.48	1.81	18.69	0.923	108.3	1.313	0.703
T-V2-131218-144 Clay	Air Dry #6	4.33	7.85	1.81	19.01	0.804	124.3	1.187	0.678
T-V2-131218-144 Clay	Oven Dry	0	0	1.81	21.76	0.524	190.9	1.118	0.468
T-V2-131218-144 Sand	Initial	6.49	11.18	1.72	20.47	0.862	116.0	1.579	0.546
T-V2-131218-144 Sand	0	20.43	35.17	1.72	18.21	1.838	54.4	2.756	0.667
T-V2-131218-144 Sand	Air Dry #1	15.23	26.23	1.72	16.02	1.462	68.4	2.140	0.683
T-V2-131218-144 Sand	Air Dry #2	8.51	14.66	1.72	17.91	1.068	93.6	1.738	0.615
T-V2-131218-144 Sand	Air Dry #3	5.63	9.70	1.72	18.74	0.753	132.8	1.307	0.576
T-V2-131218-144 Sand	Air Dry #4	4.51	7.77	1.72	20.41	0.737	135.7	1.516	0.486
T-V2-131218-144 Sand	Air Dry #5	3.03	5.21	1.72	19.19	0.480	208.3	1.627	0.295
T-V2-131218-144 Sand	Air Dry #6	2.26	3.90	1.72	19.17	0.434	230.5	1.470	0.295
T-V2-131218-144 Sand	Oven Dry	0	0	1.72	21.86	0.298	335.3	1.299	0.230

¹ Adjusted for volume changes during testing, if applicable.



Thermal Properties Results Sheet for Sample: T-V2-131218-144 Clay

Job Name: EA Engineering
 Job Number: DB18.1333.00
 Sample Number: T-V2-131218-144 Clay
 Project Name: 62735DM02 KAFB-Bulk Fuels Fac Vadose Zone
 PO Number: 18189

Instrument Description: Decagon KD2 Pro

Probe: ☐ KS-1, 6 cm length, 1.3 mm diameter, single needle
☐ TR-1, 10 cm length, 2.4 mm diameter, single needle
☒ SH-1, 3 cm length, 1.3 mm diameter, dual needle, 6 mm spacing

Test Start Date: 12/18/18

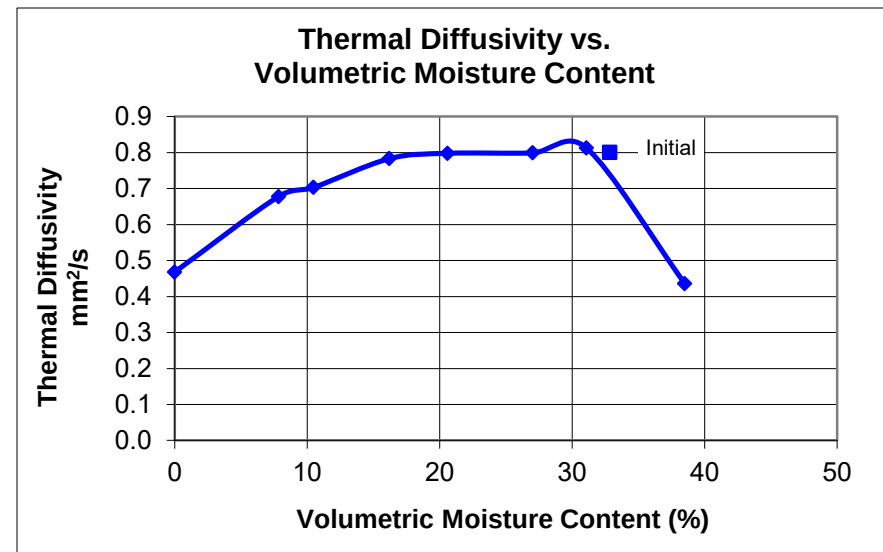
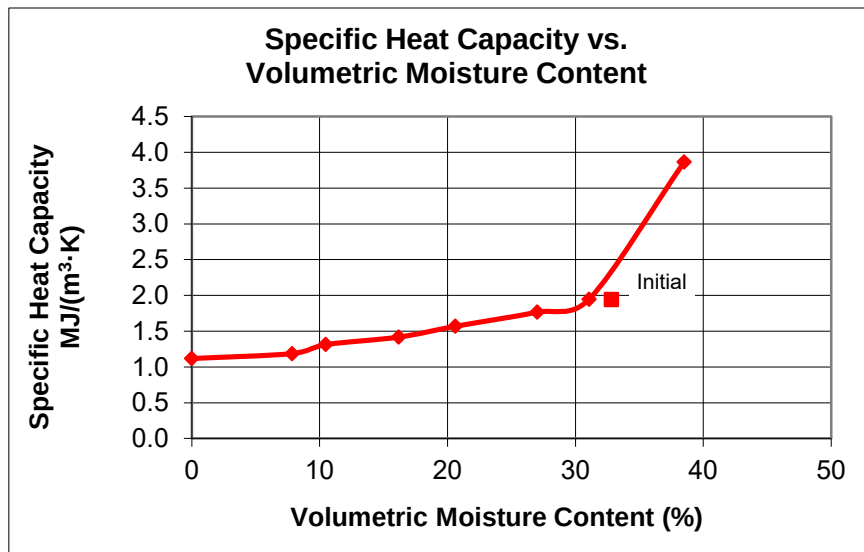
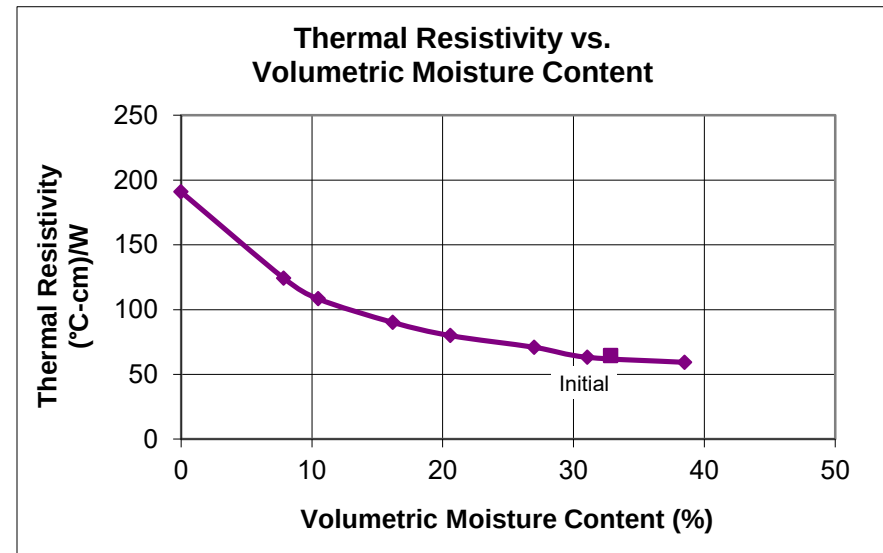
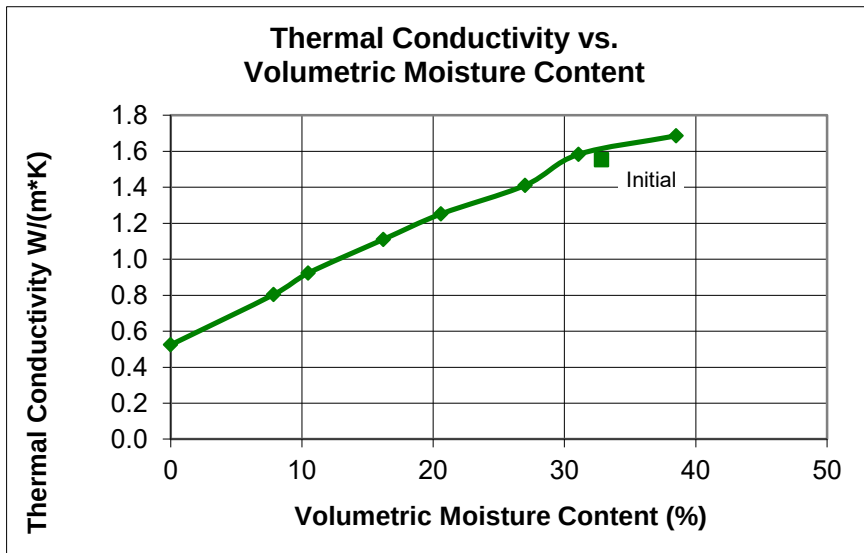
Reading	Water Potential (-cm water)	Gravimetric Moisture Content (g/g, %)	Volumetric Moisture Content ¹ (vol/vol, %)	Dry Bulk Density ¹ (g/cm ³)	Test Temperature (°C)	K Thermal Conductivity W/(m·K)	ρ Thermal Resistivity °C·cm/W	C Specific Heat Capacity MJ/(m ³ ·K)	D Thermal Diffusivity (mm ² /s)
Initial	---	19.96	32.83	1.64	20.08	1.552	64.4	1.939	0.800
Saturated	0	23.41	38.50	1.64	18.31	1.686	59.3	3.865	0.436
Air Dry #1	---	18.89	31.07	1.64	15.66	1.583	63.2	1.947	0.813
Air Dry #2	---	15.48	27.01	1.74	17.68	1.411	70.9	1.766	0.799
Air Dry #3	---	11.40	20.60	1.81	17.57	1.252	79.9	1.568	0.798
Air Dry #4	---	8.94	16.20	1.81	19.71	1.110	90.1	1.418	0.783
Air Dry #5	---	5.79	10.48	1.81	18.69	0.923	108.3	1.313	0.703
Air Dry #6	---	4.33	7.85	1.81	19.01	0.804	124.3	1.187	0.678
Oven Dry	---	0	0	1.81	21.76	0.524	190.9	1.118	0.468

--- = Value not measured.

¹ Adjusted for volume changes during testing, if applicable.



Thermal Properties Results Sheet for Sample: T-V2-131218-144 Clay Scatter Plots





Thermal Properties Data

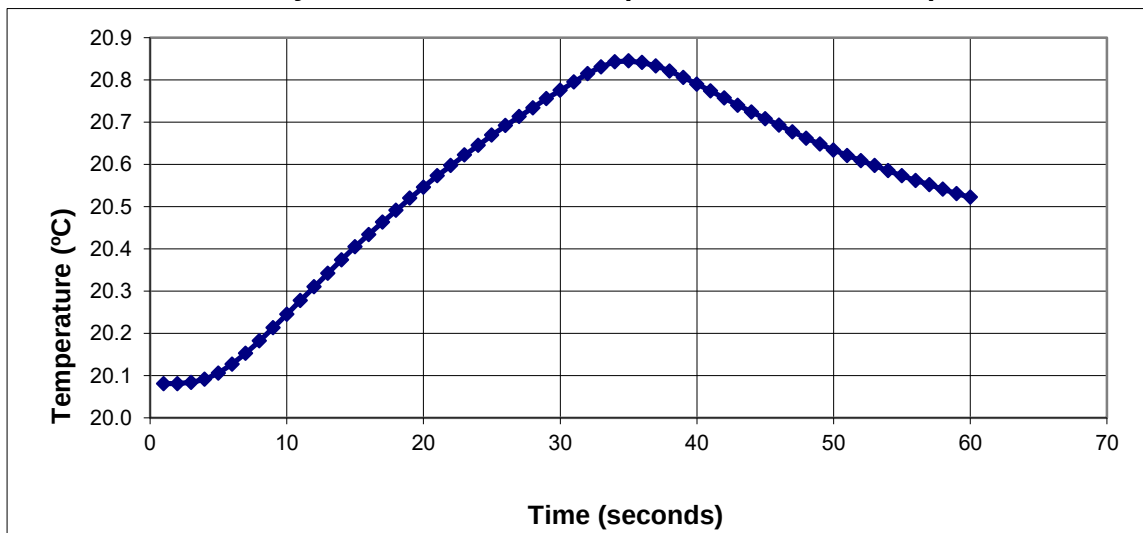
Sample Number: T-V2-131218-144 Clay
Potential (-cm water): Initial

Test Date/Time: 12/18/2018 9:04 AM	K (W/(m·K)): 1.552
Sensor: SH-1	ρ (°C·cm/W): 64.4
Test Temp.(°C): 20.1	C (MJ/(m ³ ·K)): 1.939
KD2 Pro Sample ID: CLAY-IN	D (mm ² /s): 0.800
Power (W/m): 20.610	Err: 0.0036
Current (amps): 0.141	

Raw Data

Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)	
1	20.081	16	20.434	31	20.795	46	20.693
2	20.081	17	20.463	32	20.815	47	20.677
3	20.084	18	20.491	33	20.831	48	20.662
4	20.091	19	20.520	34	20.843	49	20.648
5	20.106	20	20.546	35	20.845	50	20.634
6	20.127	21	20.573	36	20.841	51	20.621
7	20.153	22	20.597	37	20.833	52	20.609
8	20.182	23	20.622	38	20.821	53	20.597
9	20.213	24	20.645	39	20.806	54	20.585
10	20.245	25	20.669	40	20.790	55	20.573
11	20.278	26	20.692	41	20.774	56	20.562
12	20.310	27	20.713	42	20.757	57	20.552
13	20.342	28	20.734	43	20.740	58	20.541
14	20.374	29	20.756	44	20.724	59	20.531
15	20.405	30	20.775	45	20.708	60	20.522

T-V2-131218-144 Clay,Potential: Initial - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

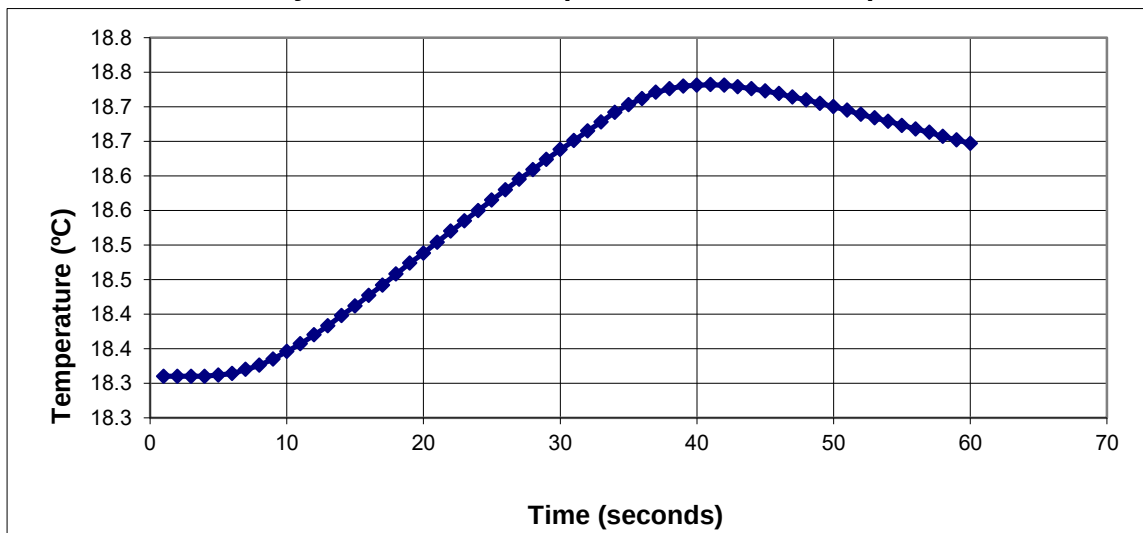
Sample Number: T-V2-131218-144 Clay
Potential (-cm water): 0

Test Date/Time: 12/19/2018 8:42 AM	K (W/(m·K)): 1.686
Sensor: SH-1	ρ (°C·cm/W): 59.3
Test Temp.(°C): 18.3	C (MJ/(m ³ ·K)): 3.865
KD2 Pro Sample ID: CLAY-SA	D (mm ² /s): 0.436
Power (W/m): 20.440	Err: 0.0009
Current (amps): 0.140	

Raw Data

Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)
1 18.310	16 18.427	31 18.651	46 18.719
2 18.310	17 18.442	32 18.665	47 18.714
3 18.310	18 18.458	33 18.678	48 18.710
4 18.310	19 18.474	34 18.692	49 18.705
5 18.312	20 18.488	35 18.703	50 18.700
6 18.314	21 18.504	36 18.712	51 18.695
7 18.320	22 18.520	37 18.721	52 18.689
8 18.326	23 18.535	38 18.726	53 18.684
9 18.335	24 18.550	39 18.730	54 18.679
10 18.346	25 18.565	40 18.731	55 18.673
11 18.357	26 18.580	41 18.732	56 18.668
12 18.370	27 18.595	42 18.731	57 18.663
13 18.383	28 18.609	43 18.729	58 18.657
14 18.398	29 18.624	44 18.726	59 18.652
15 18.412	30 18.638	45 18.723	60 18.647

T-V2-131218-144 Clay,Potential: 0 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

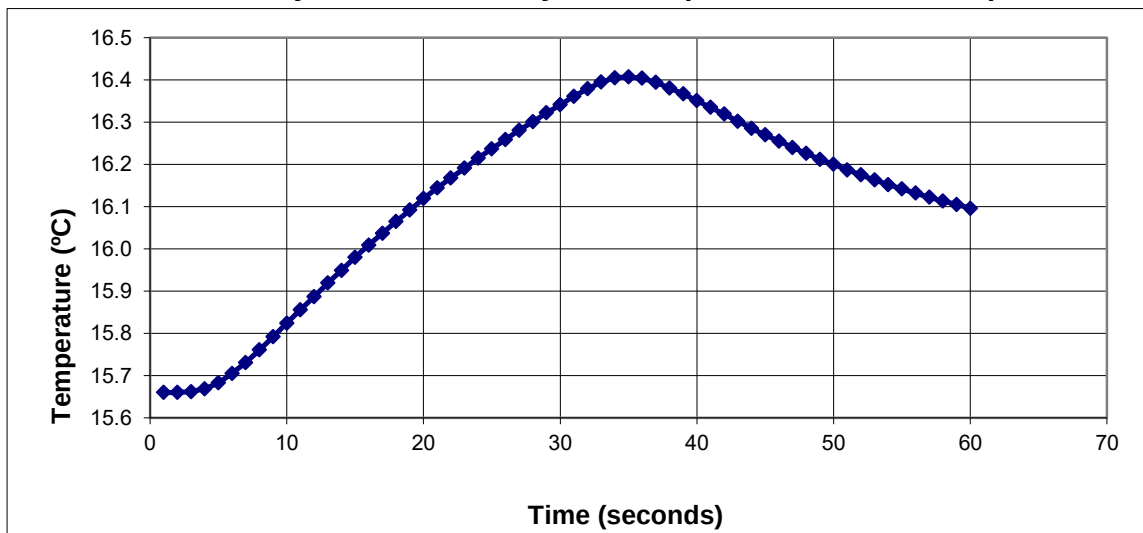
Sample Number: T-V2-131218-144 Clay
Potential (-cm water): Air Dry #1

Test Date/Time: 12/20/2018 10:43 AM	K (W/(m·K)): 1.583
Sensor: SH-1	ρ (°C·cm/W): 63.2
Test Temp.(°C): 15.7	C (MJ/(m ³ ·K)): 1.947
KD2 Pro Sample ID: CLAYAD1	D (mm ² /s): 0.813
Power (W/m): 20.150	Err: 0.0049
Current (amps): 0.139	

Raw Data

Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)
1 15.660	16 16.009	31 16.361	46 16.255
2 15.660	17 16.037	32 16.379	47 16.240
3 15.662	18 16.065	33 16.395	48 16.226
4 15.669	19 16.092	34 16.405	49 16.212
5 15.683	20 16.119	35 16.407	50 16.200
6 15.705	21 16.144	36 16.404	51 16.187
7 15.731	22 16.168	37 16.394	52 16.175
8 15.761	23 16.191	38 16.381	53 16.163
9 15.792	24 16.215	39 16.367	54 16.152
10 15.824	25 16.237	40 16.351	55 16.142
11 15.856	26 16.259	41 16.335	56 16.132
12 15.887	27 16.281	42 16.319	57 16.122
13 15.919	28 16.301	43 16.302	58 16.113
14 15.949	29 16.322	44 16.285	59 16.105
15 15.980	30 16.341	45 16.270	60 16.096

T-V2-131218-144 Clay,Potential: Air Dry #1 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

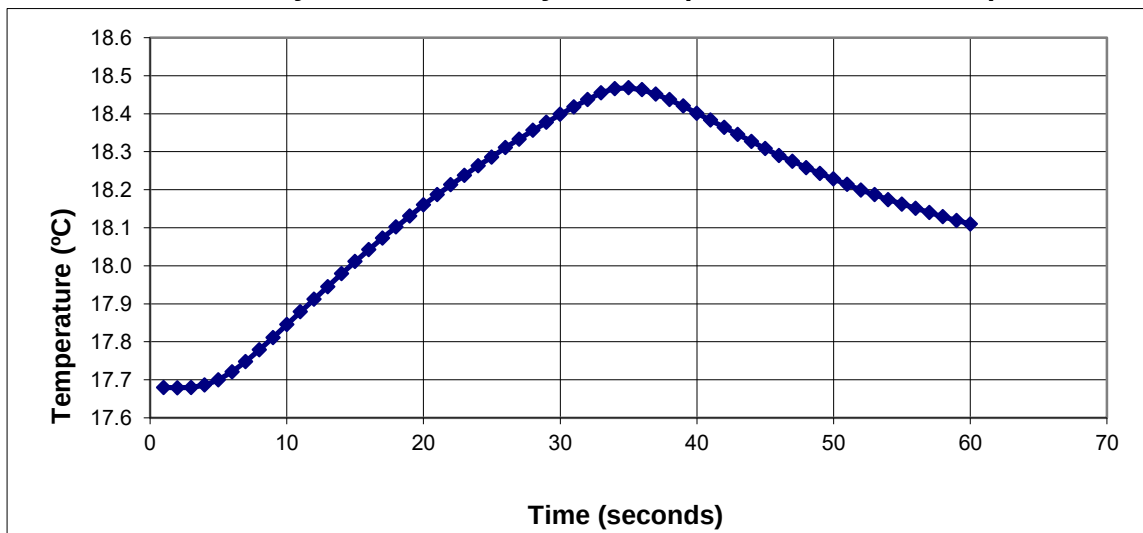
Sample Number: T-V2-131218-144 Clay
Potential (-cm water): Air Dry #2

Test Date/Time: 12/21/2018 1:36 PM	K (W/(m·K)): 1.411
Sensor: SH-1	ρ (°C·cm/W): 70.9
Test Temp.(°C): 17.7	C (MJ/(m ³ ·K)): 1.766
KD2 Pro Sample ID: CLAYAD2	D (mm ² /s): 0.799
Power (W/m): 19.930	Err: 0.0060
Current (amps): 0.138	

Raw Data

Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)
1 17.680	16 18.042	31 18.418	46 18.290
2 17.679	17 18.073	32 18.437	47 18.275
3 17.680	18 18.102	33 18.455	48 18.258
4 17.686	19 18.131	34 18.466	49 18.243
5 17.700	20 18.160	35 18.468	50 18.228
6 17.721	21 18.187	36 18.463	51 18.214
7 17.748	22 18.213	37 18.451	52 18.199
8 17.779	23 18.238	38 18.437	53 18.187
9 17.811	24 18.263	39 18.420	54 18.174
10 17.845	25 18.286	40 18.401	55 18.162
11 17.879	26 18.311	41 18.383	56 18.151
12 17.912	27 18.333	42 18.364	57 18.140
13 17.945	28 18.356	43 18.345	58 18.129
14 17.979	29 18.377	44 18.327	59 18.119
15 18.011	30 18.398	45 18.308	60 18.110

T-V2-131218-144 Clay,Potential: Air Dry #2 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

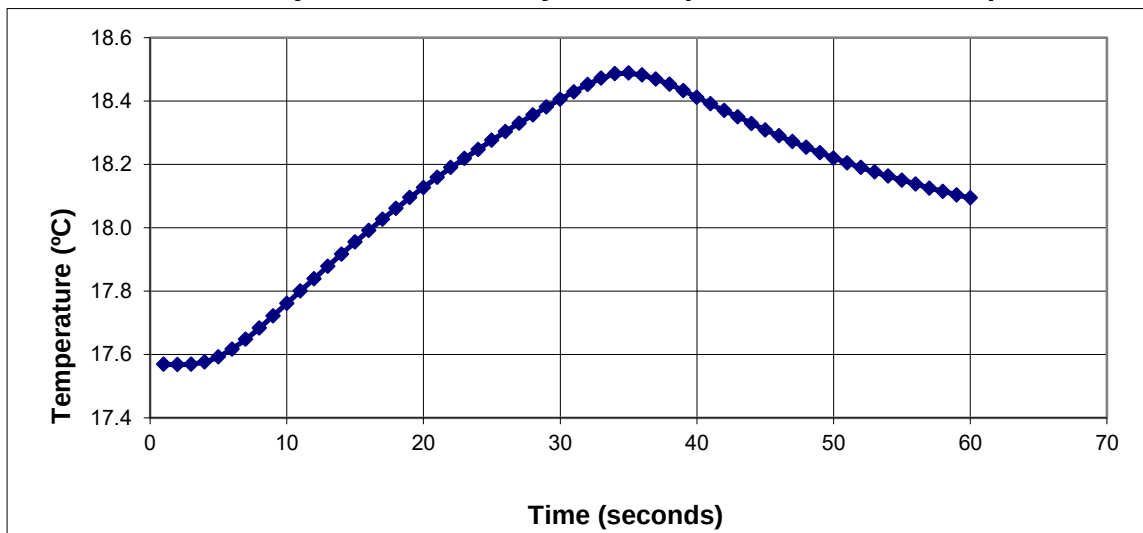
Sample Number: T-V2-131218-144 Clay
Potential (-cm water): Air Dry #3

Test Date/Time: 12/27/2018 8:10 AM	K (W/(m·K)): 1.252
Sensor: SH-1	ρ (°C·cm/W): 79.9
Test Temp.(°C): 17.6	C (MJ/(m ³ ·K)): 1.568
KD2 Pro Sample ID: CLAYAD3	D (mm ² /s): 0.798
Power (W/m): 20.200	Err: 0.0067
Current (amps): 0.139	

Raw Data

Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)
1 17.569	16 17.991	31 18.429	46 18.290
2 17.568	17 18.027	32 18.452	47 18.272
3 17.569	18 18.061	33 18.472	48 18.254
4 17.576	19 18.095	34 18.486	49 18.237
5 17.592	20 18.127	35 18.488	50 18.220
6 17.617	21 18.159	36 18.482	51 18.205
7 17.648	22 18.190	37 18.469	52 18.190
8 17.683	23 18.219	38 18.453	53 18.176
9 17.722	24 18.247	39 18.433	54 18.163
10 17.761	25 18.276	40 18.412	55 18.150
11 17.801	26 18.304	41 18.391	56 18.138
12 17.839	27 18.330	42 18.370	57 18.125
13 17.878	28 18.356	43 18.350	58 18.115
14 17.917	29 18.381	44 18.329	59 18.104
15 17.955	30 18.406	45 18.309	60 18.094

T-V2-131218-144 Clay,Potential: Air Dry #3 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

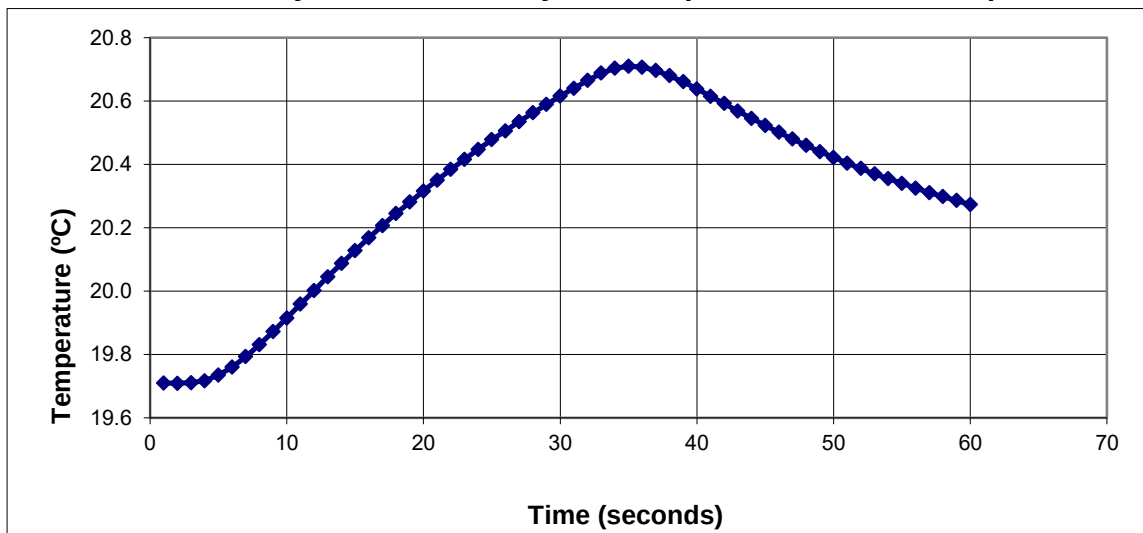
Sample Number: T-V2-131218-144 Clay
Potential (-cm water): Air Dry #4

Test Date/Time: 12/28/2018 2:56 PM	K (W/(m·K)): 1.110
Sensor: SH-1	ρ (°C·cm/W): 90.1
Test Temp.(°C): 19.7	C (MJ/(m ³ ·K)): 1.418
KD2 Pro Sample ID: CLAYAD4	D (mm ² /s): 0.783
Power (W/m): 20.080	Err: 0.0059
Current (amps): 0.139	

Raw Data

Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)
1 19.710	16 20.168	31 20.640	46 20.501
2 19.709	17 20.207	32 20.665	47 20.480
3 19.711	18 20.245	33 20.688	48 20.460
4 19.717	19 20.281	34 20.703	49 20.440
5 19.735	20 20.316	35 20.710	50 20.422
6 19.760	21 20.350	36 20.707	51 20.404
7 19.793	22 20.384	37 20.696	52 20.387
8 19.831	23 20.416	38 20.680	53 20.370
9 19.872	24 20.447	39 20.661	54 20.355
10 19.915	25 20.478	40 20.638	55 20.340
11 19.959	26 20.506	41 20.615	56 20.325
12 20.002	27 20.535	42 20.592	57 20.311
13 20.045	28 20.563	43 20.568	58 20.298
14 20.087	29 20.589	44 20.545	59 20.286
15 20.128	30 20.616	45 20.523	60 20.273

T-V2-131218-144 Clay,Potential: Air Dry #4 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

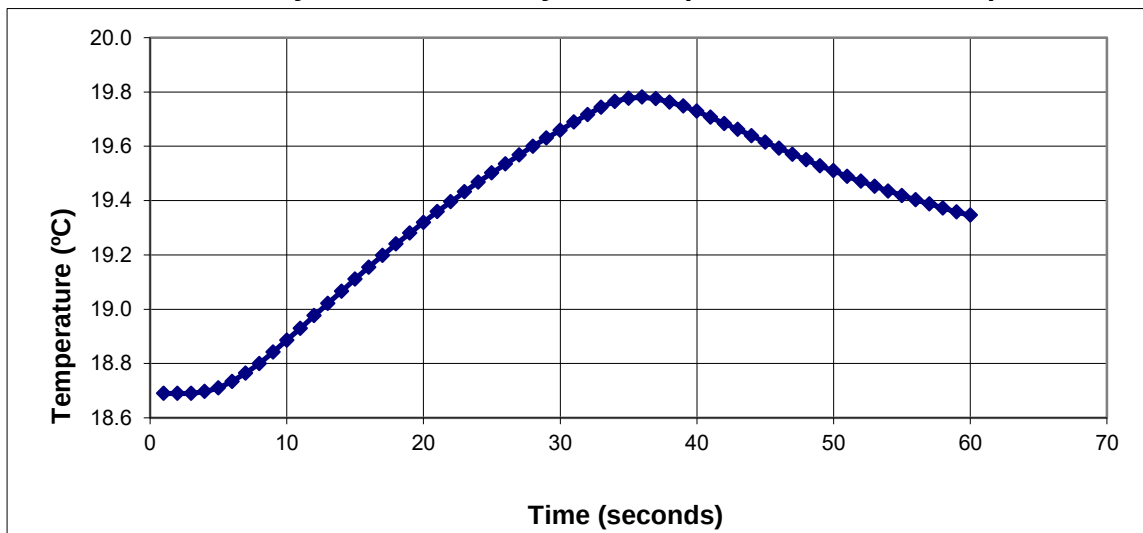
Sample Number: T-V2-131218-144 Clay
Potential (-cm water): Air Dry #5

Test Date/Time: 12/31/2018 8:37 AM	K (W/(m·K)): 0.923
Sensor: SH-1	ρ (°C·cm/W): 108.3
Test Temp.(°C): 18.7	C (MJ/(m ³ ·K)): 1.313
KD2 Pro Sample ID: CLAYAD5	D (mm ² /s): 0.703
Power (W/m): 19.820	Err: 0.0045
Current (amps): 0.138	

Raw Data

Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)
1 18.690	16 19.155	31 19.689	46 19.593
2 18.690	17 19.198	32 19.717	47 19.571
3 18.690	18 19.240	33 19.744	48 19.550
4 18.697	19 19.280	34 19.765	49 19.528
5 18.710	20 19.320	35 19.777	50 19.510
6 18.734	21 19.359	36 19.781	51 19.489
7 18.765	22 19.396	37 19.775	52 19.471
8 18.800	23 19.432	38 19.763	53 19.453
9 18.842	24 19.468	39 19.748	54 19.435
10 18.886	25 19.502	40 19.729	55 19.419
11 18.930	26 19.535	41 19.707	56 19.403
12 18.976	27 19.568	42 19.684	57 19.388
13 19.021	28 19.600	43 19.662	58 19.373
14 19.066	29 19.630	44 19.639	59 19.358
15 19.111	30 19.659	45 19.615	60 19.346

T-V2-131218-144 Clay,Potential: Air Dry #5 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

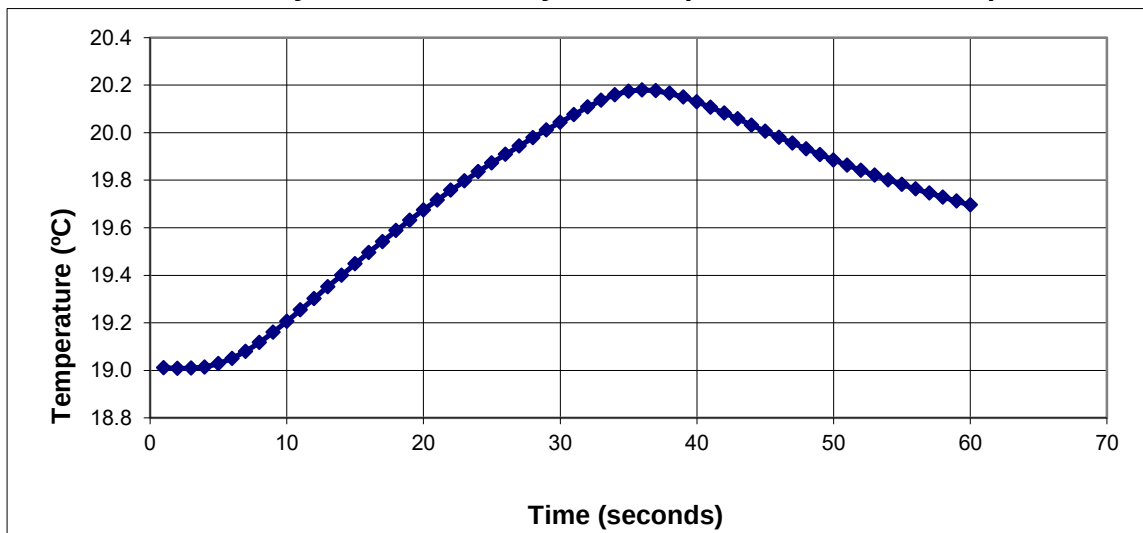
Sample Number: T-V2-131218-144 Clay
Potential (-cm water): Air Dry #6

Test Date/Time: 1/2/2019 10:04 AM	K (W/(m·K)): 0.804
Sensor: SH-1	ρ (°C·cm/W): 124.3
Test Temp.(°C): 19.0	C (MJ/(m ³ ·K)): 1.187
KD2 Pro Sample ID: CLAYAD6	D (mm ² /s): 0.678
Power (W/m): 19.610	Err: 0.0042
Current (amps): 0.137	

Raw Data

Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)
1 19.011	16 19.496	31 20.076	46 19.981
2 19.008	17 19.542	32 20.108	47 19.956
3 19.009	18 19.588	33 20.136	48 19.932
4 19.014	19 19.632	34 20.160	49 19.908
5 19.028	20 19.675	35 20.175	50 19.885
6 19.050	21 19.716	36 20.180	51 19.863
7 19.080	22 19.758	37 20.177	52 19.842
8 19.117	23 19.797	38 20.166	53 19.821
9 19.160	24 19.836	39 20.150	54 19.802
10 19.206	25 19.873	40 20.130	55 19.782
11 19.254	26 19.909	41 20.107	56 19.763
12 19.302	27 19.944	42 20.083	57 19.746
13 19.352	28 19.979	43 20.058	58 19.729
14 19.400	29 20.012	44 20.032	59 19.712
15 19.448	30 20.044	45 20.006	60 19.696

T-V2-131218-144 Clay,Potential: Air Dry #6 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

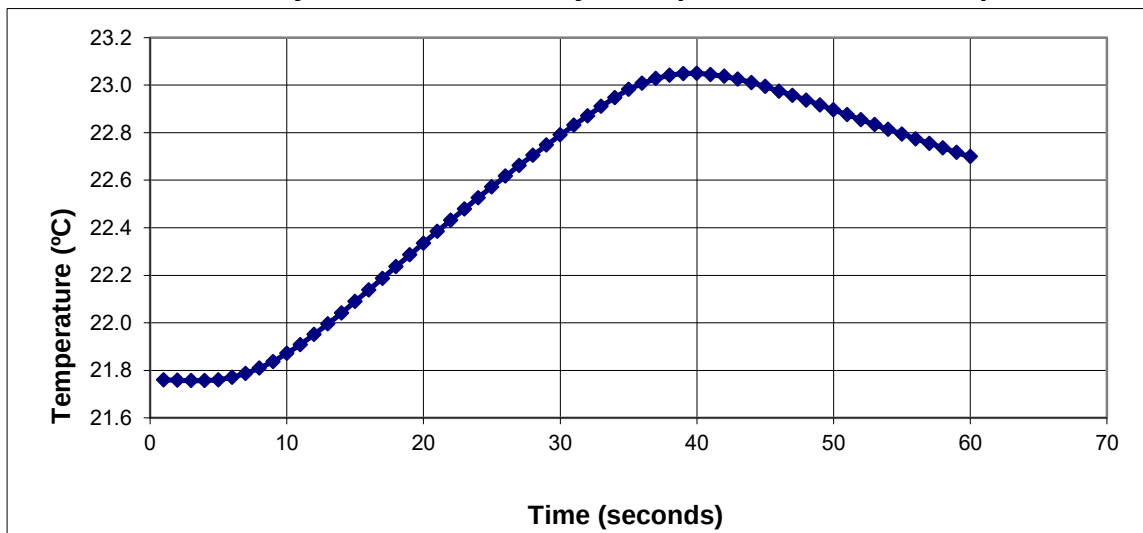
Sample Number: T-V2-131218-144 Clay
Potential (-cm water): Oven Dry

Test Date/Time:	1/7/2019 1:46 PM	K (W/(m·K)):	0.524
Sensor:	SH-1	ρ (°C·cm/W):	190.9
Test Temp.(°C):	21.8	C (MJ/(m ³ ·K)):	1.118
KD2 Pro Sample ID:	CLAYOD	D (mm ² /s):	0.468
Power (W/m):	19.390	Err:	0.0021
Current (amps):	0.137		

Raw Data

Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)	
1	21.760	16	22.138	31	22.831	46	22.975
2	21.758	17	22.187	32	22.871	47	22.957
3	21.757	18	22.237	33	22.911	48	22.937
4	21.757	19	22.286	34	22.948	49	22.917
5	21.760	20	22.335	35	22.982	50	22.896
6	21.770	21	22.384	36	23.008	51	22.876
7	21.787	22	22.432	37	23.028	52	22.855
8	21.809	23	22.479	38	23.042	53	22.834
9	21.837	24	22.526	39	23.049	54	22.814
10	21.871	25	22.572	40	23.050	55	22.794
11	21.908	26	22.617	41	23.045	56	22.774
12	21.951	27	22.662	42	23.037	57	22.755
13	21.995	28	22.705	43	23.025	58	22.736
14	22.041	29	22.748	44	23.011	59	22.717
15	22.090	30	22.791	45	22.994	60	22.699

T-V2-131218-144 Clay,Potential: Oven Dry - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Daniel B. Stephens & Associates, Inc.

Thermal Properties Results Sheet for Sample: T-V2-131218-144 Sand

Job Name: EA Engineering, Inc.
 Job Number: DB18.1333.00
 Sample Number: T-V2-131218-144 Sand
 Project Name: 62735DM02 KAFB-Bulk Fuels Fac Vadose Zone
 PO Number: 18189

Instrument Description: Decagon KD2 Pro

Probe: ☐ KS-1, 6 cm length, 1.3 mm diameter, single needle
☐ TR-1, 10 cm length, 2.4 mm diameter, single needle
☒ SH-1, 3 cm length, 1.3 mm diameter, dual needle, 6 mm spacing

Test Start Date: 12/18/18

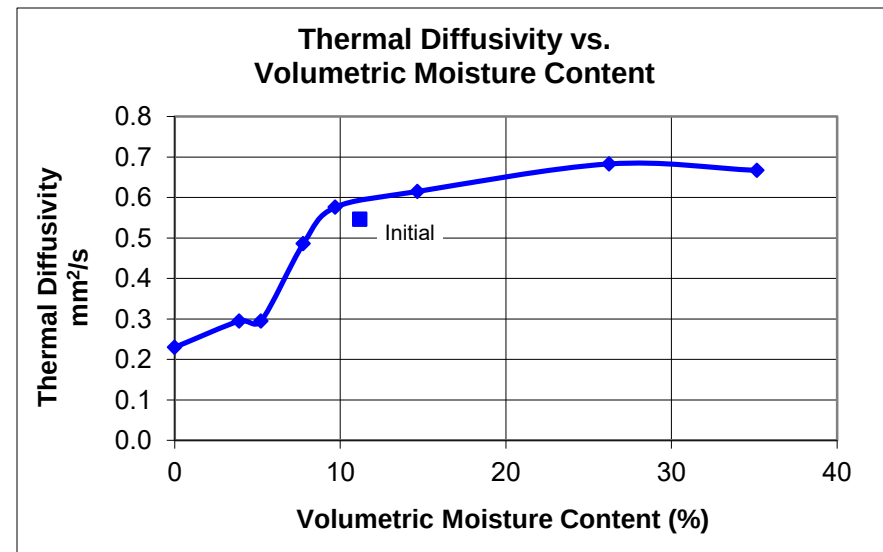
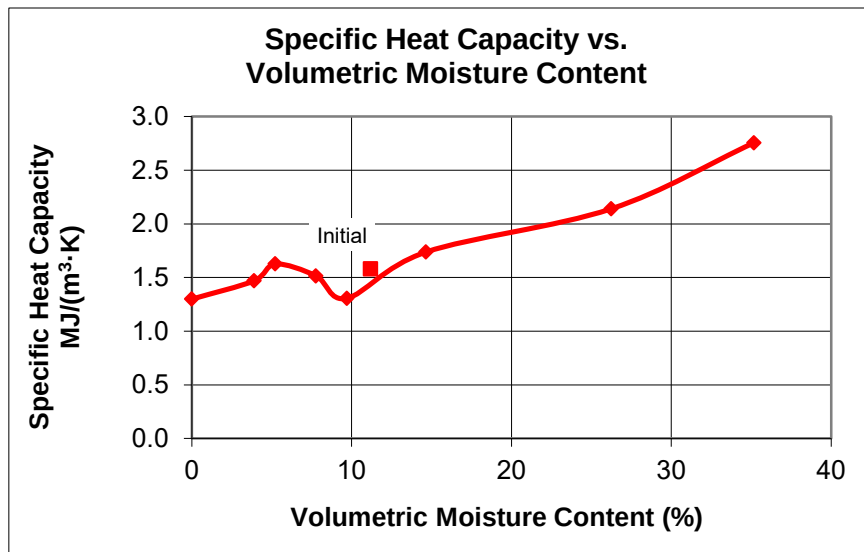
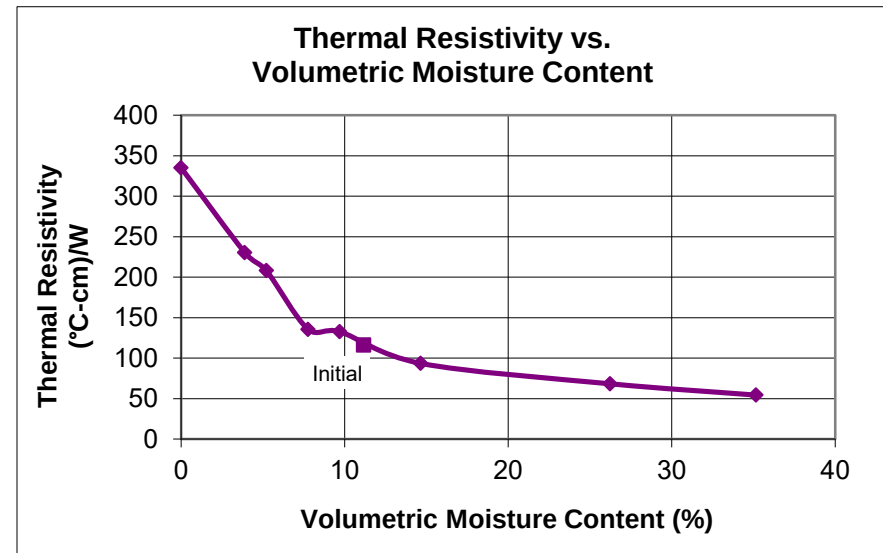
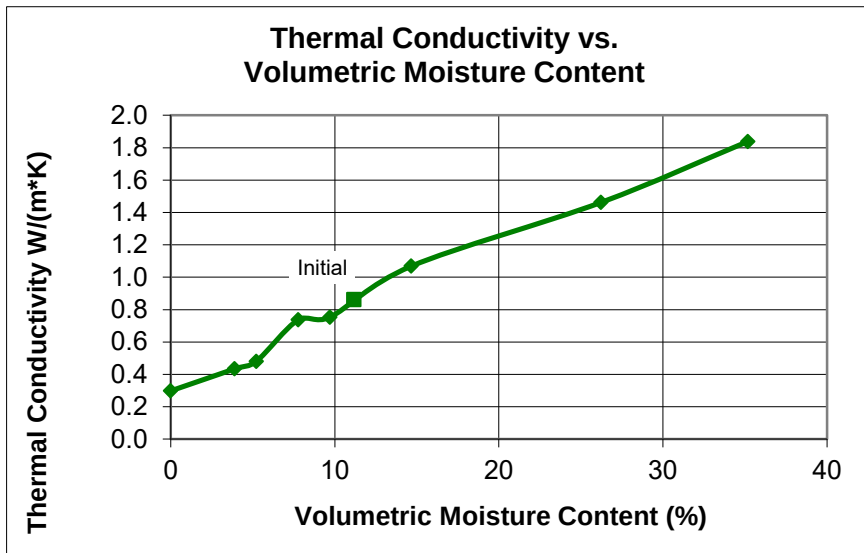
Reading	Water Potential (-cm water)	Gravimetric Moisture Content (g/g, %)	Volumetric Moisture Content ¹ (vol/vol, %)	Dry Bulk Density ¹ (g/cm ³)	Test Temperature (°C)	K Thermal Conductivity W/(m·K)	ρ Thermal Resistivity °C·cm/W	C Specific Heat Capacity MJ/(m ³ ·K)	D Thermal Diffusivity (mm ² /s)
Initial	---	6.49	11.18	1.72	20.47	0.862	116.0	1.579	0.546
Saturated	0	20.43	35.17	1.72	18.21	1.838	54.4	2.756	0.667
Air Dry #1	---	15.23	26.23	1.72	16.02	1.462	68.4	2.140	0.683
Air Dry #2	---	8.51	14.66	1.72	17.91	1.068	93.6	1.738	0.615
Air Dry #3	---	5.63	9.70	1.72	18.74	0.753	132.8	1.307	0.576
Air Dry #4	---	4.51	7.77	1.72	20.41	0.737	135.7	1.516	0.486
Air Dry #5	---	3.03	5.21	1.72	19.19	0.480	208.3	1.627	0.295
Air Dry #6	---	2.26	3.90	1.72	19.17	0.434	230.5	1.470	0.295
Oven Dry	---	0	0	1.72	21.86	0.298	335.3	1.299	0.230

--- = Value not measured.

¹ Adjusted for volume changes during testing, if applicable.



Thermal Properties Results Sheet for Sample: T-V2-131218-144 Sand Scatter Plots





Thermal Properties Data

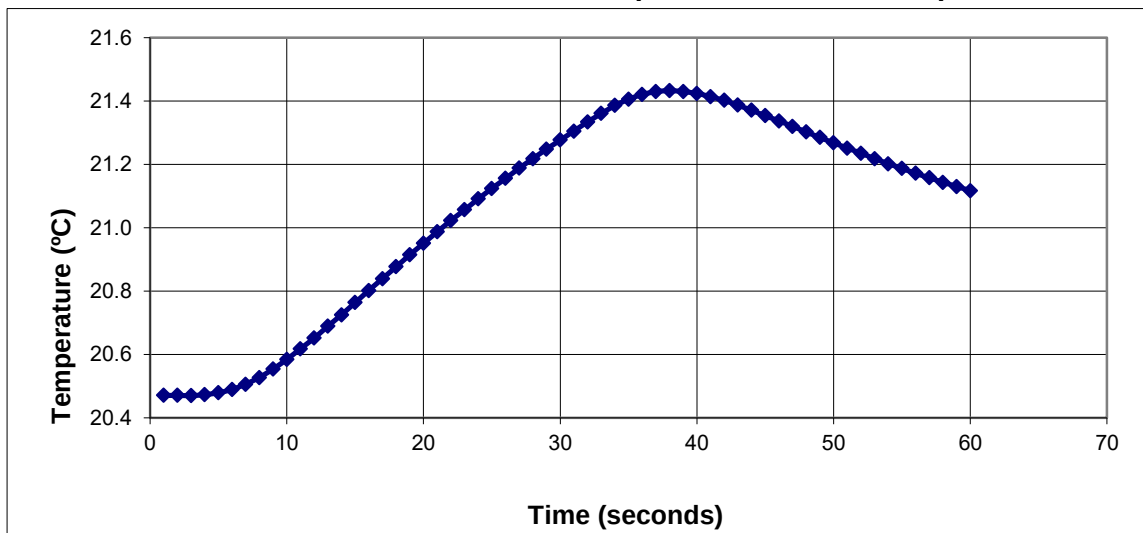
Sample Number: T-V2-131218-144 Sand
Potential (-cm water): Initial

Test Date/Time: 12/18/2018 9:12 AM	K (W/(m·K)): 0.862
Sensor: SH-1	ρ (°C·cm/W): 116.0
Test Temp.(°C): 20.5	C (MJ/(m ³ ·K)): 1.579
KD2 Pro Sample ID: SAND-IN	D (mm ² /s): 0.546
Power (W/m): 20.600	Err: 0.0030
Current (amps): 0.141	

Raw Data

Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)
1 20.471	16 20.802	31 21.305	46 21.337
2 20.471	17 20.839	32 21.334	47 21.320
3 20.470	18 20.877	33 21.361	48 21.302
4 20.473	19 20.915	34 21.386	49 21.285
5 20.479	20 20.951	35 21.406	50 21.268
6 20.489	21 20.987	36 21.421	51 21.251
7 20.506	22 21.023	37 21.430	52 21.235
8 20.527	23 21.057	38 21.433	53 21.218
9 20.554	24 21.091	39 21.430	54 21.202
10 20.584	25 21.124	40 21.424	55 21.187
11 20.618	26 21.156	41 21.414	56 21.172
12 20.652	27 21.188	42 21.402	57 21.158
13 20.689	28 21.218	43 21.387	58 21.143
14 20.725	29 21.248	44 21.371	59 21.130
15 20.764	30 21.277	45 21.354	60 21.117

T-V2-131218-144 Sand,Potential: Initial - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

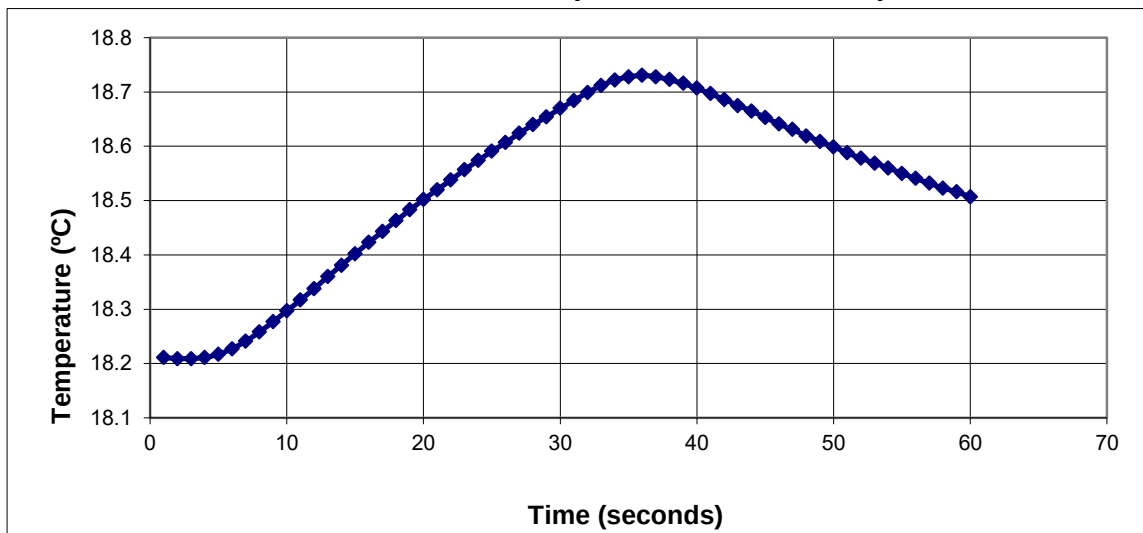
Sample Number: T-V2-131218-144 Sand
Potential (-cm water): 0

Test Date/Time: 12/19/2018 8:30 AM	K (W/(m·K)): 1.838
Sensor: SH-1	ρ (°C·cm/W): 54.4
Test Temp.(°C): 18.2	C (MJ/(m ³ ·K)): 2.756
KD2 Pro Sample ID: SAND-SA	D (mm ² /s): 0.667
Power (W/m): 20.390	Err: 0.0025
Current (amps): 0.140	

Raw Data

Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)	
1	18.211	16	18.423	31	18.684	46	18.641
2	18.209	17	18.443	32	18.699	47	18.631
3	18.209	18	18.463	33	18.712	48	18.619
4	18.211	19	18.483	34	18.722	49	18.609
5	18.217	20	18.502	35	18.728	50	18.599
6	18.227	21	18.520	36	18.731	51	18.588
7	18.241	22	18.538	37	18.728	52	18.578
8	18.258	23	18.557	38	18.723	53	18.569
9	18.277	24	18.574	39	18.716	54	18.560
10	18.297	25	18.591	40	18.707	55	18.550
11	18.317	26	18.607	41	18.697	56	18.541
12	18.338	27	18.624	42	18.686	57	18.532
13	18.360	28	18.640	43	18.675	58	18.523
14	18.381	29	18.654	44	18.665	59	18.516
15	18.402	30	18.670	45	18.653	60	18.507

T-V2-131218-144 Sand,Potential: 0 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

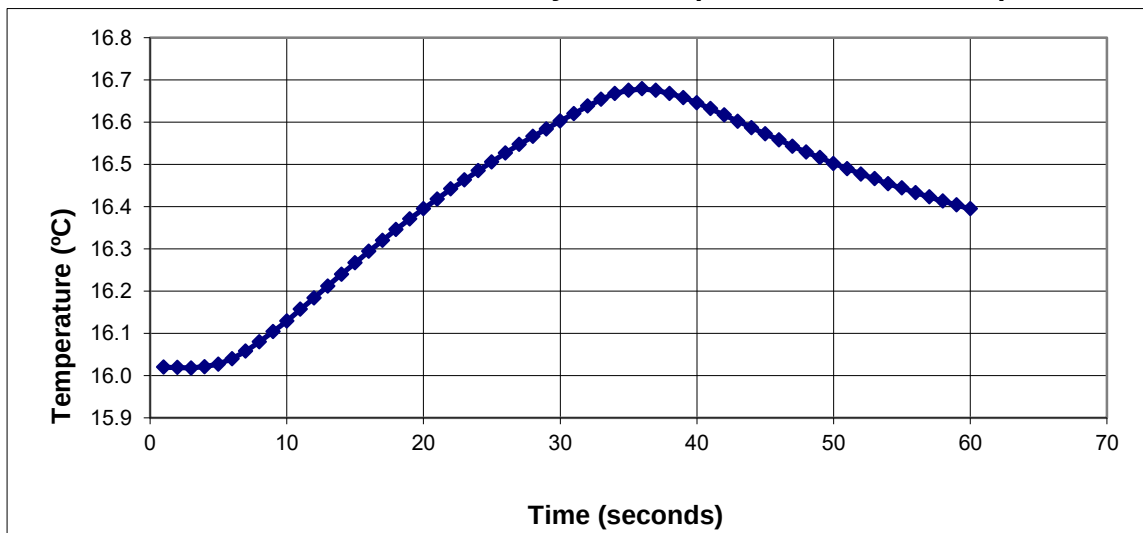
Sample Number: T-V2-131218-144 Sand
Potential (-cm water): Air Dry #1

Test Date/Time: 12/20/2018 10:39 AM	K (W/(m·K)): 1.462
Sensor: SH-1	ρ (°C·cm/W): 68.4
Test Temp.(°C): 16.0	C (MJ/(m ³ ·K)): 2.140
KD2 Pro Sample ID: SANDAD1	D (mm ² /s): 0.683
Power (W/m): 20.110	Err: 0.0047
Current (amps): 0.139	

Raw Data

Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)
1 16.020	16 16.294	31 16.620	46 16.558
2 16.019	17 16.320	32 16.638	47 16.543
3 16.018	18 16.346	33 16.654	48 16.529
4 16.021	19 16.371	34 16.668	49 16.516
5 16.027	20 16.395	35 16.675	50 16.502
6 16.040	21 16.418	36 16.679	51 16.490
7 16.058	22 16.442	37 16.675	52 16.477
8 16.080	23 16.463	38 16.668	53 16.466
9 16.104	24 16.485	39 16.658	54 16.454
10 16.129	25 16.506	40 16.646	55 16.444
11 16.157	26 16.527	41 16.632	56 16.433
12 16.184	27 16.547	42 16.617	57 16.423
13 16.212	28 16.566	43 16.602	58 16.413
14 16.240	29 16.584	44 16.587	59 16.404
15 16.267	30 16.603	45 16.572	60 16.395

T-V2-131218-144 Sand,Potential: Air Dry #1 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

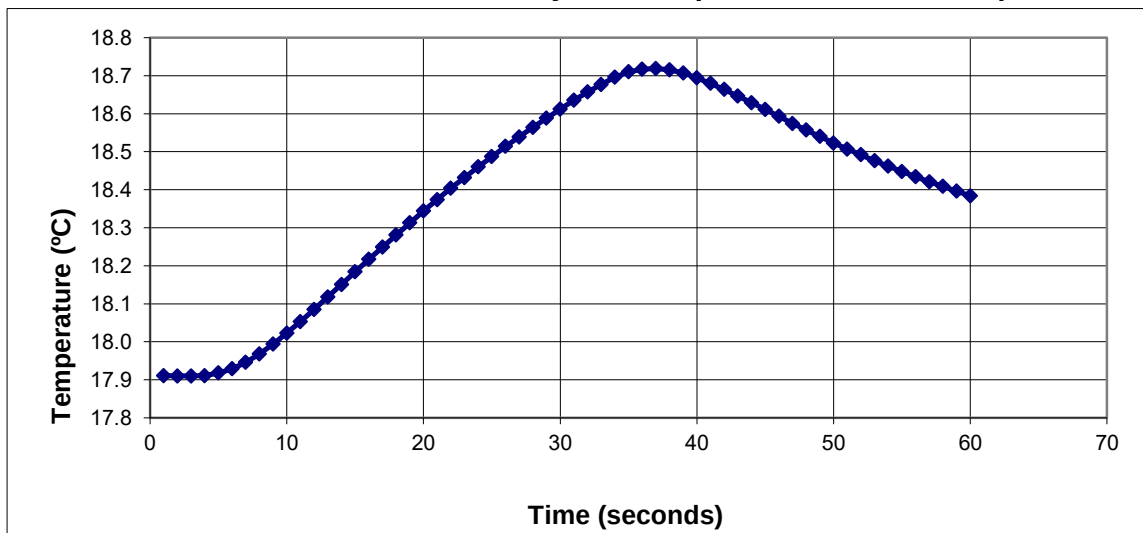
Sample Number: T-V2-131218-144 Sand
Potential (-cm water): Air Dry #2

Test Date/Time: 12/21/2018 1:33 PM	K (W/(m·K)): 1.068
Sensor: SH-1	ρ (°C·cm/W): 93.6
Test Temp.(°C): 17.9	C (MJ/(m ³ ·K)): 1.738
KD2 Pro Sample ID: SANDAD2	D (mm ² /s): 0.615
Power (W/m): 20.190	Err: 0.0039
Current (amps): 0.139	

Raw Data

Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)
1 17.911	16 18.217	31 18.635	46 18.593
2 17.910	17 18.249	32 18.657	47 18.574
3 17.910	18 18.281	33 18.677	48 18.557
4 17.911	19 18.313	34 18.696	49 18.540
5 17.918	20 18.344	35 18.710	50 18.523
6 17.929	21 18.374	36 18.717	51 18.507
7 17.946	22 18.404	37 18.719	52 18.492
8 17.968	23 18.432	38 18.715	53 18.476
9 17.994	24 18.460	39 18.707	54 18.462
10 18.023	25 18.487	40 18.694	55 18.448
11 18.053	26 18.514	41 18.680	56 18.434
12 18.085	27 18.539	42 18.664	57 18.421
13 18.118	28 18.564	43 18.646	58 18.409
14 18.151	29 18.588	44 18.629	59 18.396
15 18.184	30 18.612	45 18.611	60 18.384

T-V2-131218-144 Sand,Potential: Air Dry #2 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

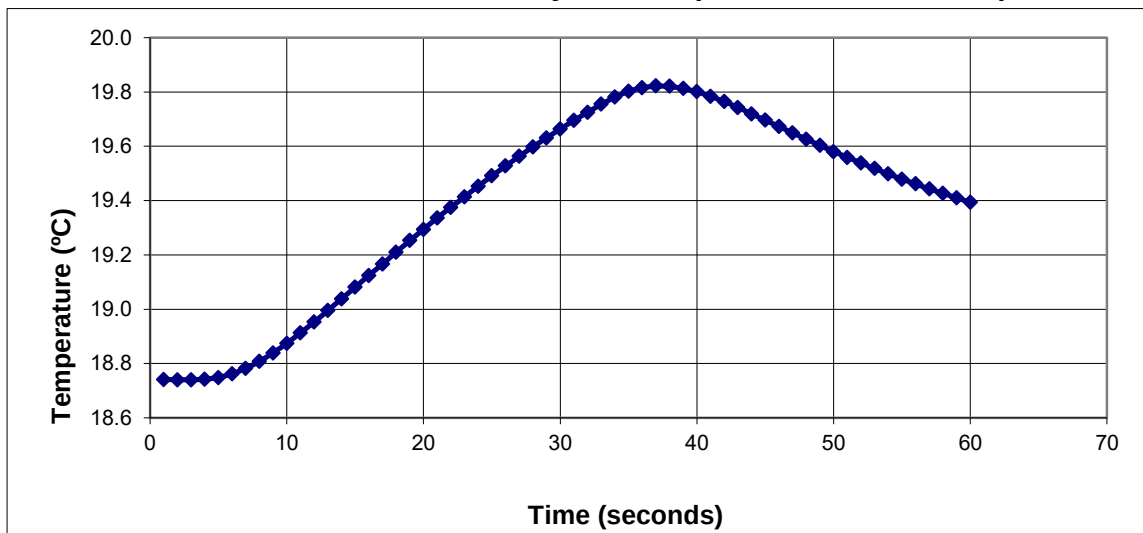
Sample Number: T-V2-131218-144 Sand
Potential (-cm water): Air Dry #3

Test Date/Time: 12/27/2018 8:13 AM	K (W/(m·K)): 0.753
Sensor: SH-1	ρ (°C·cm/W): 132.8
Test Temp.(°C): 18.7	C (MJ/(m ³ ·K)): 1.307
KD2 Pro Sample ID: SANDAD3	D (mm ² /s): 0.576
Power (W/m): 20.310	Err: 0.0037
Current (amps): 0.140	

Raw Data

Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)
1 18.741	16 19.124	31 19.695	46 19.673
2 18.740	17 19.166	32 19.725	47 19.649
3 18.740	18 19.210	33 19.755	48 19.626
4 18.742	19 19.253	34 19.781	49 19.603
5 18.748	20 19.294	35 19.803	50 19.580
6 18.762	21 19.336	36 19.816	51 19.559
7 18.782	22 19.375	37 19.822	52 19.539
8 18.808	23 19.414	38 19.821	53 19.518
9 18.839	24 19.453	39 19.813	54 19.498
10 18.874	25 19.491	40 19.801	55 19.479
11 18.913	26 19.528	41 19.784	56 19.462
12 18.953	27 19.563	42 19.765	57 19.443
13 18.995	28 19.598	43 19.743	58 19.427
14 19.038	29 19.631	44 19.719	59 19.410
15 19.081	30 19.664	45 19.696	60 19.394

T-V2-131218-144 Sand,Potential: Air Dry #3 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

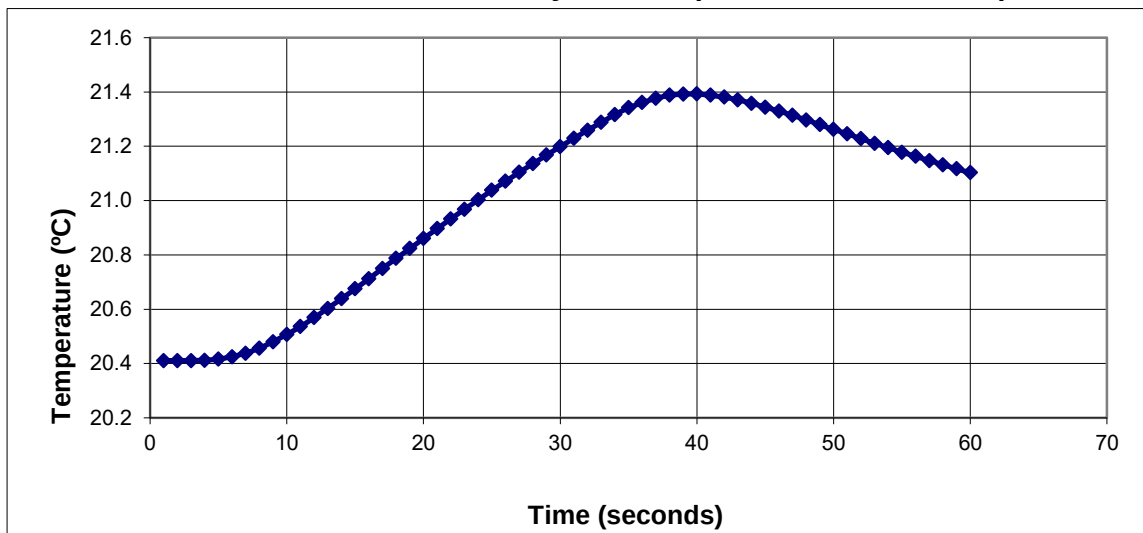
Sample Number: T-V2-131218-144 Sand
Potential (-cm water): Air Dry #4

Test Date/Time: 12/28/2018 2:44 PM	K (W/(m·K)): 0.737
Sensor: SH-1	ρ (°C·cm/W): 135.7
Test Temp.(°C): 20.4	C (MJ/(m ³ ·K)): 1.516
KD2 Pro Sample ID: SANDAD4	D (mm ² /s): 0.486
Power (W/m): 20.000	Err: 0.0038
Current (amps): 0.139	

Raw Data

Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)
1 20.411	16 20.712	31 21.229	46 21.330
2 20.411	17 20.750	32 21.259	47 21.314
3 20.411	18 20.787	33 21.288	48 21.297
4 20.412	19 20.824	34 21.317	49 21.280
5 20.416	20 20.860	35 21.342	50 21.262
6 20.425	21 20.897	36 21.361	51 21.246
7 20.438	22 20.933	37 21.377	52 21.228
8 20.456	23 20.968	38 21.388	53 21.211
9 20.480	24 21.003	39 21.392	54 21.195
10 20.507	25 21.038	40 21.393	55 21.178
11 20.536	26 21.071	41 21.388	56 21.163
12 20.569	27 21.104	42 21.381	57 21.147
13 20.603	28 21.136	43 21.371	58 21.132
14 20.639	29 21.168	44 21.358	59 21.117
15 20.675	30 21.199	45 21.344	60 21.103

T-V2-131218-144 Sand,Potential: Air Dry #4 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

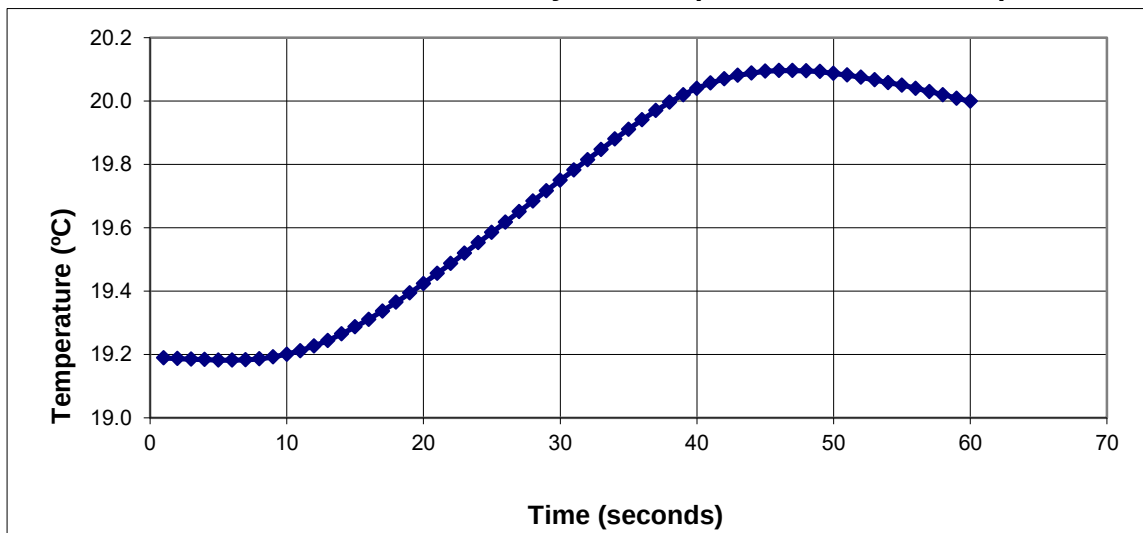
Sample Number: T-V2-131218-144 Sand
Potential (-cm water): Air Dry #5

Test Date/Time: 12/31/2018 8:34 AM	K (W/(m·K)): 0.480
Sensor: SH-1	ρ (°C·cm/W): 208.3
Test Temp.(°C): 19.2	C (MJ/(m ³ ·K)): 1.627
KD2 Pro Sample ID: SANDAD5	D (mm ² /s): 0.295
Power (W/m): 19.830	Err: 0.0016
Current (amps): 0.138	

Raw Data

Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)
1 19.189	16 19.311	31 19.782	46 20.096
2 19.187	17 19.337	32 19.815	47 20.096
3 19.185	18 19.365	33 19.847	48 20.095
4 19.184	19 19.394	34 19.880	49 20.093
5 19.182	20 19.424	35 19.911	50 20.087
6 19.182	21 19.456	36 19.941	51 20.082
7 19.183	22 19.487	37 19.970	52 20.075
8 19.186	23 19.520	38 19.996	53 20.067
9 19.192	24 19.553	39 20.020	54 20.058
10 19.201	25 19.585	40 20.040	55 20.050
11 19.212	26 19.618	41 20.057	56 20.040
12 19.227	27 19.651	42 20.070	57 20.030
13 19.244	28 19.684	43 20.081	58 20.020
14 19.265	29 19.717	44 20.088	59 20.009
15 19.287	30 19.750	45 20.094	60 19.999

T-V2-131218-144 Sand,Potential: Air Dry #5 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

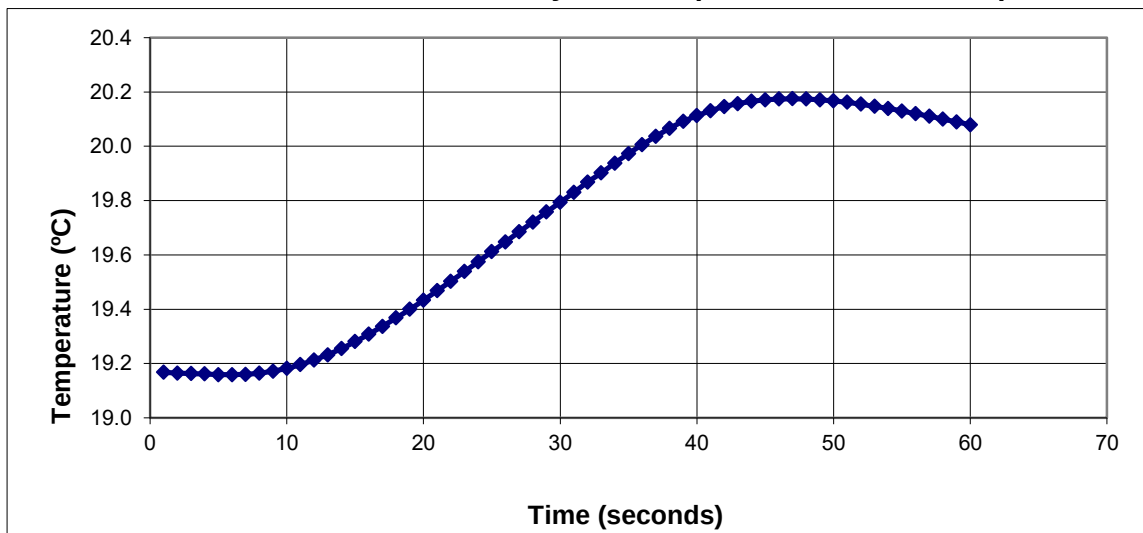
Sample Number: T-V2-131218-144 Sand
Potential (-cm water): Air Dry #6

Test Date/Time: 1/2/2019 9:58 AM	K (W/(m·K)): 0.434
Sensor: SH-1	ρ (°C·cm/W): 230.5
Test Temp.(°C): 19.2	C (MJ/(m ³ ·K)): 1.470
KD2 Pro Sample ID: SANDAD6	D (mm ² /s): 0.295
Power (W/m): 19.630	Err: 0.0009
Current (amps): 0.137	

Raw Data

Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)
1 19.168	16 19.308	31 19.830	46 20.174
2 19.165	17 19.337	32 19.868	47 20.175
3 19.163	18 19.368	33 19.902	48 20.174
4 19.162	19 19.400	34 19.938	49 20.171
5 19.159	20 19.433	35 19.973	50 20.167
6 19.159	21 19.468	36 20.006	51 20.162
7 19.160	22 19.503	37 20.037	52 20.155
8 19.164	23 19.539	38 20.066	53 20.147
9 19.172	24 19.575	39 20.092	54 20.139
10 19.182	25 19.612	40 20.113	55 20.130
11 19.196	26 19.648	41 20.131	56 20.120
12 19.213	27 19.685	42 20.146	57 20.111
13 19.232	28 19.721	43 20.157	58 20.100
14 19.255	29 19.758	44 20.166	59 20.090
15 19.281	30 19.794	45 20.171	60 20.079

T-V2-131218-144 Sand,Potential: Air Dry #6 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

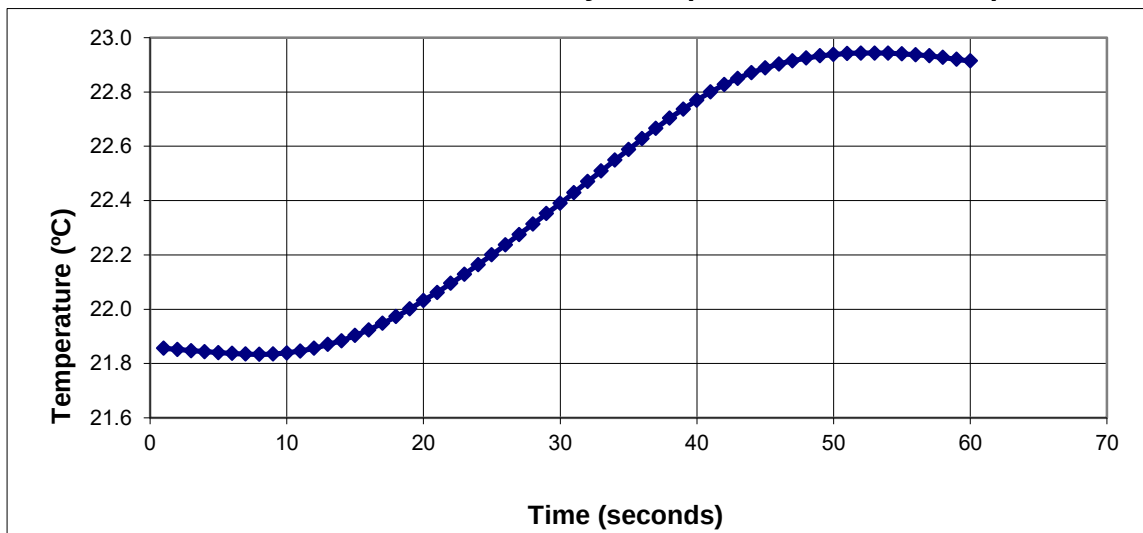
Sample Number: T-V2-131218-144 Sand
Potential (-cm water): Oven Dry

Test Date/Time: 1/7/2019 1:44 PM	K (W/(m·K)): 0.298
Sensor: SH-1	ρ (°C·cm/W): 335.3
Test Temp.(°C): 21.9	C (MJ/(m ³ ·K)): 1.299
KD2 Pro Sample ID: SANDOD	D (mm ² /s): 0.230
Power (W/m): 19.500	Err: 0.0003
Current (amps): 0.137	

Raw Data

Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)	
1	21.856	16	21.924	31	22.429	46	22.903
2	21.852	17	21.948	32	22.470	47	22.915
3	21.847	18	21.973	33	22.509	48	22.925
4	21.844	19	22.001	34	22.549	49	22.933
5	21.840	20	22.032	35	22.588	50	22.938
6	21.837	21	22.062	36	22.628	51	22.942
7	21.835	22	22.095	37	22.666	52	22.943
8	21.834	23	22.129	38	22.703	53	22.943
9	21.835	24	22.164	39	22.737	54	22.943
10	21.839	25	22.200	40	22.770	55	22.940
11	21.846	26	22.237	41	22.800	56	22.937
12	21.856	27	22.275	42	22.827	57	22.933
13	21.870	28	22.313	43	22.850	58	22.927
14	21.884	29	22.352	44	22.871	59	22.920
15	21.903	30	22.390	45	22.888	60	22.914

T-V2-131218-144 Sand,Potential: Oven Dry - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines

Laboratory Tests and Methods



Tests and Methods

Dry Bulk Density:	ASTM D7263
Moisture Content:	ASTM D7263, ASTM D2216
Calculated Porosity:	ASTM D7263
Thermal Properties:	ASTM D5334



COC1-V2-DBJ-20181213

~~CoCT-DBS-JRL~~

[illegible]

Laboratory Report for EA Engineering, Science and Technology, Inc.

KAFB-Bulk Fuels Fac Vadose Zone

Project #: 62735DM02, PO: 18189

February 21, 2019



Daniel B. Stephens & Associates, Inc.

4400 Alameda Blvd. NE, Suite C • Albuquerque, New Mexico 87113



February 21, 2019

Bernie Bockisch
EA Engineering, Science and Technology, Inc.
320 Gold Ave SW, Ste. 1300
Albuquerque, NM 87102
(505) 234-1105

Re: DBS&A Laboratory Report for the EA Engineering, Science, and Technology, Inc.
62735DM02 KAFB-Bulk Fuels Fac Vadose Zone Project

Dear Mr. Bockisch:

Enclosed is the report for the EA Engineering, Science, and Technology, Inc. 62735DM02 KAFB-Bulk Fuels Fac Vadose Zone project samples. Please review this report and provide any comments as samples will be held for a maximum of 30 days. After 30 days samples will be returned or disposed of in an appropriate manner.

All testing results were evaluated subjectively for consistency and reasonableness, and the results appear to be reasonably representative of the material tested. However, DBS&A does not assume any responsibility for interpretations or analyses based on the data enclosed, nor can we guarantee that these data are fully representative of the undisturbed materials at the field site. We recommend that careful evaluation of these laboratory results be made for your particular application.

The testing utilized to generate the enclosed report employs methods that are standard for the industry. The results do not constitute a professional opinion by DBS&A, nor can the results affect any professional or expert opinions rendered with respect thereto by DBS&A. You have acknowledged that all the testing undertaken by us, and the report provided, constitutes mere test results using standardized methods, and cannot be used to disqualify DBS&A from rendering any professional or expert opinion, having waived any claim of conflict of interest by DBS&A.

We are pleased to provide this service to EA Engineering and look forward to future laboratory testing on other projects. If you have any questions about the enclosed data, please do not hesitate to call.

Sincerely,

DANIEL B. STEPHENS & ASSOCIATES, INC.
SOIL TESTING & RESEARCH LABORATORY

Joleen Hines
Laboratory Manager

Enclosure

Daniel B. Stephens & Associates, Inc.
Soil Testing & Research Laboratory

4400 Alameda Blvd. NE, Suite C
Albuquerque, NM 87113

505-889-7752
FAX 505-889-0258

Summaries



Summary of Tests Performed

Laboratory Sample Number	Initial Soil Properties ¹			Saturated Hydraulic Conductivity ²			Moisture Characteristics ³								Particle Size ⁴			Specific Gravity ⁵		Air Perm- eability	FOC ⁶	Thermal Properties
	G	VM	VD	CH	FH	FW	HC	PP	FP	DPP	RH	EP	WHC	K _{unsat}	DS	WS	H	F	C			
T-247-300119-199	X	X																				X
FOC-247-310119-120																					X	
FOC-247-310119-164																					X	

¹ G = Gravimetric Moisture Content, VM = Volume Measurement Method, VD = Volume Displacement Method

² CH = Constant Head Rigid Wall, FH = Falling Head Rigid Wall, FW = Falling Head Rising Tail Flexible Wall

³ HC = Hanging Column, PP = Pressure Plate, FP = Filter Paper, DPP = Dew Point Potentiometer, RH = Relative Humidity Box, EP = Effective Porosity, WHC = Water Holding Capacity, K_{unsat} = Calculated Unsaturated Hydraulic Conductivity

⁴ DS = Dry Sieve, WS = Wet Sieve, H = Hydrometer

⁵ F = Fine (<4.75mm), C = Coarse (>4.75mm)

⁶ FOC = Fractional Organic Carbon testing was performed by Hall Environmental Analysis Laboratory in Albuquerque, NM.



Notes

Sample Receipt:

A total of three samples were hand-delivered on January 31, 2019. One sample was received in a full 6" x 24" plastic bag sealed with duct tape inside a cardboard box. The remaining two samples were received, each in an 8 oz. glass jar sealed with a lid inside a quart Ziploc bag. The two jar samples were received inside a cooler surrounded with ice. All samples were received in good order.

Sample Preparation and Testing Notes:

A Sub-sample was obtained by gently advancing a ring into the specimen for sample T-247-300119-199. The subsample was then subjected to thermal properties testing at the initial moisture content, the saturated moisture content, at several air drying moisture contents, and at the oven dry state.

Each thermal properties reading was obtained in the same location, whenever possible.

Samples FOC-247-310119-120 and FOC-247-310119-164 were subjected to Fractional Organic Carbon (FOC) testing. The FOC testing was performed by Hall Environmental Analysis Laboratory in Albuquerque, NM.

Porosity calculations are based on the use of an assumed specific gravity value of 2.65.

Volumetric water contents were adjusted for changes in volume, where applicable. Due to the irregularities formed on the sample surfaces during settling, volume measurements obtained after the initial reading should be considered estimates.



Summary of Sample Preparation/Volume Changes

Sample Number	Initial Sample Data ¹		Volume Change Post Saturation ²			Volume Change Post Oven Drying ³		
	Moisture Content (%, g/g)	Dry Bulk Density (g/cm ³)	Dry Bulk Density (g/cm ³)	% Volume Change (%)	% of Initial Density (%)	Dry Bulk Density (g/cm ³)	% Volume Change (%)	% of Initial Density (%)
T-247-300119-199	16.4	1.48	1.48	---	100.0%	1.52	-2.9%	103.0%

¹Initial Sample Data: The 'as received' dry bulk density and moisture content.

²Volume Change Post Saturation: Volume change measurements were obtained after saturation.

³Volume Change Post Oven Drying: Volume change measurements were obtained after oven drying.

Notes:

"+" indicates sample swelling, "-" indicates sample settling, and "---" indicates no volume change occurred.



**Summary of Initial Moisture Content, Dry Bulk Density
Wet Bulk Density and Calculated Porosity**

Sample Number	Moisture Content				Dry Bulk Density (g/cm ³)	Wet Bulk Density (g/cm ³)	Calculated Porosity (%)
	As Received		Remolded				
	Gravimetric	Volumetric	Gravimetric	Volumetric			
	(%, g/g)	(%, cm ³ /cm ³)	(%, g/g)	(%, cm ³ /cm ³)			
T-247-300119-199	16.4	24.3	---	---	1.48	1.72	44.1

NA = Not analyzed

--- = This sample was not remolded



Summary of Thermal Properties

Sample	Reading	Gravimetric Moisture Content (g/g, %)	Volumetric Moisture Content ¹ (vol/vol, %)	Dry Bulk Density ¹ (g/cm ³)	Temp °C	K W/(m·K)	ρ °C·cm/W	C MJ/(m ³ ·K)	D mm ² /s
T-247-300119-199	Initial	16.41	24.29	1.48	19.72	1.172	85.4	2.095	0.559
T-247-300119-199	0	30.66	45.39	1.48	18.35	1.605	62.3	2.910	0.552
T-247-300119-199	Air Dry #1	26.93	39.88	1.48	20.32	1.567	63.8	2.732	0.573
T-247-300119-199	Air Dry #2	24.28	36.71	1.51	19.96	1.351	74.0	2.373	0.569
T-247-300119-199	Air Dry #3	18.57	28.32	1.52	15.75	1.122	89.1	2.162	0.519
T-247-300119-199	Air Dry #4	13.13	20.02	1.52	17.53	0.846	118.2	2.021	0.419
T-247-300119-199	Air Dry #5	6.58	10.03	1.52	18.63	0.766	130.5	1.804	0.425
T-247-300119-199	Oven Dry	0	0	1.52	20.48	0.358	279.2	1.312	0.273

¹ Adjusted for volume changes during testing, if applicable.



Summary of Fractional Organic Carbon Tests

Sample Number	Fractional Organic Carbon (%)
FOC-247-310119-120	ND
FOC-247-310119-164	ND

ND= not detected at the reporting limit of 0.1% C

Analysis provided by Hall Environmental, Albuquerque, NM.

Initial Properties



**Summary of Initial Moisture Content, Dry Bulk Density
Wet Bulk Density and Calculated Porosity**

Sample Number	Moisture Content				Dry Bulk Density (g/cm ³)	Wet Bulk Density (g/cm ³)	Calculated Porosity (%)
	As Received		Remolded				
	Gravimetric	Volumetric	Gravimetric	Volumetric			
	(%, g/g)	(%, cm ³ /cm ³)	(%, g/g)	(%, cm ³ /cm ³)			
T-247-300119-199	16.4	24.3	---	---	1.48	1.72	44.1

NA = Not analyzed

--- = This sample was not remolded



**Data for Initial Moisture Content,
Bulk Density, Porosity, and Percent Saturation**

Job Name: EA Engineering, Science, and Technology, Inc.
Job Number: DB18.1333.00
Sample Number: T-247-300119-199
Project Name: 62735DM02 KAFB-Bulk Fuels Fac Vadose Zone
PO Number: 18189

	<u>As Received</u>	<u>Remolded</u>
Test Date:	1-Feb-19	---
Field weight* of sample (g):	522.23	
Tare weight, ring (g):	138.36	
Tare weight, pan/plate (g):	0.00	
Tare weight, other (g):	0.00	
Dry weight of sample (g):	329.76	
Sample volume (cm ³):	222.72	
Assumed particle density (g/cm ³):	2.65	
Gravimetric Moisture Content (% g/g):	16.4	
Volumetric Moisture Content (% vol):	24.3	
Dry bulk density (g/cm ³):	1.48	
Wet bulk density (g/cm ³):	1.72	
Calculated Porosity (% vol):	44.1	
Percent Saturation:	55.1	

Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines

Comments:

* Weight including tares
NA = Not analyzed
--- = This sample was not remolded

Thermal Properties



Summary of Thermal Properties

Sample	Reading	Gravimetric Moisture Content (g/g, %)	Volumetric Moisture Content ¹ (vol/vol, %)	Dry Bulk Density ¹ (g/cm ³)	Temp °C	K W/(m·K)	ρ °C·cm/W	C MJ/(m ³ ·K)	D mm ² /s
T-247-300119-199	Initial	16.41	24.29	1.48	19.72	1.172	85.4	2.095	0.559
T-247-300119-199	0	30.66	45.39	1.48	18.35	1.605	62.3	2.910	0.552
T-247-300119-199	Air Dry #1	26.93	39.88	1.48	20.32	1.567	63.8	2.732	0.573
T-247-300119-199	Air Dry #2	24.28	36.71	1.51	19.96	1.351	74.0	2.373	0.569
T-247-300119-199	Air Dry #3	18.57	28.32	1.52	15.75	1.122	89.1	2.162	0.519
T-247-300119-199	Air Dry #4	13.13	20.02	1.52	17.53	0.846	118.2	2.021	0.419
T-247-300119-199	Air Dry #5	6.58	10.03	1.52	18.63	0.766	130.5	1.804	0.425
T-247-300119-199	Oven Dry	0	0	1.52	20.48	0.358	279.2	1.312	0.273

¹ Adjusted for volume changes during testing, if applicable.



Daniel B. Stephens & Associates, Inc.

Thermal Properties Results Sheet for Sample: T-247-300119-199

Job Name: EA Engineering, Inc.

Job Number: DB18.1333.00

Sample Number: T-247-300119-199

Project Name: 62735DM02 KAFB-Bulk Fuels Fac Vadose Zone

PO Number: 18189

Instrument Description: Decagon KD2 Pro

Probe: ☐ KS-1, 6 cm length, 1.3 mm diameter, single needle

☐ TR-1, 10 cm length, 2.4 mm diameter, single needle

☒ SH-1, 3 cm length, 1.3 mm diameter, dual needle, 6 mm spacing

Test Start Date: 2/1/19

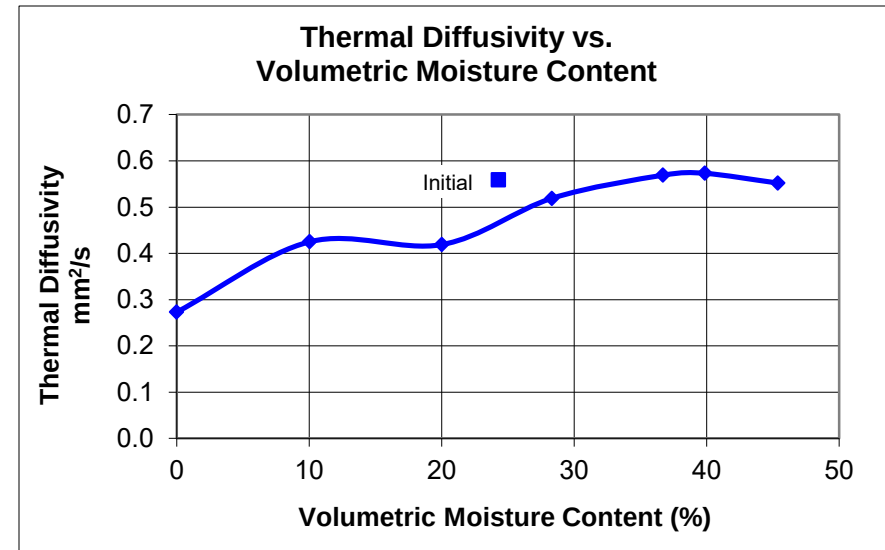
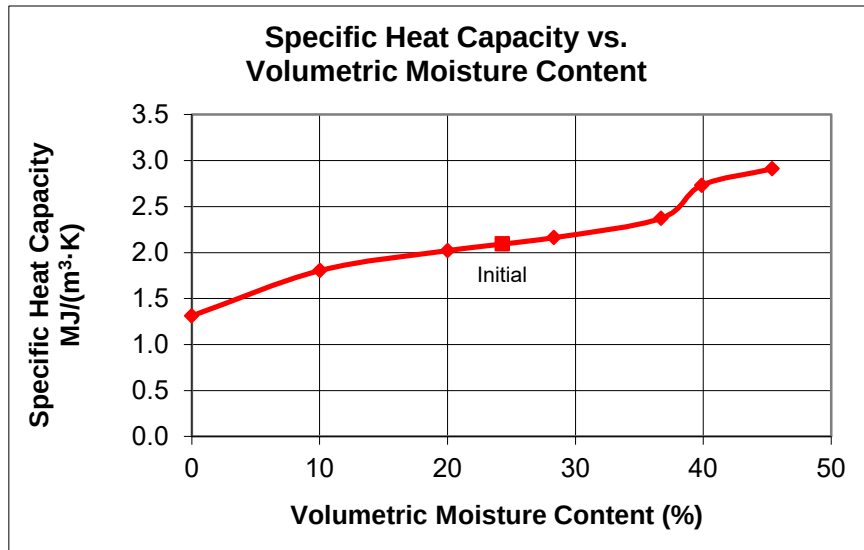
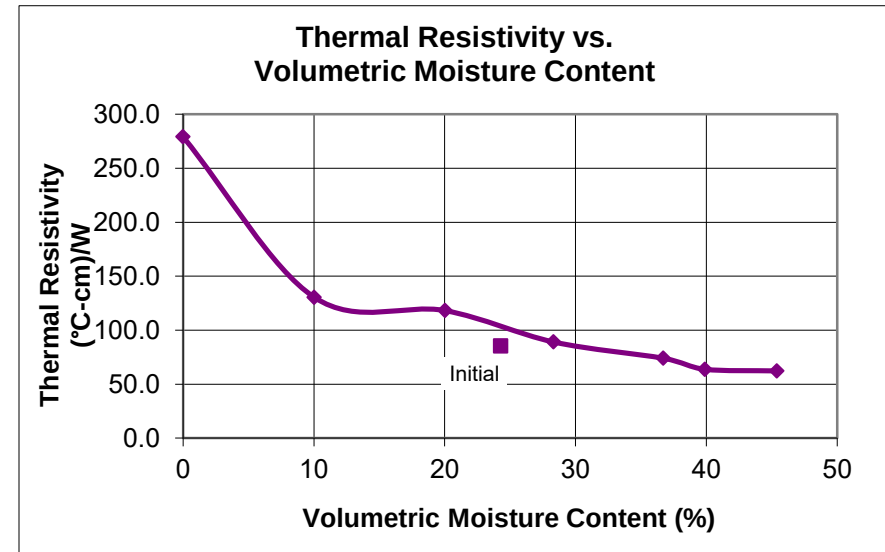
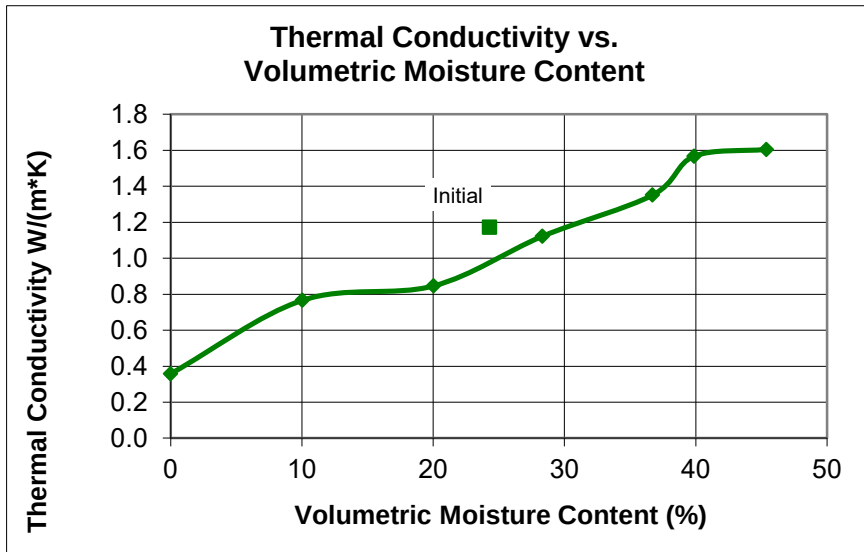
Reading	Water Potential (-cm water)	Gravimetric Moisture Content (g/g, %)	Volumetric Moisture Content ¹ (vol/vol, %)	Dry Bulk Density ¹ (g/cm ³)	Test Temperature (°C)	K Thermal Conductivity W/(m·K)	ρ Thermal Resistivity °C·cm/W	C Specific Heat Capacity MJ/(m ³ ·K)	D Thermal Diffusivity (mm ² /s)
Initial	---	16.41	24.29	1.48	19.72	1.172	85.4	2.095	0.559
Saturated	0	30.66	45.39	1.48	18.35	1.605	62.3	2.910	0.552
Air Dry #1	---	26.93	39.88	1.48	20.32	1.567	63.8	2.732	0.573
Air Dry #2	---	24.28	36.71	1.51	19.96	1.351	74.0	2.373	0.569
Air Dry #3	---	18.57	28.32	1.52	15.75	1.122	89.1	2.162	0.519
Air Dry #4	---	13.13	20.02	1.52	17.53	0.846	118.2	2.021	0.419
Air Dry #5	---	6.58	10.03	1.52	18.63	0.766	130.5	1.804	0.425
Oven Dry	---	0	0	1.52	20.48	0.358	279.2	1.312	0.273

--- = Value not measured.

¹ Adjusted for volume changes during testing, if applicable.



Thermal Properties Results Sheet for Sample: T-247-300119-199
Scatter Plots





Thermal Properties Data

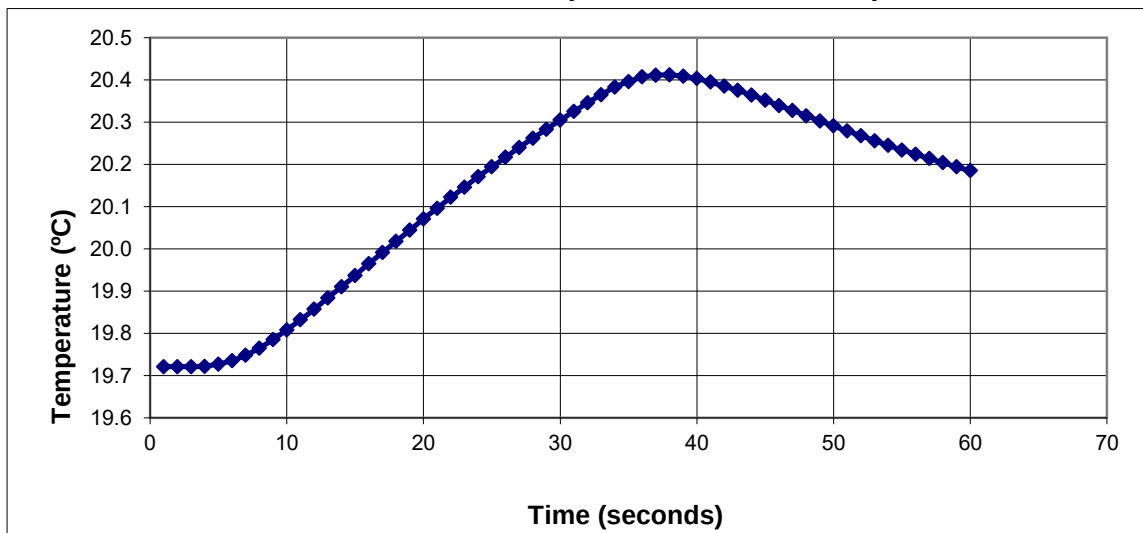
Sample Number: T-247-300119-199
Potential (-cm water): Initial

Test Date/Time: 2/1/2019 12:44 PM	K (W/(m·K)): 1.172
Sensor: SH-1	ρ (°C·cm/W): 85.4
Test Temp.(°C): 19.7	C (MJ/(m ³ ·K)): 2.095
KD2 Pro Sample ID: 199-AR	D (mm ² /s): 0.559
Power (W/m): 19.490	Err: 0.0027
Current (amps): 0.137	

Raw Data

Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)	
1	19.721	16	19.965	31	20.325	46	20.339
2	19.721	17	19.991	32	20.346	47	20.328
3	19.721	18	20.018	33	20.365	48	20.315
4	19.722	19	20.044	34	20.383	49	20.303
5	19.727	20	20.071	35	20.396	50	20.291
6	19.735	21	20.096	36	20.407	51	20.279
7	19.748	22	20.122	37	20.411	52	20.268
8	19.765	23	20.146	38	20.412	53	20.256
9	19.785	24	20.171	39	20.409	54	20.245
10	19.808	25	20.194	40	20.403	55	20.234
11	19.832	26	20.217	41	20.395	56	20.224
12	19.857	27	20.240	42	20.385	57	20.214
13	19.884	28	20.262	43	20.375	58	20.204
14	19.910	29	20.283	44	20.364	59	20.194
15	19.937	30	20.305	45	20.352	60	20.185

T-247-300119-199,Potential: Initial - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

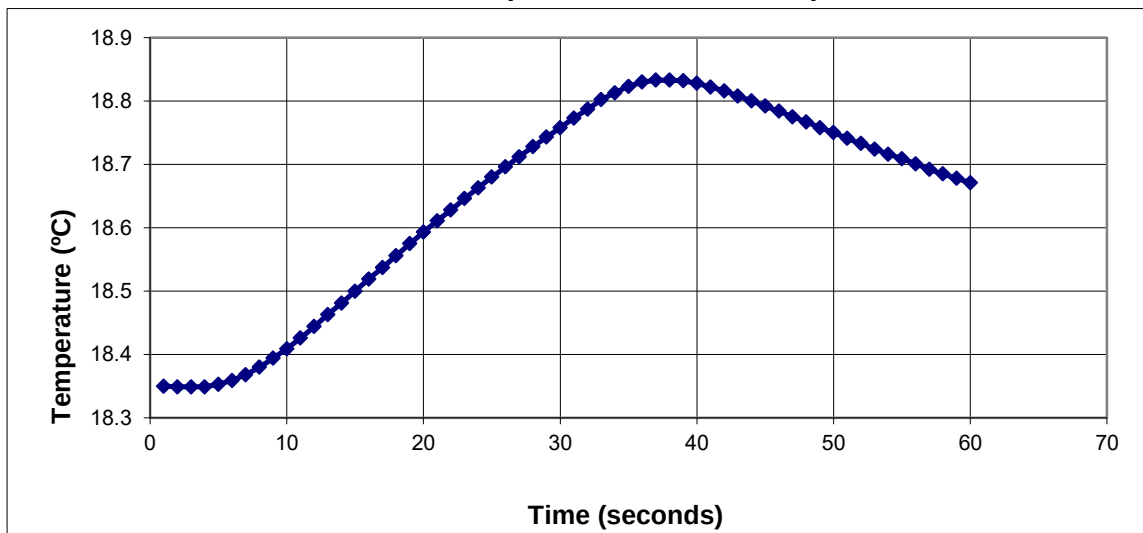
Sample Number: T-247-300119-199
Potential (-cm water): 0

Test Date/Time: 2/4/2019 8:55 AM	K (W/(m·K)): 1.605
Sensor: SH-1	ρ (°C·cm/W): 62.3
Test Temp.(°C): 18.4	C (MJ/(m ³ ·K)): 2.910
KD2 Pro Sample ID: 199-SA	D (mm ² /s): 0.552
Power (W/m): 19.130	Err: 0.0020
Current (amps): 0.136	

Raw Data

Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)	
1	18.350	16	18.519	31	18.773	46	18.784
2	18.349	17	18.537	32	18.787	47	18.775
3	18.349	18	18.556	33	18.802	48	18.767
4	18.349	19	18.575	34	18.813	49	18.758
5	18.353	20	18.593	35	18.823	50	18.750
6	18.359	21	18.611	36	18.830	51	18.741
7	18.368	22	18.628	37	18.833	52	18.733
8	18.380	23	18.646	38	18.833	53	18.724
9	18.394	24	18.663	39	18.832	54	18.716
10	18.409	25	18.680	40	18.828	55	18.709
11	18.426	26	18.696	41	18.822	56	18.701
12	18.444	27	18.712	42	18.816	57	18.692
13	18.463	28	18.728	43	18.808	58	18.685
14	18.481	29	18.743	44	18.800	59	18.678
15	18.500	30	18.758	45	18.792	60	18.671

T-247-300119-199,Potential: 0 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

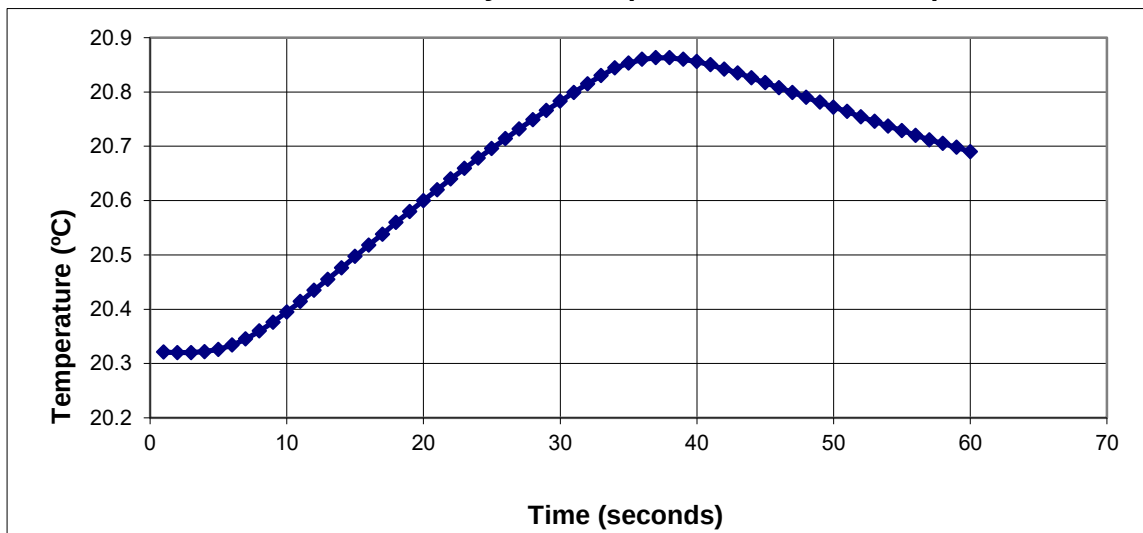
Sample Number: T-247-300119-199
Potential (-cm water): Air Dry #1

Test Date/Time:	2/5/2019 8:12 AM	K (W/(m·K)):	1.567
Sensor:	SH-1	ρ (°C·cm/W):	63.8
Test Temp.(°C):	20.3	C (MJ/(m ³ ·K)):	2.732
KD2 Pro Sample ID:	199-AD1	D (mm ² /s):	0.573
Power (W/m):	19.600	Err:	0.0020
Current (amps):	0.137		

Raw Data

Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)	
1	20.321	16	20.518	31	20.799	46	20.808
2	20.320	17	20.538	32	20.815	47	20.799
3	20.320	18	20.560	33	20.830	48	20.790
4	20.322	19	20.580	34	20.844	49	20.781
5	20.326	20	20.600	35	20.853	50	20.772
6	20.334	21	20.620	36	20.860	51	20.764
7	20.345	22	20.640	37	20.863	52	20.754
8	20.360	23	20.659	38	20.863	53	20.746
9	20.376	24	20.678	39	20.860	54	20.737
10	20.395	25	20.696	40	20.856	55	20.729
11	20.414	26	20.714	41	20.850	56	20.720
12	20.435	27	20.732	42	20.842	57	20.712
13	20.455	28	20.749	43	20.835	58	20.705
14	20.476	29	20.766	44	20.826	59	20.698
15	20.497	30	20.783	45	20.817	60	20.690

T-247-300119-199,Potential: Air Dry #1 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

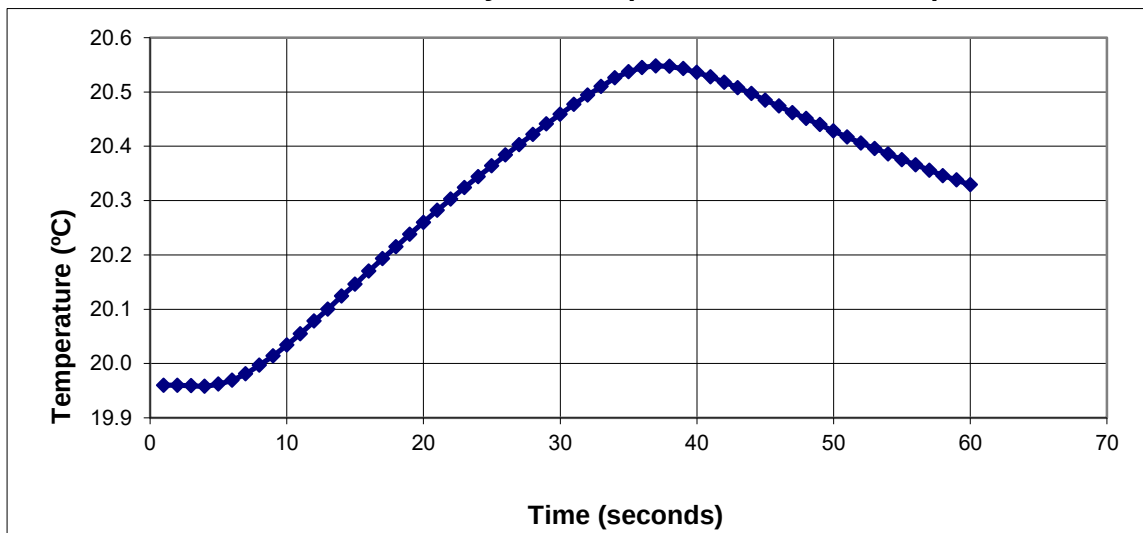
Sample Number: T-247-300119-199
Potential (-cm water): Air Dry #2

Test Date/Time:	2/6/2019 7:45 AM	K (W/(m·K)):	1.351
Sensor:	SH-1	ρ (°C·cm/W):	74.0
Test Temp.(°C):	20.0	C (MJ/(m ³ ·K)):	2.373
KD2 Pro Sample ID:	199-AD2	D (mm ² /s):	0.569
Power (W/m):	19.490	Err:	0.0026
Current (amps):	0.137		

Raw Data

Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)	
1	19.960	16	20.170	31	20.477	46	20.474
2	19.960	17	20.193	32	20.494	47	20.462
3	19.959	18	20.215	33	20.510	48	20.451
4	19.958	19	20.238	34	20.526	49	20.440
5	19.962	20	20.260	35	20.537	50	20.428
6	19.969	21	20.282	36	20.545	51	20.417
7	19.981	22	20.303	37	20.548	52	20.406
8	19.997	23	20.324	38	20.547	53	20.396
9	20.014	24	20.344	39	20.543	54	20.386
10	20.034	25	20.364	40	20.536	55	20.375
11	20.055	26	20.384	41	20.528	56	20.366
12	20.078	27	20.403	42	20.518	57	20.356
13	20.100	28	20.422	43	20.508	58	20.346
14	20.124	29	20.441	44	20.497	59	20.338
15	20.146	30	20.459	45	20.485	60	20.329

T-247-300119-199,Potential: Air Dry #2 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

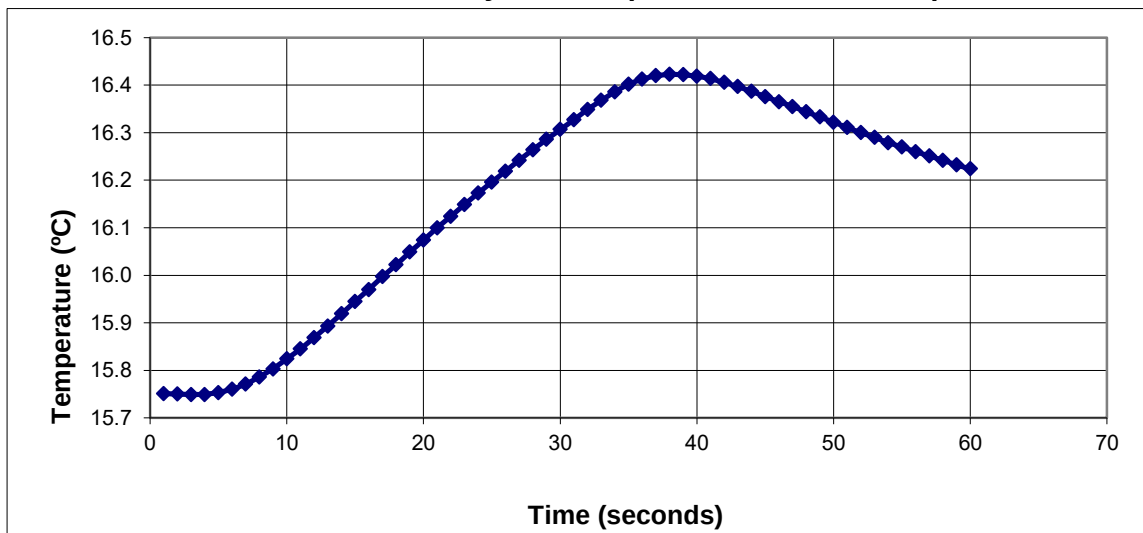
Sample Number: T-247-300119-199
Potential (-cm water): Air Dry #3

Test Date/Time: 2/7/2019 7:52 AM	K (W/(m·K)): 1.122
Sensor: SH-1	ρ (°C·cm/W): 89.1
Test Temp.(°C): 15.8	C (MJ/(m ³ ·K)): 2.162
KD2 Pro Sample ID: 199-AD3	D (mm ² /s): 0.519
Power (W/m): 19.380	Err: 0.0022
Current (amps): 0.137	

Raw Data

Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)	
1	15.751	16	15.970	31	16.327	46	16.365
2	15.750	17	15.997	32	16.349	47	16.355
3	15.749	18	16.022	33	16.368	48	16.344
4	15.749	19	16.049	34	16.386	49	16.333
5	15.753	20	16.074	35	16.402	50	16.322
6	15.760	21	16.100	36	16.413	51	16.311
7	15.771	22	16.124	37	16.420	52	16.300
8	15.786	23	16.149	38	16.423	53	16.290
9	15.803	24	16.173	39	16.422	54	16.279
10	15.824	25	16.196	40	16.419	55	16.270
11	15.845	26	16.219	41	16.414	56	16.260
12	15.869	27	16.242	42	16.406	57	16.251
13	15.893	28	16.264	43	16.397	58	16.242
14	15.919	29	16.286	44	16.387	59	16.232
15	15.945	30	16.307	45	16.376	60	16.224

T-247-300119-199,Potential: Air Dry #3 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

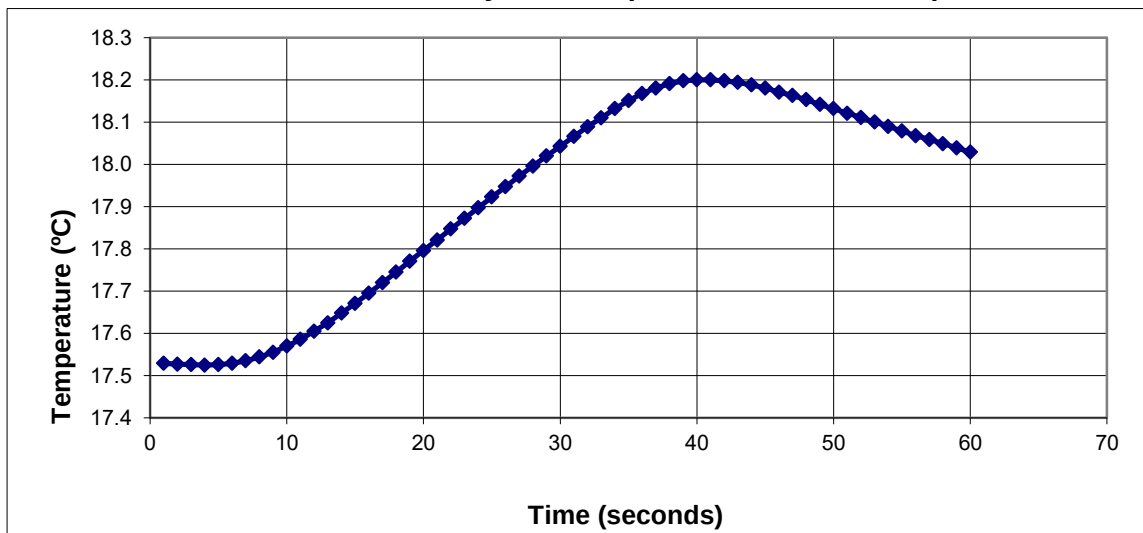
Sample Number: T-247-300119-199
Potential (-cm water): Air Dry #4

Test Date/Time:	2/8/2019 8:37 AM	K (W/(m·K)):	0.846
Sensor:	SH-1	ρ (°C·cm/W):	118.2
Test Temp.(°C):	17.5	C (MJ/(m ³ ·K)):	2.021
KD2 Pro Sample ID:	199-AD4	D (mm ² /s):	0.419
Power (W/m):	18.830	Err:	0.0017
Current (amps):	0.135		

Raw Data

Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)	
1	17.529	16	17.695	31	18.066	46	18.171
2	17.527	17	17.720	32	18.089	47	18.163
3	17.526	18	17.745	33	18.110	48	18.153
4	17.525	19	17.771	34	18.132	49	18.142
5	17.526	20	17.796	35	18.151	50	18.132
6	17.529	21	17.821	36	18.168	51	18.121
7	17.535	22	17.847	37	18.181	52	18.111
8	17.544	23	17.872	38	18.191	53	18.100
9	17.555	24	17.897	39	18.198	54	18.090
10	17.570	25	17.923	40	18.200	55	18.079
11	17.586	26	17.947	41	18.200	56	18.068
12	17.605	27	17.972	42	18.198	57	18.059
13	17.625	28	17.996	43	18.194	58	18.049
14	17.648	29	18.020	44	18.188	59	18.039
15	17.671	30	18.043	45	18.181	60	18.029

T-247-300119-199,Potential: Air Dry #4 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

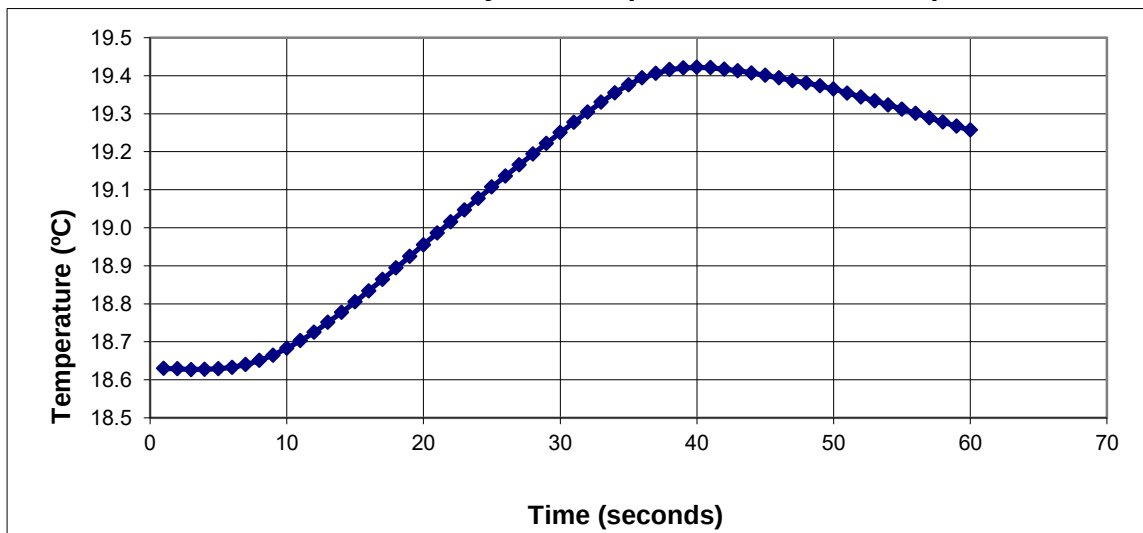
Sample Number: T-247-300119-199
Potential (-cm water): Air Dry #5

Test Date/Time: 2/11/2019 8:08 AM	K (W/(m·K)): 0.766
Sensor: SH-1	ρ (°C·cm/W): 130.5
Test Temp.(°C): 18.6	C (MJ/(m ³ ·K)): 1.804
KD2 Pro Sample ID: 199-AD5	D (mm ² /s): 0.425
Power (W/m): 18.830	Err: 0.0029
Current (amps): 0.135	

Raw Data

Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)
1 18.630	16 18.834	31 19.277	46 19.394
2 18.629	17 18.864	32 19.304	47 19.387
3 18.627	18 18.894	33 19.330	48 19.381
4 18.628	19 18.925	34 19.355	49 19.373
5 18.629	20 18.955	35 19.376	50 19.365
6 18.633	21 18.986	36 19.394	51 19.354
7 18.640	22 19.016	37 19.406	52 19.344
8 18.651	23 19.047	38 19.416	53 19.334
9 18.665	24 19.077	39 19.420	54 19.323
10 18.683	25 19.107	40 19.422	55 19.312
11 18.703	26 19.136	41 19.421	56 19.301
12 18.725	27 19.165	42 19.417	57 19.289
13 18.751	28 19.194	43 19.413	58 19.278
14 18.777	29 19.222	44 19.407	59 19.267
15 18.805	30 19.250	45 19.401	60 19.257

T-247-300119-199,Potential: Air Dry #5 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

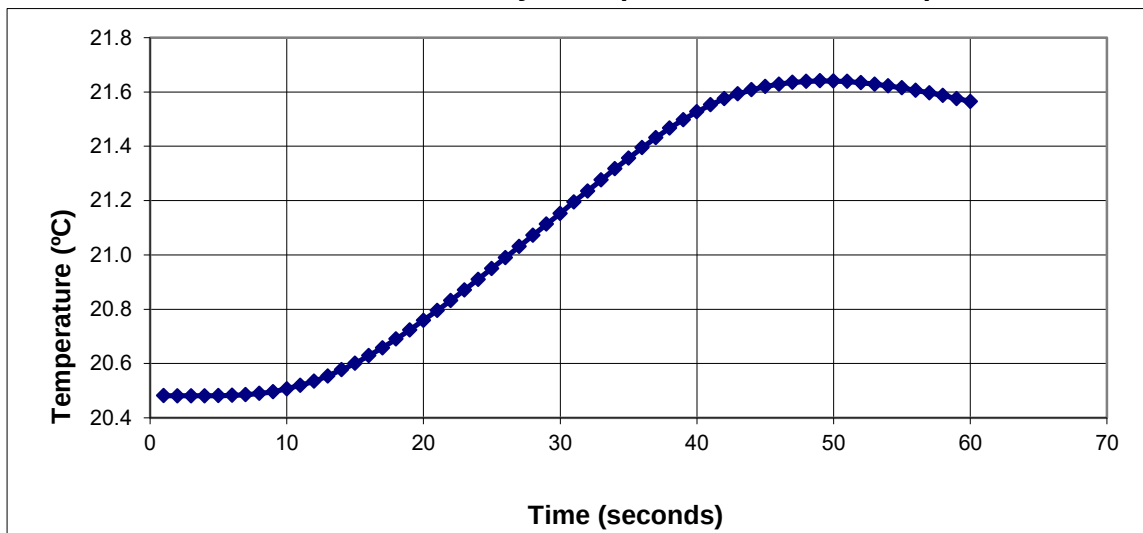
Sample Number: T-247-300119-199
Potential (-cm water): Oven Dry

Test Date/Time: 2/13/2019 1:34 PM	K (W/(m·K)): 0.358
Sensor: SH-1	ρ (°C·cm/W): 279.2
Test Temp.(°C): 20.5	C (MJ/(m ³ ·K)): 1.312
KD2 Pro Sample ID: 199-OD	D (mm ² /s): 0.273
Power (W/m): 18.870	Err: 0.0021
Current (amps): 0.135	

Raw Data

Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)	
1	20.482	16	20.629	31	21.195	46	21.629
2	20.481	17	20.658	32	21.235	47	21.635
3	20.481	18	20.690	33	21.276	48	21.639
4	20.481	19	20.724	34	21.317	49	21.641
5	20.482	20	20.759	35	21.356	50	21.640
6	20.483	21	20.795	36	21.395	51	21.639
7	20.485	22	20.832	37	21.432	52	21.634
8	20.490	23	20.871	38	21.467	53	21.629
9	20.496	24	20.910	39	21.498	54	21.622
10	20.507	25	20.950	40	21.527	55	21.615
11	20.520	26	20.990	41	21.553	56	21.606
12	20.535	27	21.031	42	21.575	57	21.597
13	20.554	28	21.072	43	21.593	58	21.587
14	20.577	29	21.113	44	21.608	59	21.576
15	20.601	30	21.153	45	21.620	60	21.565

T-247-300119-199,Potential: Oven Dry - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines

Fractional Organic Carbon



Summary of Fractional Organic Carbon Tests

Sample Number	Fractional Organic Carbon (%)
FOC-247-310119-120	ND
FOC-247-310119-164	ND

ND= not detected at the reporting limit of 0.1% C

Analysis provided by Hall Environmental, Albuquerque, NM.

Analytical ReportLab Order: **1902031**Date Reported: **2/11/2019****Hall Environmental Analysis Laboratory, Inc.****CLIENT:** Daniel B. Stephens & Assoc.**Lab Order:** 1902031**Project:** DBSA EA Engineering**Lab ID:** 1902031-001**Collection Date:** 1/31/2019 2:30:00 PM**Client Sample ID:** FOC-247-300119-120**Matrix:** SOIL

Analyses	Result	RL	Qual	Units	DF	Date Analyzed	Batch ID
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WALKLEY BLACK TOC/FOC/OMAnalyst: **JRR**

FOC	ND	0.097		% C	1	2/7/2019 11:23:00 AM	43029
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Lab ID: 1902031-002**Collection Date:** 1/31/2019 2:30:00 PM**Client Sample ID:** FOC-247-310119-164**Matrix:** SOIL

Analyses	Result	RL	Qual	Units	DF	Date Analyzed	Batch ID
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WALKLEY BLACK TOC/FOC/OMAnalyst: **JRR**

FOC	ND	0.098		% C	1	2/7/2019 11:23:00 AM	43029
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Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

Qualifiers:	*	Value exceeds Maximum Contaminant Level.	B	Analyte detected in the associated Method Blank
	D	Sample Diluted Due to Matrix	E	Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limits
	ND	Not Detected at the Reporting Limit	P	Sample pH Not In Range
	PQL	Practical Quantitative Limit	RL	Reporting Detection Limit

Page 1 of 2

QC SUMMARY REPORT

Hall Environmental Analysis Laboratory, Inc.

WO#: 1902031

11-Feb-19

Client: Daniel B. Stephens & Assoc.

Project: DBSA EA Engineering

Sample ID	MB-43029		SampType: MBLK		TestCode: Walkley Black TOC/FOC/OM					
Client ID:	PBS		Batch ID: 43029		RunNo: 57534					
Prep Date:	2/7/2019		Analysis Date: 2/7/2019		SeqNo: 1924845		Units: % C			
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	%RPD	RPDLimit	Qual
FOC	ND	0.098								

Sample ID	LCS-43029		SampType: LCS		TestCode: Walkley Black TOC/FOC/OM					
Client ID:	LCSS		Batch ID: 43029		RunNo: 57534					
Prep Date:	2/7/2019		Analysis Date: 2/7/2019		SeqNo: 1924846		Units: % C			
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	%RPD	RPDLimit	Qual
FOC	2.1	0.10	2.100	0	101	80	120			

Sample ID	1902031-002AMS		SampType: MS		TestCode: Walkley Black TOC/FOC/OM					
Client ID:	FOC-247-310119-16		Batch ID: 43029		RunNo: 57534					
Prep Date:	2/7/2019		Analysis Date: 2/7/2019		SeqNo: 1924849		Units: % C			
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	%RPD	RPDLimit	Qual
FOC	2.0	0.096	2.100	0	97.6	75	125			

Sample ID	1902031-002AMSD		SampType: MSD		TestCode: Walkley Black TOC/FOC/OM					
Client ID:	FOC-247-310119-16		Batch ID: 43029		RunNo: 57534					
Prep Date:	2/7/2019		Analysis Date: 2/7/2019		SeqNo: 1924850		Units: % C			
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	%RPD	RPDLimit	Qual
FOC	2.0	0.097	2.100	0	97.6	75	125	0	20	

Qualifiers:

* Value exceeds Maximum Contaminant Level.
D Sample Diluted Due to Matrix
H Holding times for preparation or analysis exceeded
ND Not Detected at the Reporting Limit
PQL Practical Quantitative Limit
S % Recovery outside of range due to dilution or matrix

B Analyte detected in the associated Method Blank
E Value above quantitation range
J Analyte detected below quantitation limits
P Sample pH Not In Range
RL Reporting Detection Limit
W Sample container temperature is out of limit as specified

Page 2 of 2

Laboratory Tests and Methods



Tests and Methods

Dry Bulk Density:	ASTM D7263
Moisture Content:	ASTM D7263, ASTM D2216
Calculated Porosity:	ASTM D7263
TOC/FOC:	Walkley Black; Nelson, D. and L. Sommers. 1996. Total Carbon, Organic Carbon, and Organic Matter. Chp. 34, pp. 995-1001, Walkley Black Method, in D. Sparks (ed.), Methods of Soil Analysis. Part 3. American Society of Agronomy, Madison, WI
Thermal Properties:	ASTM D5334



Daniel B. Stephens & Associates, Inc.

Copy

COC-247-DBS-20190131

Daniel B. Stephens & Associates, Inc.
Hydraulic Properties Soils Laboratory
4400 Alameda Blvd. NE, Suite C
ALBUQUERQUE, NM 87113
Ph: (505) 889-7752 (800) 933-3105
FAX (505) 889-0258

Name/Address/Phone for report:

Bernie Bockisch (bbockisch@eaest.com)
320 Gold Ave SW Ste. 1300
Albuquerque, NM 87102

cc: Name/Address/Phone for invoice: (or same as report)

Dallas Accounts Payable
email to: dallasap@eaest.com

Project Name or P.O. #	# of containers	Type of Container	As Received	Gravimetric Moisture Content	Initial Soil Properties (density, moisture, porosity)	Saturated Hydraulic Conductivity (rigid wall or flexible wall)	Hydraulic Properties Package (HPP) with Flexible Wall Ksat	Hydraulic Properties Package (HPP) with Rigid Wall Ksat	Particle Size Analysis (wet sieve & hydrometer)	Proctor Compaction (standard or modified)	Specific Gravity, Fines	Specific Gravity, Coarse	Atterberg Limits	Thermal Properties, 1-point	Thermal Properties, 5-6-point	REMARKS
02/30/19 T-247-300119-144	1	BAG														
02/30/19 FOC-247-300119-120	1	803 JPR														
01/31/19 FOC-247-310119-164	↓	↓														
JR L																
RELINQUISHED BY: (SIGNATURE) DATE RECEIVED BY: (SIGNATURE) DATE																
Comments:																

Laboratory Report for EA Engineering, Science and Technology, Inc.

KAFB-Bulk Fuels Fac Vadose Zone

Project #: 62735DM02, PO: 18189

February 28, 2019



Daniel B. Stephens & Associates, Inc.

4400 Alameda Blvd. NE, Suite C • Albuquerque, New Mexico 87113



February 28, 2019

Bernie Bockisch
EA Engineering, Science and Technology, Inc.
320 Gold Ave SW, Ste. 1300
Albuquerque, NM 87102
(505) 234-1105

Re: DBS&A Laboratory Report for the EA Engineering, Science, and Technology, Inc.
62735DM02 KAFB-Bulk Fuels Fac Vadose Zone Project

Dear Mr. Bockisch:

Enclosed is the report for the EA Engineering, Science, and Technology, Inc. 62735DM02 KAFB-Bulk Fuels Fac Vadose Zone project samples. Please review this report and provide any comments as samples will be held for a maximum of 30 days. After 30 days samples will be returned or disposed of in an appropriate manner.

All testing results were evaluated subjectively for consistency and reasonableness, and the results appear to be reasonably representative of the material tested. However, DBS&A does not assume any responsibility for interpretations or analyses based on the data enclosed, nor can we guarantee that these data are fully representative of the undisturbed materials at the field site. We recommend that careful evaluation of these laboratory results be made for your particular application.

The testing utilized to generate the enclosed report employs methods that are standard for the industry. The results do not constitute a professional opinion by DBS&A, nor can the results affect any professional or expert opinions rendered with respect thereto by DBS&A. You have acknowledged that all the testing undertaken by us, and the report provided, constitutes mere test results using standardized methods, and cannot be used to disqualify DBS&A from rendering any professional or expert opinion, having waived any claim of conflict of interest by DBS&A.

We are pleased to provide this service to EA Engineering and look forward to future laboratory testing on other projects. If you have any questions about the enclosed data, please do not hesitate to call.

Sincerely,

DANIEL B. STEPHENS & ASSOCIATES, INC.
SOIL TESTING & RESEARCH LABORATORY

A handwritten signature in black ink, appearing to read 'Joleen Hines', is written over a light blue horizontal line.

Joleen Hines
Laboratory Manager

Enclosure

Daniel B. Stephens & Associates, Inc.
Soil Testing & Research Laboratory

4400 Alameda Blvd. NE, Suite C
Albuquerque, NM 87113

505-889-7752
FAX 505-889-0258

Summaries



Summary of Tests Performed

Laboratory Sample Number	Initial Soil Properties ¹			Saturated Hydraulic Conductivity ²			Moisture Characteristics ³								Particle Size ⁴			Specific Gravity ⁵		Air Perm- eability	FOC ⁶	Thermal Properties
	G	VM	VD	CH	FH	FW	HC	PP	FP	DPP	RH	EP	WHC	K _{unsat}	DS	WS	H	F	C			
T-247-050219-490	X	X																				X
FOC-247-040219-474																					X	
FOC-247-050219-480																					X	
FOC-247-050219-489																					X	

¹ G = Gravimetric Moisture Content, VM = Volume Measurement Method, VD = Volume Displacement Method

² CH = Constant Head Rigid Wall, FH = Falling Head Rigid Wall, FW = Falling Head Rising Tail Flexible Wall

³ HC = Hanging Column, PP = Pressure Plate, FP = Filter Paper, DPP = Dew Point Potentiometer, RH = Relative Humidity Box, EP = Effective Porosity, WHC = Water Holding Capacity, K_{unsat} = Calculated Unsaturated Hydraulic Conductivity

⁴ DS = Dry Sieve, WS = Wet Sieve, H = Hydrometer

⁵ F = Fine (<4.75mm), C = Coarse (>4.75mm)

⁶FOC = Fractional Organic Carbon testing was performed by Hall Environmental Analysis Laboratory in Albuquerque, NM.



Notes

Sample Receipt:

A total of four samples were hand-delivered on February 5, 2019. One sample, was received in a full 6" x 24" plastic bag sealed with duct tape inside a cardboard box. The remaining three samples were received, each in an 8oz glass jar sealed with a lid inside a quart resealable bag. The three samples were received inside a cooler surrounded with ice. All samples were received in good order.

Sample Preparation and Testing Notes:

A sub-sample was obtained for sample T-247-050219-490 by gently advancing a ring into the specimen and lightly packing approximating 85% of standard proctor compactive effort.

The sub-sample was then subjected to thermal properties testing at the initial moisture content, the saturated moisture content, at several air drying moisture contents, and at the oven dry state.

Each thermal properties reading was obtained in the same location.

Samples FOC-247-040219-474, FOC-247-050219-480 and FOC-247-050219-489 were subjected to Fractional Organic Carbon (FOC) testing. The FOC testing was performed by Hall Environmental Analysis Laboratory in Albuquerque, NM.

Porosity calculations are based on the use of an assumed specific gravity value of 2.65.



Summary of Sample Preparation/Volume Changes

Sample Number	Initial Sample Data ¹		Volume Change Post Saturation ²			Volume Change Post Oven Drying ³		
	Moisture Content (%, g/g)	Dry Bulk Density (g/cm ³)	Dry Bulk Density (g/cm ³)	% Volume Change (%)	% of Initial Density (%)	Dry Bulk Density (g/cm ³)	% Volume Change (%)	% of Initial Density (%)
T-247-050219-490	9.3	1.93	1.93	---	100.0%	1.93	---	100.0%

¹Initial Sample Data: The 'as received' moisture content and lightly packing approximating 85% of standard proctor compactive effort.

²Volume Change Post Saturation: Volume change measurements were obtained after saturation.

³Volume Change Post Oven Drying: Volume change measurements were obtained after oven drying.

Notes:

"+" indicates sample swelling, "-" indicates sample settling, and "---" indicates no volume change occurred.



**Summary of Initial Moisture Content, Dry Bulk Density
Wet Bulk Density and Calculated Porosity**

Sample Number	Moisture Content				Dry Bulk Density (g/cm ³)	Wet Bulk Density (g/cm ³)	Calculated Porosity (%)
	As Received		Remolded				
	Gravimetric	Volumetric	Gravimetric	Volumetric			
	(%, g/g)	(%, cm ³ /cm ³)	(%, g/g)	(%, cm ³ /cm ³)			
T-247-050219-490	NA	NA	9.3	17.9	1.93	2.11	27.3

NA = Not analyzed

--- = This sample was not remolded



Summary of Thermal Properties

Sample	Reading	Gravimetric Moisture Content (g/g, %)	Volumetric Moisture Content ¹ (vol/vol, %)	Dry Bulk Density ¹ (g/cm ³)	Temp °C	K W/(m·K)	ρ °C·cm/W	C MJ/(m ³ ·K)	D mm ² /s
T-247-050219-490	Initial	9.28	17.88	1.93	19.76	1.460	68.5	1.676	0.871
T-247-050219-490	0	13.92	26.82	1.93	17.90	1.629	61.4	2.469	0.660
T-247-050219-490	Air Dry #1	10.32	19.88	1.93	16.19	1.307	76.5	2.411	0.542
T-247-050219-490	Air Dry #2	7.33	14.12	1.93	18.37	0.862	116.0	1.934	0.446
T-247-050219-490	Air Dry #3	4.64	8.94	1.93	18.03	0.609	164.3	1.645	0.370
T-247-050219-490	Air Dry #4	2.02	3.90	1.93	19.02	0.464	215.4	1.489	0.312
T-247-050219-490	Air Dry #5	1.69	3.26	1.93	19.51	0.459	218.0	1.781	0.257
T-247-050219-490	Oven Dry	0	0	1.93	20.89	0.387	258.7	1.982	0.195

¹ Adjusted for volume changes during testing, if applicable.



Summary of Fractional Organic Carbon Tests

Sample Number	Fractional Organic Carbon (%)
FOC-247-040219-474	ND
FOC-247-050219-480	ND
FOC-247-050219-489	ND

ND= not detected at the reporting limit of 0.1% C

Analysis provided by Hall Environmental, Albuquerque, NM.

Initial Properties



**Summary of Initial Moisture Content, Dry Bulk Density
Wet Bulk Density and Calculated Porosity**

Sample Number	Moisture Content				Dry Bulk Density (g/cm ³)	Wet Bulk Density (g/cm ³)	Calculated Porosity (%)
	As Received		Remolded				
	Gravimetric	Volumetric	Gravimetric	Volumetric			
	(%, g/g)	(%, cm ³ /cm ³)	(%, g/g)	(%, cm ³ /cm ³)			
T-247-050219-490	NA	NA	9.3	17.9	1.93	2.11	27.3

NA = Not analyzed

--- = This sample was not remolded



**Data for Initial Moisture Content,
Bulk Density, Porosity, and Percent Saturation**

Job Name: EA Engineering, Science, and Technology, Inc.
Job Number: DB18.1333.00
Sample Number: T-247-050219-490
Project Name: 62735DM02 KAFB-Bulk Fuels Fac Vadose Zone
PO Number: 18189

	<u>As Received</u>	<u>Remolded</u>
<i>Test Date:</i>	NA	11-Feb-19
<i>Field weight* of sample (g):</i>		884.42
<i>Tare weight, ring (g):</i>		214.78
<i>Tare weight, pan/plate (g):</i>		0.00
<i>Tare weight, other (g):</i>		0.00
<i>Dry weight of sample (g):</i>		612.78
<i>Sample volume (cm³):</i>		318.03
<i>Assumed particle density (g/cm³):</i>		2.65
<i>Gravimetric Moisture Content (% g/g):</i>		9.3
<i>Volumetric Moisture Content (% vol):</i>		17.9
<i>Dry bulk density (g/cm³):</i>		1.93
<i>Wet bulk density (g/cm³):</i>		2.11
<i>Calculated Porosity (% vol):</i>		27.3
<i>Percent Saturation:</i>		65.5
<i>Laboratory analysis by:</i>	D. O'Dowd	
<i>Data entered by:</i>	A. Bland	
<i>Checked by:</i>	J. Hines	

Comments:

* Weight including tares
NA = Not analyzed
--- = This sample was not remolded

Thermal Properties



Summary of Thermal Properties

Sample	Reading	Gravimetric Moisture Content (g/g, %)	Volumetric Moisture Content ¹ (vol/vol, %)	Dry Bulk Density ¹ (g/cm ³)	Temp °C	K W/(m·K)	ρ °C·cm/W	C MJ/(m ³ ·K)	D mm ² /s
T-247-050219-490	Initial	9.28	17.88	1.93	19.76	1.460	68.5	1.676	0.871
T-247-050219-490	0	13.92	26.82	1.93	17.90	1.629	61.4	2.469	0.660
T-247-050219-490	Air Dry #1	10.32	19.88	1.93	16.19	1.307	76.5	2.411	0.542
T-247-050219-490	Air Dry #2	7.33	14.12	1.93	18.37	0.862	116.0	1.934	0.446
T-247-050219-490	Air Dry #3	4.64	8.94	1.93	18.03	0.609	164.3	1.645	0.370
T-247-050219-490	Air Dry #4	2.02	3.90	1.93	19.02	0.464	215.4	1.489	0.312
T-247-050219-490	Air Dry #5	1.69	3.26	1.93	19.51	0.459	218.0	1.781	0.257
T-247-050219-490	Oven Dry	0	0	1.93	20.89	0.387	258.7	1.982	0.195

¹ Adjusted for volume changes during testing, if applicable.



Thermal Properties Results Sheet for Sample: T-247-050219-490

Job Name: EA Engineering, Inc.

Job Number: DB18.1333.00

Sample Number: T-247-050219-490

Project Name: 62735DM02 KAFB-Bulk Fuels Fac Vadose Zone

PO Number: 18189

Instrument Description: Decagon KD2 Pro

Probe: ☐ KS-1, 6 cm length, 1.3 mm diameter, single needle

☐ TR-1, 10 cm length, 2.4 mm diameter, single needle

☒ SH-1, 3 cm length, 1.3 mm diameter, dual needle, 6 mm spacing

Test Start Date: 2/11/19

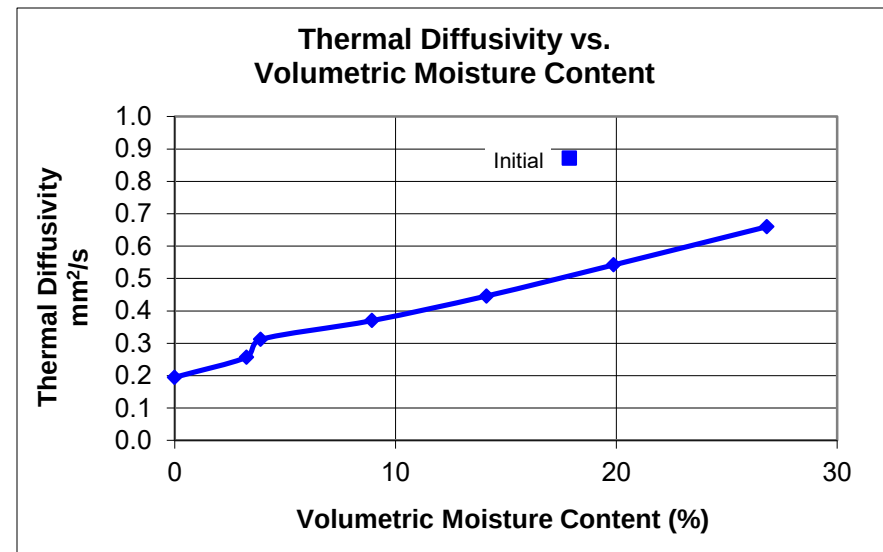
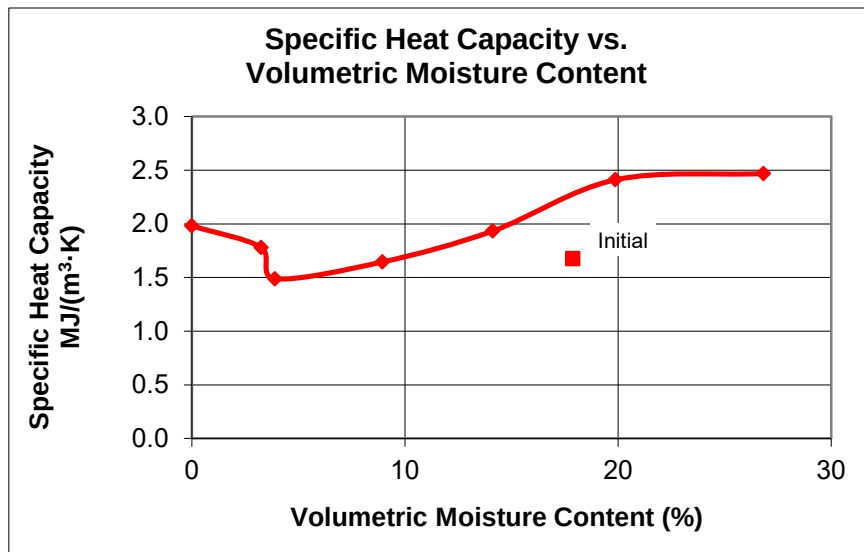
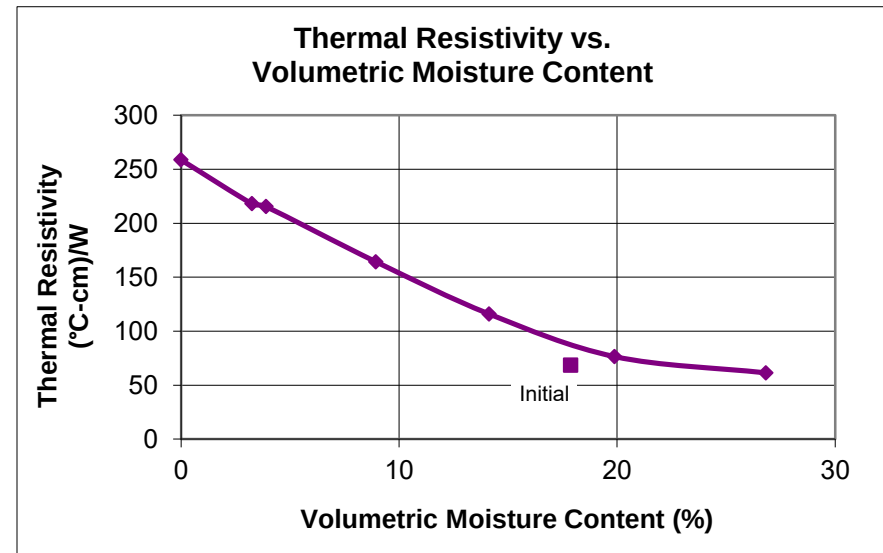
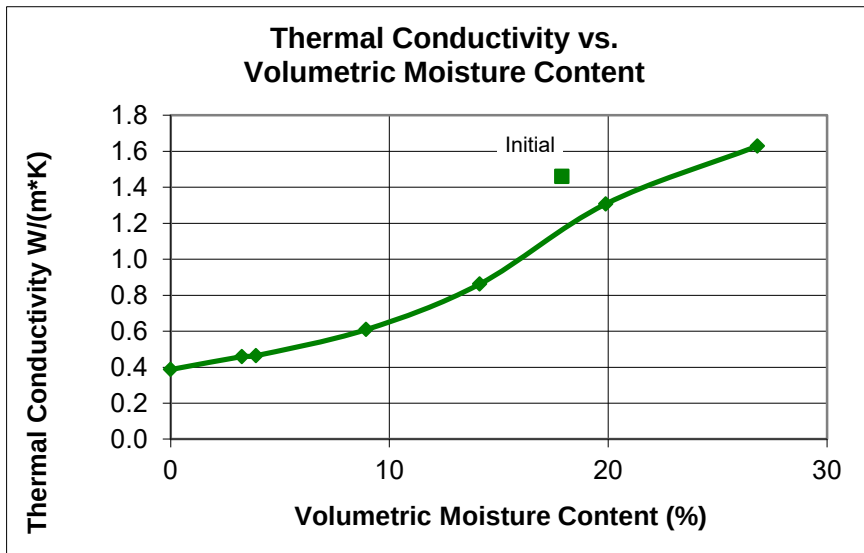
Reading	Water Potential (-cm water)	Gravimetric Moisture Content (g/g, %)	Volumetric Moisture Content ¹ (vol/vol, %)	Dry Bulk Density ¹ (g/cm ³)	Test Temperature (°C)	K Thermal Conductivity W/(m·K)	ρ Thermal Resistivity °C·cm/W	C Specific Heat Capacity MJ/(m ³ ·K)	D Thermal Diffusivity (mm ² /s)
Initial	---	9.28	17.88	1.93	19.76	1.460	68.5	1.676	0.871
Saturated	0	13.92	26.82	1.93	17.90	1.629	61.4	2.469	0.660
Air Dry #1	---	10.32	19.88	1.93	16.19	1.307	76.5	2.411	0.542
Air Dry #2	---	7.33	14.12	1.93	18.37	0.862	116.0	1.934	0.446
Air Dry #3	---	4.64	8.94	1.93	18.03	0.609	164.3	1.645	0.370
Air Dry #4	---	2.02	3.90	1.93	19.02	0.464	215.4	1.489	0.312
Air Dry #5	---	1.69	3.26	1.93	19.51	0.459	218.0	1.781	0.257
Oven Dry	---	0	0	1.93	20.89	0.387	258.7	1.982	0.195

--- = Value not measured.

¹ Adjusted for volume changes during testing, if applicable.



Thermal Properties Results Sheet for Sample: T-247-050219-490
Scatter Plots





Thermal Properties Data

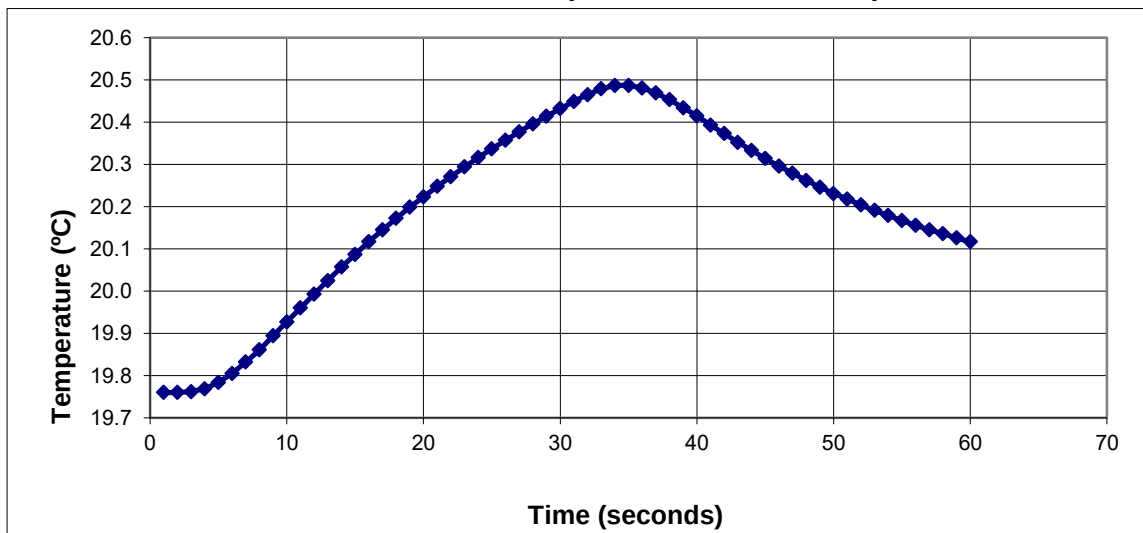
Sample Number: T-247-050219-490
Potential (-cm water): Initial

Test Date/Time: 2/11/2019 9:20 AM	K (W/(m·K)): 1.460
Sensor: SH-1	ρ (°C·cm/W): 68.5
Test Temp.(°C): 19.8	C (MJ/(m ³ ·K)): 1.676
KD2 Pro Sample ID: 490-AR	D (mm ² /s): 0.871
Power (W/m): 18.360	Err: 0.0077
Current (amps): 0.133	

Raw Data

Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)
1 19.760	16 20.117	31 20.449	46 20.296
2 19.760	17 20.145	32 20.465	47 20.279
3 19.762	18 20.172	33 20.479	48 20.262
4 19.769	19 20.199	34 20.487	49 20.246
5 19.784	20 20.223	35 20.487	50 20.231
6 19.805	21 20.248	36 20.481	51 20.218
7 19.832	22 20.271	37 20.469	52 20.204
8 19.861	23 20.294	38 20.453	53 20.191
9 19.894	24 20.316	39 20.434	54 20.179
10 19.927	25 20.337	40 20.415	55 20.167
11 19.960	26 20.357	41 20.393	56 20.156
12 19.993	27 20.377	42 20.373	57 20.145
13 20.025	28 20.396	43 20.352	58 20.136
14 20.057	29 20.414	44 20.333	59 20.126
15 20.087	30 20.432	45 20.314	60 20.117

T-247-050219-490,Potential: Initial - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: A. Bland
Checked by: J. Hines



Thermal Properties Data

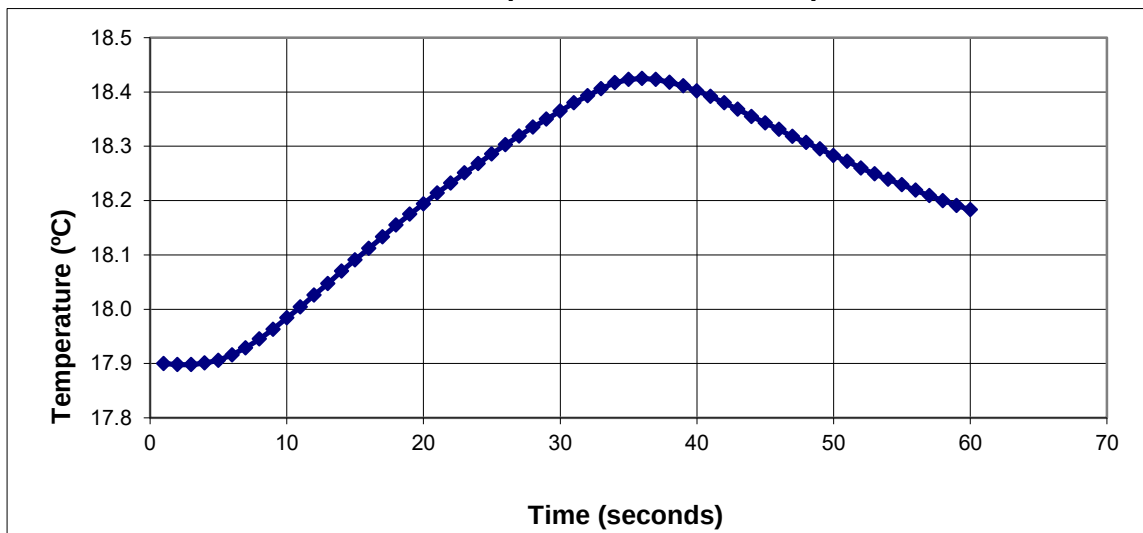
Sample Number: T-247-050219-490
Potential (-cm water): 0

Test Date/Time: 2/13/2019 8:16 AM	K (W/(m·K)): 1.629
Sensor: SH-1	ρ (°C·cm/W): 61.4
Test Temp.(°C): 17.9	C (MJ/(m ³ ·K)): 2.469
KD2 Pro Sample ID: 490-SA	D (mm ² /s): 0.660
Power (W/m): 19.140	Err: 0.0030
Current (amps): 0.136	

Raw Data

Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)
1 17.900	16 18.112	31 18.380	46 18.331
2 17.898	17 18.133	32 18.393	47 18.318
3 17.898	18 18.155	33 18.406	48 18.307
4 17.901	19 18.175	34 18.417	49 18.295
5 17.906	20 18.194	35 18.423	50 18.283
6 17.916	21 18.214	36 18.425	51 18.272
7 17.929	22 18.232	37 18.423	52 18.260
8 17.945	23 18.251	38 18.418	53 18.249
9 17.963	24 18.268	39 18.411	54 18.239
10 17.984	25 18.286	40 18.402	55 18.229
11 18.004	26 18.303	41 18.392	56 18.219
12 18.026	27 18.319	42 18.380	57 18.209
13 18.047	28 18.335	43 18.368	58 18.200
14 18.070	29 18.350	44 18.355	59 18.191
15 18.091	30 18.365	45 18.343	60 18.183

T-247-050219-490,Potential: 0 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: A. Bland
Checked by: J. Hines



Thermal Properties Data

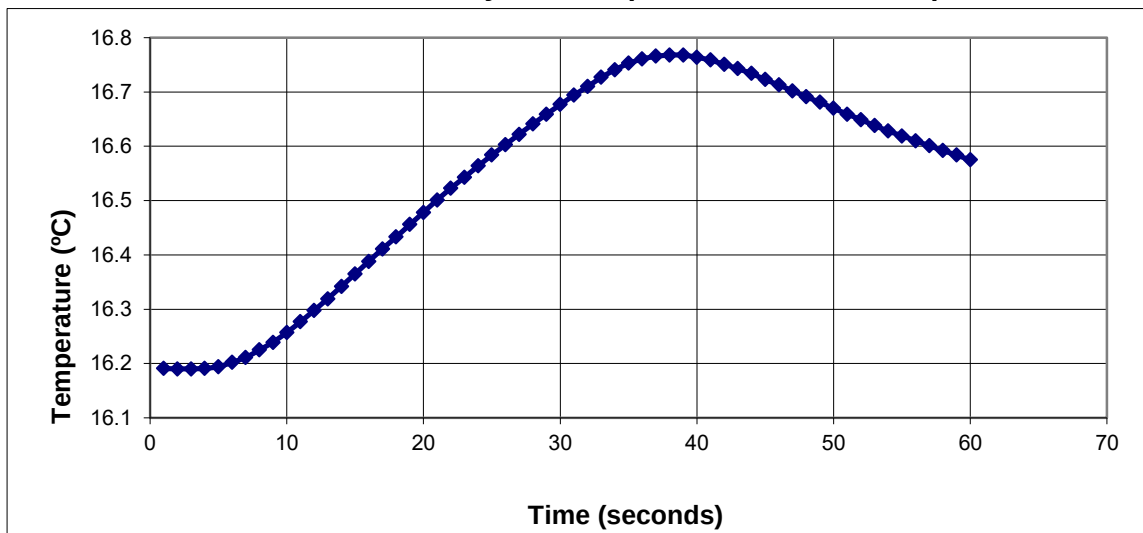
Sample Number: T-247-050219-490
Potential (-cm water): Air Dry #1

Test Date/Time:	2/13/2019 9:12 AM	K (W/(m·K)):	1.307
Sensor:	SH-1	ρ (°C·cm/W):	76.5
Test Temp. (°C):	16.2	C (MJ/(m ³ ·K)):	2.411
KD2 Pro Sample ID:	490-AD1	D (mm ² /s):	0.542
Power (W/m):	19.110	Err:	0.0032
Current (amps):	0.136		

Raw Data

Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)	
1	16.191	16	16.388	31	16.694	46	16.713
2	16.190	17	16.411	32	16.710	47	16.702
3	16.190	18	16.433	33	16.727	48	16.691
4	16.191	19	16.456	34	16.741	49	16.681
5	16.194	20	16.478	35	16.753	50	16.670
6	16.202	21	16.501	36	16.761	51	16.659
7	16.211	22	16.523	37	16.766	52	16.649
8	16.225	23	16.543	38	16.768	53	16.638
9	16.239	24	16.564	39	16.768	54	16.628
10	16.257	25	16.584	40	16.764	55	16.619
11	16.277	26	16.603	41	16.759	56	16.610
12	16.298	27	16.622	42	16.751	57	16.601
13	16.319	28	16.641	43	16.743	58	16.592
14	16.342	29	16.659	44	16.734	59	16.584
15	16.365	30	16.677	45	16.723	60	16.575

T-247-050219-490,Potential: Air Dry #1 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: A. Bland
Checked by: J. Hines



Thermal Properties Data

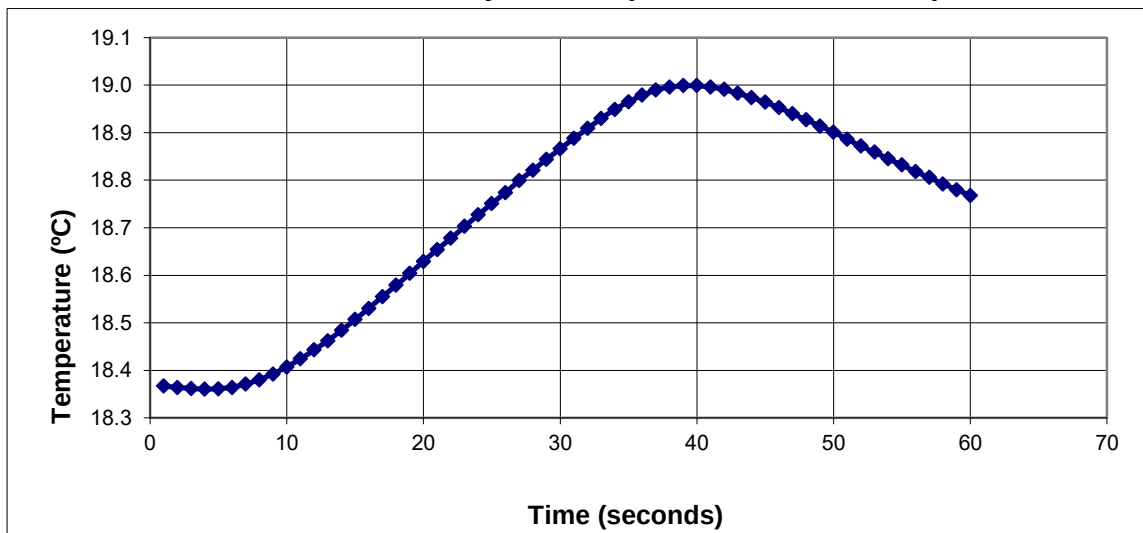
Sample Number: T-247-050219-490
Potential (-cm water): Air Dry #2

Test Date/Time:	2/14/2019 8:52 AM	K (W/(m·K)):	0.862
Sensor:	SH-1	ρ (°C·cm/W):	116.0
Test Temp.(°C):	18.4	C (MJ/(m ³ ·K)):	1.934
KD2 Pro Sample ID:	490-AD2	D (mm ² /s):	0.446
Power (W/m):	18.660	Err:	0.0023
Current (amps):	0.134		

Raw Data

Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)	
1	18.367	16	18.530	31	18.888	46	18.953
2	18.364	17	18.555	32	18.909	47	18.940
3	18.362	18	18.579	33	18.930	48	18.927
4	18.360	19	18.604	34	18.949	49	18.914
5	18.361	20	18.629	35	18.965	50	18.901
6	18.364	21	18.654	36	18.979	51	18.886
7	18.371	22	18.678	37	18.990	52	18.872
8	18.380	23	18.703	38	18.996	53	18.859
9	18.392	24	18.727	39	18.999	54	18.845
10	18.407	25	18.751	40	18.999	55	18.832
11	18.424	26	18.774	41	18.996	56	18.818
12	18.443	27	18.799	42	18.991	57	18.806
13	18.462	28	18.821	43	18.983	58	18.792
14	18.484	29	18.844	44	18.974	59	18.780
15	18.507	30	18.866	45	18.964	60	18.768

T-247-050219-490,Potential: Air Dry #2 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: A. Bland
Checked by: J. Hines



Thermal Properties Data

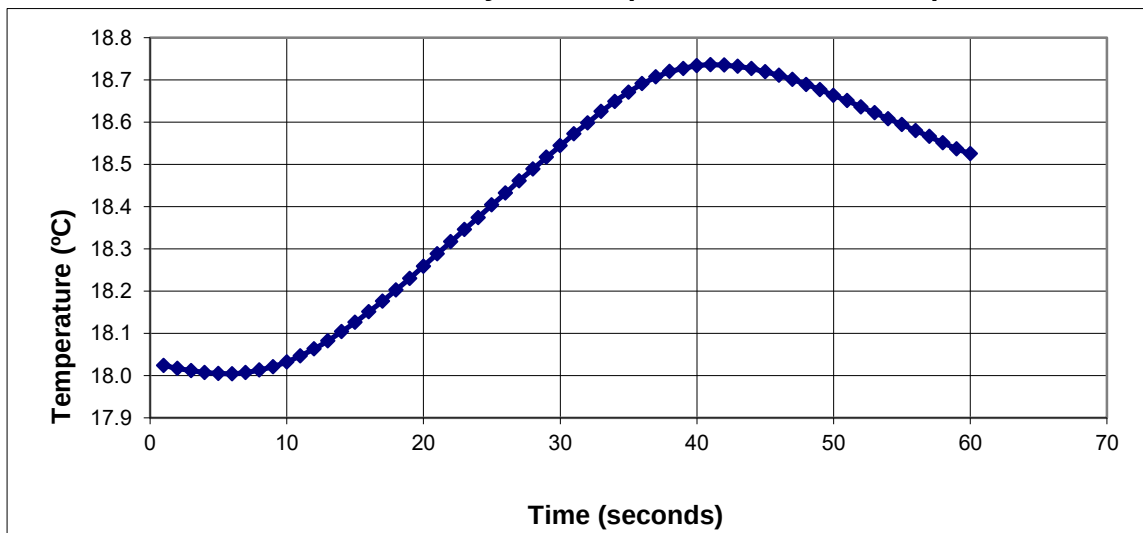
Sample Number: T-247-050219-490
Potential (-cm water): Air Dry #3

Test Date/Time: 2/15/2019 7:35 AM	K (W/(m·K)): 0.609
Sensor: SH-1	ρ (°C·cm/W): 164.3
Test Temp.(°C): 18.0	C (MJ/(m ³ ·K)): 1.645
KD2 Pro Sample ID: 490-AD3	D (mm ² /s): 0.370
Power (W/m): 18.870	Err: 0.0013
Current (amps): 0.135	

Raw Data

Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)
1 18.024	16 18.151	31 18.572	46 18.711
2 18.017	17 18.176	32 18.598	47 18.701
3 18.012	18 18.203	33 18.625	48 18.689
4 18.007	19 18.230	34 18.649	49 18.677
5 18.005	20 18.259	35 18.671	50 18.663
6 18.004	21 18.288	36 18.691	51 18.651
7 18.007	22 18.317	37 18.707	52 18.636
8 18.013	23 18.346	38 18.720	53 18.622
9 18.021	24 18.374	39 18.727	54 18.608
10 18.032	25 18.404	40 18.734	55 18.594
11 18.047	26 18.432	41 18.736	56 18.580
12 18.063	27 18.461	42 18.735	57 18.566
13 18.082	28 18.489	43 18.732	58 18.551
14 18.104	29 18.517	44 18.727	59 18.537
15 18.126	30 18.544	45 18.719	60 18.525

T-247-050219-490,Potential: Air Dry #3 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: A. Bland
Checked by: J. Hines



Thermal Properties Data

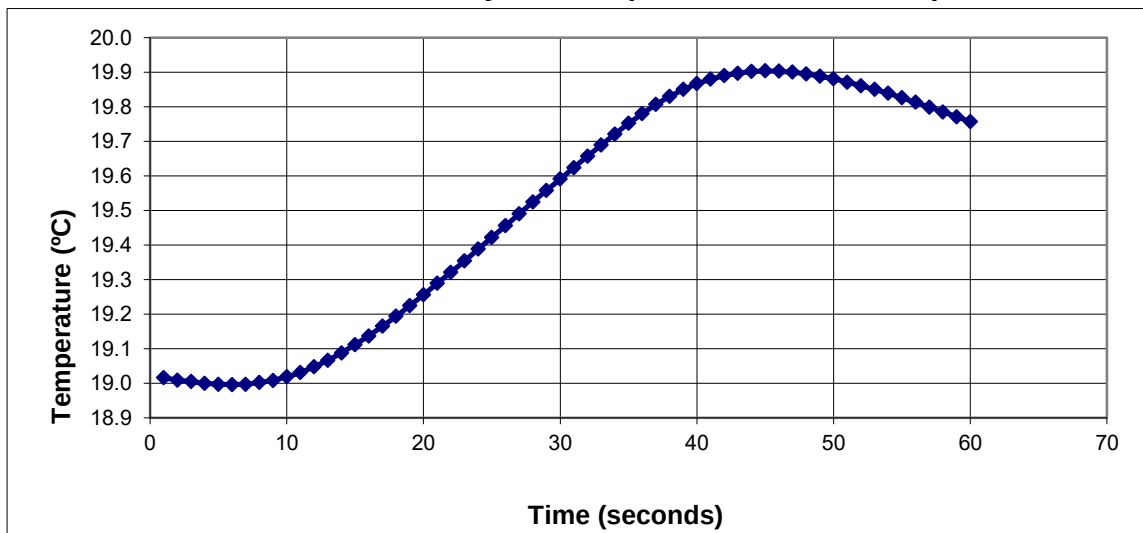
Sample Number: T-247-050219-490
Potential (-cm water): Air Dry #4

Test Date/Time: 2/18/2019 7:55 AM	K (W/(m·K)): 0.464
Sensor: SH-1	ρ (°C·cm/W): 215.4
Test Temp.(°C): 19.0	C (MJ/(m ³ ·K)): 1.489
KD2 Pro Sample ID: 490-AD4	D (mm ² /s): 0.312
Power (W/m): 19.360	Err: 0.0011
Current (amps): 0.136	

Raw Data

Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)
1 19.016	16 19.137	31 19.624	46 19.903
2 19.009	17 19.165	32 19.657	47 19.900
3 19.005	18 19.194	33 19.689	48 19.895
4 19.000	19 19.225	34 19.721	49 19.888
5 18.997	20 19.256	35 19.752	50 19.881
6 18.996	21 19.289	36 19.780	51 19.871
7 18.997	22 19.321	37 19.807	52 19.861
8 19.002	23 19.354	38 19.830	53 19.850
9 19.008	24 19.388	39 19.850	54 19.839
10 19.019	25 19.422	40 19.867	55 19.826
11 19.031	26 19.456	41 19.880	56 19.813
12 19.048	27 19.490	42 19.890	57 19.799
13 19.066	28 19.525	43 19.897	58 19.785
14 19.088	29 19.558	44 19.902	59 19.771
15 19.112	30 19.591	45 19.904	60 19.757

T-247-050219-490,Potential: Air Dry #4 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: A. Bland
Checked by: J. Hines



Thermal Properties Data

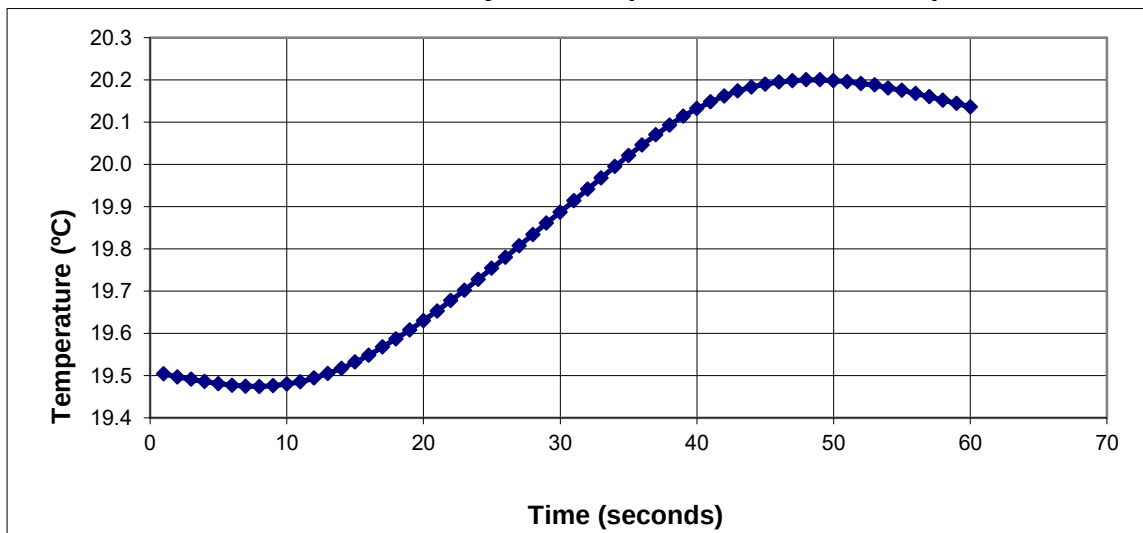
Sample Number: T-247-050219-490
Potential (-cm water): Air Dry #5

Test Date/Time: 2/19/2019 7:43 AM	K (W/(m·K)): 0.459
Sensor: SH-1	ρ (°C·cm/W): 218.0
Test Temp.(°C): 19.5	C (MJ/(m ³ ·K)): 1.781
KD2 Pro Sample ID: 490-AD5	D (mm ² /s): 0.257
Power (W/m): 19.340	Err: 0.0008
Current (amps): 0.136	

Raw Data

Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)
1 19.504	16 19.548	31 19.914	46 20.195
2 19.497	17 19.568	32 19.941	47 20.198
3 19.491	18 19.587	33 19.968	48 20.200
4 19.486	19 19.608	34 19.995	49 20.200
5 19.481	20 19.630	35 20.021	50 20.198
6 19.477	21 19.653	36 20.046	51 20.196
7 19.475	22 19.678	37 20.070	52 20.191
8 19.474	23 19.702	38 20.093	53 20.188
9 19.476	24 19.728	39 20.114	54 20.181
10 19.480	25 19.754	40 20.132	55 20.175
11 19.485	26 19.780	41 20.148	56 20.168
12 19.494	27 19.807	42 20.162	57 20.160
13 19.505	28 19.834	43 20.174	58 20.152
14 19.517	29 19.861	44 20.183	59 20.144
15 19.532	30 19.887	45 20.190	60 20.136

T-247-050219-490,Potential: Air Dry #5 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: A. Bland
Checked by: J. Hines



Thermal Properties Data

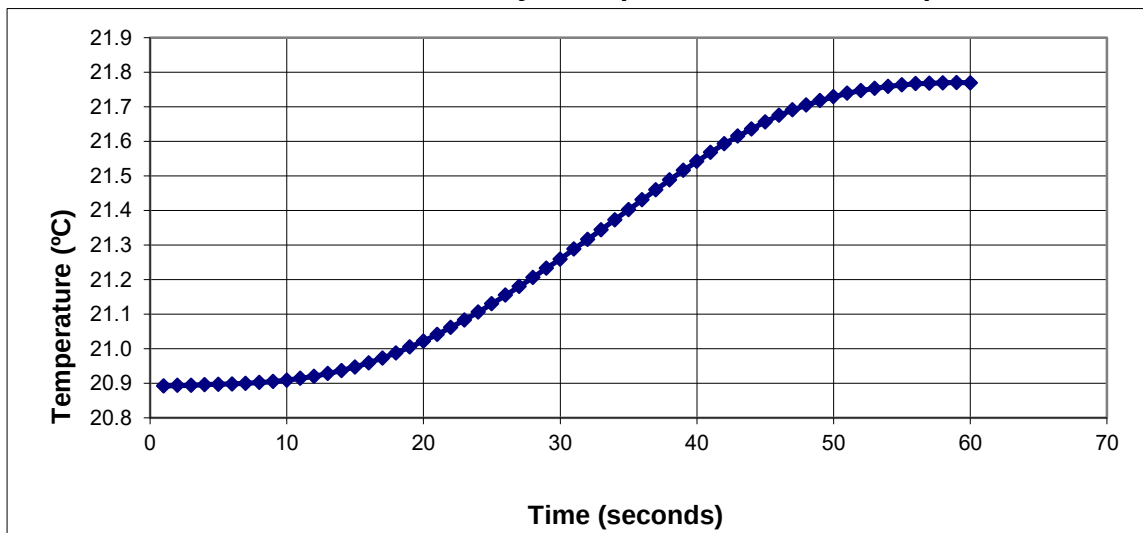
Sample Number: T-247-050219-490
Potential (-cm water): Oven Dry

Test Date/Time: 2/20/2019 12:34 PM	K (W/(m·K)): 0.387
Sensor: SH-1	ρ (°C·cm/W): 258.7
Test Temp.(°C): 20.9	C (MJ/(m ³ ·K)): 1.982
KD2 Pro Sample ID: 490-OD	D (mm ² /s): 0.195
Power (W/m): 19.340	Err: 0.0014
Current (amps): 0.136	

Raw Data

Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)	
1	20.892	16	20.959	31	21.288	46	21.675
2	20.894	17	20.973	32	21.316	47	21.691
3	20.894	18	20.988	33	21.344	48	21.705
4	20.896	19	21.005	34	21.373	49	21.718
5	20.897	20	21.022	35	21.402	50	21.729
6	20.898	21	21.041	36	21.431	51	21.739
7	20.900	22	21.062	37	21.460	52	21.747
8	20.902	23	21.083	38	21.488	53	21.753
9	20.905	24	21.106	39	21.516	54	21.759
10	20.909	25	21.130	40	21.542	55	21.763
11	20.914	26	21.155	41	21.568	56	21.767
12	20.920	27	21.180	42	21.593	57	21.768
13	20.928	28	21.206	43	21.615	58	21.769
14	20.937	29	21.233	44	21.636	59	21.770
15	20.947	30	21.259	45	21.656	60	21.769

T-247-050219-490,Potential: Oven Dry - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: A. Bland
Checked by: J. Hines

Fractional Organic Carbon



Summary of Fractional Organic Carbon Tests

Sample Number	Fractional Organic Carbon (%)
FOC-247-040219-474	ND
FOC-247-050219-480	ND
FOC-247-050219-489	ND

ND= not detected at the reporting limit of 0.1% C

Analysis provided by Hall Environmental, Albuquerque, NM.

Analytical ReportLab Order: **1902412**Date Reported: **2/14/2019****Hall Environmental Analysis Laboratory, Inc.****CLIENT:** Daniel B. Stephens & Assoc.**Lab Order:** 1902412**Project:** DBSA EA Engineering**Lab ID:** 1902412-001**Collection Date:** 2/4/2019 12:00:00 PM**Client Sample ID:** FOC-247-040219-474**Matrix:** SOIL

Analyses	Result	RL	Qual	Units	DF	Date Analyzed	Batch ID
----------	--------	----	------	-------	----	---------------	----------

WALKLEY BLACK TOC/FOC/OMAnalyst: **JRR**

FOC	ND	0.096		% C	1	2/14/2019 9:52:00 AM	43158
-----	----	-------	--	-----	---	----------------------	-------

Lab ID: 1902412-002**Collection Date:** 2/5/2019 12:00:00 PM**Client Sample ID:** FOC-247-050219-480**Matrix:** SOIL

Analyses	Result	RL	Qual	Units	DF	Date Analyzed	Batch ID
----------	--------	----	------	-------	----	---------------	----------

WALKLEY BLACK TOC/FOC/OMAnalyst: **JRR**

FOC	ND	0.10		% C	1	2/14/2019 9:52:00 AM	43158
-----	----	------	--	-----	---	----------------------	-------

Lab ID: 1902412-003**Collection Date:** 2/5/2019 12:00:00 PM**Client Sample ID:** FOC-247-050219-489**Matrix:** SOIL

Analyses	Result	RL	Qual	Units	DF	Date Analyzed	Batch ID
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WALKLEY BLACK TOC/FOC/OMAnalyst: **JRR**

FOC	ND	0.097		% C	1	2/14/2019 9:52:00 AM	43158
-----	----	-------	--	-----	---	----------------------	-------

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

Qualifiers:	*	Value exceeds Maximum Contaminant Level.	B	Analyte detected in the associated Method Blank
	D	Sample Diluted Due to Matrix	E	Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limits
	ND	Not Detected at the Reporting Limit	P	Sample pH Not In Range
	PQL	Practical Quantitative Limit	RL	Reporting Detection Limit

Page 1 of 2

QC SUMMARY REPORT

Hall Environmental Analysis Laboratory, Inc.

WO#: 1902412

14-Feb-19

Client: Daniel B. Stephens & Assoc.

Project: DBSA EA Engineering

Sample ID	MB-43158		SampType:	MBLK		TestCode:	Walkley Black TOC/FOC/OM				
Client ID:	PBS		Batch ID:	43158		RunNo:	57706				
Prep Date:	2/14/2019		Analysis Date:	2/14/2019		SeqNo:	1931164		Units: % C		
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	%RPD	RPDLimit	Qual	
FOC	ND	0.10									

Sample ID	LCS-43158		SampType: LCS		TestCode: Walkley Black TOC/FOC/OM					
Client ID:	LCSS		Batch ID: 43158		RunNo: 57706					
Prep Date:	2/14/2019		Analysis Date: 2/14/2019		SeqNo: 1931165		Units: % C			
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	%RPD	RPDLimit	Qual
FOC	2.2	0.098	2.059	0	106	80	120			

Sample ID	1902412-003AMS		SampType:	MS		TestCode:	Walkley Black TOC/FOC/OM				
Client ID:	FOC-247-050219-48		Batch ID:	43158		RunNo:	57706				
Prep Date:	2/14/2019		Analysis Date:	2/14/2019		SeqNo:	1931169		Units: % C		
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	%RPD	RPDLimit	Qual	
FOC	2.2	0.10	2.100	0	104	75	125				

Sample ID	1902412-003AMSD		SampType:	MSD		TestCode:	Walkley Black TOC/FOC/OM				
Client ID:	FOC-247-050219-48		Batch ID:	43158		RunNo:	57706				
Prep Date:	2/14/2019		Analysis Date:	2/14/2019		SeqNo:	1931170		Units: % C		
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	%RPD	RPDLimit	Qual	
FOC	2.2	0.10	2.100	0	104	75	125	0.458	20		

Qualifiers:

* Value exceeds Maximum Contaminant Level.
D Sample Diluted Due to Matrix
H Holding times for preparation or analysis exceeded
ND Not Detected at the Reporting Limit
PQL Practical Quantitative Limit
S % Recovery outside of range due to dilution or matrix

B Analyte detected in the associated Method Blank
E Value above quantitation range
J Analyte detected below quantitation limits
P Sample pH Not In Range
RL Reporting Detection Limit
W Sample container temperature is out of limit as specified

Page 2 of 2

Laboratory Tests and Methods



Tests and Methods

Dry Bulk Density:	ASTM D7263
Moisture Content:	ASTM D7263, ASTM D2216
Calculated Porosity:	ASTM D7263
TOC/FOC:	Walkley Black; Nelson, D. and L. Sommers. 1996. Total Carbon, Organic Carbon, and Organic Matter. Chp. 34, pp. 995-1001, Walkley Black Method, in D. Sparks (ed.), Methods of Soil Analysis. Part 3. American Society of Agronomy, Madison, WI
Thermal Properties:	ASTM D5334



Copy
Daniel B. Stephens & Associates, Inc.

COC1-247-DBS-20190205

Daniel B. Stephens & Associates, Inc.
Hydraulic Properties Soils Laboratory
4400 Alameda Blvd. NE, Suite C
ALBUQUERQUE, NM 87113
Ph: (505) 889-7752 (800) 933-3105
FAX (505) 889-0258

Name/Address/Phone for report:

Bernie Bockisch (bbockisch@eaest.com)
320 Gold Ave SW Ste. 1300
Albuquerque, NM 87102

cc: Name/Address/Phone for invoice: (or same as report)

Dallas Accounts Payable
email to: dallasap@eaest.com

Project Name or P.O. # **18189**
Kirtland Vadose Zone
Project Contact & Phone:
Bernie Bockisch
505-234-1105

DATE	SAMPLE IDENTIFICATION	# of containers	Type of Container	As Received Gravimetric Moisture Content	Initial Soil Properties (density, moisture, porosity)	Saturated Hydraulic Conductivity (rigid wall or flexible wall)	Hydraulic Properties Package (HPP) with Flexible Wall Ksat	Hydraulic Properties Package (HPP) with Rigid Wall Ksat	Particle Size Analysis (wet sieve & hydrometer)	Proctor Compaction (standard or modified)	Specific Gravity, Fines	Specific Gravity, Coarse	Atterberg Limits	Thermal Properties, 1-point	Thermal Properties, 5-6-point	FOC-Walkley-Black Method	REMARKS
02/05/19	T-247-050219-490	1	BAG	—	—	—	—	—	—	—	—	—	—	—	—	—	—
02/04/19	FOC-247-040219-474	1	507	—	—	—	—	—	—	—	—	—	—	—	—	—	—
02/05/19	FOC-247-050219-480	↓	↓	—	—	—	—	—	—	—	—	—	—	—	—	—	—
02/05/19	FOC-247-050219-489	↓	↓	—	—	—	—	—	—	—	—	—	—	—	—	—	—
JPL																	
RELINQUISHED BY: (SIGNATURE)		DATE		RECEIVED BY: (SIGNATURE)		DATE		Comments:									
John Kim		02/05/19															

Laboratory Report for EA Engineering, Science and Technology, Inc.

KAFB-Bulk Fuels Fac Vadose Zone

Project #: 62735DM02, PO: 18189

**November 8, 2018
Rev. January 10, 2019**



Daniel B. Stephens & Associates, Inc.

4400 Alameda Blvd. NE, Suite C • Albuquerque, New Mexico 87113



November 8, 2018 Rev. January 10, 2019

Bernie Bockisch
EA Engineering, Science and Technology
320 Gold Ave SW, Ste. 1300
Albuquerque, NM 87102
(505) 234-1105

Re: DBS&A Laboratory Report for the EA Engineering, Science, and Technology, Inc.
62735DM02 KAFB-Bulk Fuels Fac Vadose Zone Project

Dear Mr. Bockisch:

Enclosed is the report for the EA Engineering, Science, and Technology, Inc. 62735DM02 KAFB-Bulk Fuels Fac Vadose Zone project samples. Please review this report and provide any comments as samples will be held for a maximum of 30 days. After 30 days samples will be returned or disposed of in an appropriate manner.

All testing results were evaluated subjectively for consistency and reasonableness, and the results appear to be reasonably representative of the material tested. However, DBS&A does not assume any responsibility for interpretations or analyses based on the data enclosed, nor can we guarantee that these data are fully representative of the undisturbed materials at the field site. We recommend that careful evaluation of these laboratory results be made for your particular application.

The testing utilized to generate the enclosed report employs methods that are standard for the industry. The results do not constitute a professional opinion by DBS&A, nor can the results affect any professional or expert opinions rendered with respect thereto by DBS&A. You have acknowledged that all the testing undertaken by us, and the report provided, constitutes mere test results using standardized methods, and cannot be used to disqualify DBS&A from rendering any professional or expert opinion, having waived any claim of conflict of interest by DBS&A.

We are pleased to provide this service to EA Engineering and look forward to future laboratory testing on other projects. If you have any questions about the enclosed data, please do not hesitate to call.

Sincerely,

DANIEL B. STEPHENS & ASSOCIATES, INC.
SOIL TESTING & RESEARCH LABORATORY

A handwritten signature in cursive script, appearing to read 'Joleen Hines'.

Joleen Hines
Laboratory Manager

Enclosure

Daniel B. Stephens & Associates, Inc.
Soil Testing & Research Laboratory

4400 Alameda Blvd. NE, Suite C
Albuquerque, NM 87113

505-889-7752
FAX 505-889-0258

Summaries



Summary of Tests Performed

Laboratory Sample Number	Initial Soil Properties ¹			Saturated Hydraulic Conductivity ²			Moisture Characteristics ³								Particle Size ⁴			Specific Gravity ⁵		Air Perm- eability	Atterberg Limits	Thermal Properties
	G	VM	VD	CH	FH	FW	HC	PP	FP	DPP	RH	EP	WHC	K _{unsat}	DS	WS	H	F	C			
W-S9-161018-435	X																					
WT-S9-171018-459	X	X																				X
WT-S9-171018-464	X	X																				X
WT-S5-231018-490	X	X																				X

¹ G = Gravimetric Moisture Content, VM = Volume Measurement Method, VD = Volume Displacement Method

² CH = Constant Head Rigid Wall, FH = Falling Head Rigid Wall, FW = Falling Head Rising Tail Flexible Wall

³ HC = Hanging Column, PP = Pressure Plate, FP = Filter Paper, DPP = Dew Point Potentiometer, RH = Relative Humidity Box, EP = Effective Porosity, WHC = Water Holding Capacity, K_{unsat} = Calculated Unsaturated Hydraulic Conductivity

⁴ DS = Dry Sieve, WS = Wet Sieve, H = Hydrometer

⁵ F = Fine (<4.75mm), C = Coarse (>4.75mm)



Notes

Revision January 10, 2019:

The correct units of $^{\circ}\text{C}\cdot\text{cm}/\text{W}$ were included for ρ on the sample data sheets in the original report dated November 8, 2018; however, a typo was identified in the units of ρ on the Summary of Thermal Properties page. The typo for the units of ρ on the summary page was corrected from $^{\circ}\text{C}\cdot\text{m}/\text{W}$ to $^{\circ}\text{C}\cdot\text{cm}/\text{W}$ in this revised report. No other changes were made to the original data reported.

Sample Receipt:

Four samples, each in a full 6" x 24" plastic bag sealed with duct tape, were hand-delivered; three samples arrived on October 19, 2018, and one sample arrived on October 24, 2018. Each of the samples was received inside a cardboard box with a lid. All samples were received in good order.

Sample Preparation and Testing Notes:

All of the samples were subjected to as received gravimetric moisture content determination.

Sub-samples were obtained by gently advancing a ring into the specimen for samples WT-S9-171018-459 and WT-S9-171018-464. Sample WT-S5-231018-490 was remolded into a test ring to approximate 85% of maximum density as specified by the client.

The subsamples were then subjected to thermal properties testing at the initial moisture content, the saturated moisture content, at several air drying moisture contents, and at the oven dry state.

Each thermal properties reading was obtained in the same location, whenever possible.

Porosity calculations are based on the use of an assumed specific gravity value of 2.65.

Volumetric water contents were adjusted for changes in volume, where applicable. Due to the irregularities formed on the sample surfaces during settling or swelling, volume measurements obtained after the initial reading should be considered estimates.



**Summary of Initial Moisture Content, Dry Bulk Density
Wet Bulk Density and Calculated Porosity**

Sample Number	Moisture Content				Dry Bulk Density (g/cm ³)	Wet Bulk Density (g/cm ³)	Calculated Porosity (%)
	As Received		Remolded				
	Gravimetric (%, g/g)	Volumetric (%, cm ³ /cm ³)	Gravimetric (%, g/g)	Volumetric (%, cm ³ /cm ³)			
W-S9-161018-435	4.4	NA	---	---	NA	NA	NA
WT-S9-171018-459	24.5	NA	25.6	37.5	1.46	1.84	44.7
WT-S9-171018-464	3.8	NA	4.1	7.4	1.83	1.90	31.0
WT-S5-231018-490	13.1	NA	13.1	24.9	1.90	2.14	28.5

NA = Not analyzed

--- = This sample was not remolded



Summary of Thermal Properties

Sample	Reading	Gravimetric Moisture Content (g/g, %)	Volumetric Moisture Content ¹ (vol/vol, %)	Dry Bulk Density ¹ (g/cm ³)	Temp °C	K W/(m·K)	ρ °C·cm/W	C MJ/(m ³ ·K)	D mm ² /s
WT-S9-171018-459	Initial	25.60	37.49	1.46	22.55	1.429	70.0	3.952	0.362
WT-S9-171018-459	0	31.82	45.30	1.42	21.52	1.429	70.0	3.805	0.376
WT-S9-171018-459	Air Dry #1	27.44	39.45	1.44	18.86	1.370	73.0	3.373	0.406
WT-S9-171018-459	Air Dry #2	23.24	33.93	1.46	18.24	1.360	73.5	3.241	0.420
WT-S9-171018-459	Air Dry #3	18.58	29.18	1.57	19.07	1.293	77.3	3.166	0.408
WT-S9-171018-459	Air Dry #4	13.08	21.03	1.61	19.84	1.060	94.3	2.670	0.397
WT-S9-171018-459	Air Dry #5	11.67	18.78	1.61	20.24	0.996	100.4	2.607	0.382
WT-S9-171018-459	Air Dry #6	10.45	16.82	1.61	20.60	0.941	106.2	2.502	0.376
WT-S9-171018-459	Oven Dry	0	0	1.61	21.22	0.547	182.7	1.554	0.352
WT-S9-171018-464	Initial	4.07	7.44	1.83	22.20	0.681	146.8	1.559	0.437
WT-S9-171018-464	0	16.45	30.06	1.83	22.59	1.686	59.3	3.038	0.555
WT-S9-171018-464	Air Dry #1	14.53	27.13	1.87	21.31	1.676	59.7	2.631	0.637
WT-S9-171018-464	Air Dry #2	12.19	23.21	1.90	20.15	1.587	63.0	2.096	0.757
WT-S9-171018-464	Air Dry #3	8.23	16.09	1.95	17.87	1.167	85.7	1.793	0.651
WT-S9-171018-464	Air Dry #4	6.18	12.07	1.95	18.79	1.001	99.9	2.082	0.481
WT-S9-171018-464	Air Dry #5	4.57	8.94	1.95	18.76	0.810	123.5	1.619	0.500
WT-S9-171018-464	Air Dry #6	3.76	7.35	1.95	20.63	0.656	152.3	1.886	0.348
WT-S9-171018-464	Oven Dry	0	0	1.95	21.47	0.262	381.1	1.471	0.178

¹ Adjusted for volume changes during testing, if applicable.



Summary of Thermal Properties (Continued)

Sample	Reading	Gravimetric Moisture Content (g/g, %)	Volumetric Moisture Content ¹ (vol/vol, %)	Dry Bulk Density ¹ (g/cm ³)	Temp °C	K W/(m·K)	ρ °C·cm/W	C MJ/(m ³ ·K)	D mm ² /s
WT-S5-231018-490	Initial	13.13	24.87	1.90	21.72	1.907	52.4	4.024	0.474
WT-S5-231018-490	0	14.44	27.36	1.90	20.78	1.825	54.8	2.988	0.611
WT-S5-231018-490	Air Dry #1	11.75	22.88	1.95	18.34	1.554	64.3	2.778	0.559
WT-S5-231018-490	Air Dry #2	9.08	17.73	1.95	18.66	1.432	69.8	2.582	0.555
WT-S5-231018-490	Air Dry #3	6.74	13.17	1.95	18.80	0.922	108.5	2.023	0.456
WT-S5-231018-490	Air Dry #4	6.22	12.16	1.95	18.89	0.785	127.4	1.756	0.447
WT-S5-231018-490	Air Dry #5	5.72	11.18	1.95	18.83	0.658	152.1	2.106	0.312
WT-S5-231018-490	Oven Dry	0	0	1.95	21.76	0.282	355.0	0.921	0.306

¹ Adjusted for volume changes during testing, if applicable.

Initial Properties



**Summary of Initial Moisture Content, Dry Bulk Density
Wet Bulk Density and Calculated Porosity**

Sample Number	Moisture Content				Dry Bulk Density (g/cm ³)	Wet Bulk Density (g/cm ³)	Calculated Porosity (%)
	As Received		Remolded				
	Gravimetric (%) (%, g/g)	Volumetric (%) (%, cm ³ /cm ³)	Gravimetric (%) (%, g/g)	Volumetric (%) (%, cm ³ /cm ³)			
W-S9-161018-435	4.4	NA	---	---	NA	NA	NA
WT-S9-171018-459	24.5	NA	25.6	37.5	1.46	1.84	44.7
WT-S9-171018-464	3.8	NA	4.1	7.4	1.83	1.90	31.0
WT-S5-231018-490	13.1	NA	13.1	24.9	1.90	2.14	28.5

NA = Not analyzed

--- = This sample was not remolded



**Data for Initial Moisture Content,
Bulk Density, Porosity, and Percent Saturation**

Job Name: EA Engineering, Science, and Technology, Inc.
Job Number: DB18.1333.00
Sample Number: W-S9-161018-435
Project Name: 62735DM02 KAFB-Bulk Fuels Fac Vadose Zone
PO Number: 18189

	<u>As Received</u>	<u>Remolded</u>
<i>Test Date:</i>	22-Oct-18	---
<i>Field weight* of sample (g):</i>	1727.93	
<i>Tare weight, ring (g):</i>	0.00	
<i>Tare weight, pan/plate (g):</i>	298.95	
<i>Tare weight, other (g):</i>	0.00	
<i>Dry weight of sample (g):</i>	1369.16	
<i>Sample volume (cm³):</i>	NA	
<i>Assumed particle density (g/cm³):</i>	2.65	
<i>Gravimetric Moisture Content (% g/g):</i>	4.4	
<i>Volumetric Moisture Content (% vol):</i>	NA	
<i>Dry bulk density (g/cm³):</i>	NA	
<i>Wet bulk density (g/cm³):</i>	NA	
<i>Calculated Porosity (% vol):</i>	NA	
<i>Percent Saturation:</i>	NA	

Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines

Comments:

* Weight including tares
NA = Not analyzed
--- = This sample was not remolded



**Data for Initial Moisture Content,
Bulk Density, Porosity, and Percent Saturation**

Job Name: EA Engineering, Science, and Technology, Inc.
Job Number: DB18.1333.00
Sample Number: WT-S9-171018-459
Project Name: 62735DM02 KAFB-Bulk Fuels Fac Vadose Zone
PO Number: 18189

	<u>As Received</u>	<u>Remolded</u>
<i>Test Date:</i>	22-Oct-18	22-Oct-18
<i>Field weight* of sample (g):</i>	1151.09	547.46
<i>Tare weight, ring (g):</i>	0.00	137.44
<i>Tare weight, pan/plate (g):</i>	298.96	0.00
<i>Tare weight, other (g):</i>	0.00	0.00
<i>Dry weight of sample (g):</i>	684.30	326.44
<i>Sample volume (cm³):</i>	NA	222.96
<i>Assumed particle density (g/cm³):</i>	2.65	2.65
<i>Gravimetric Moisture Content (% g/g):</i>	24.5	25.6
<i>Volumetric Moisture Content (% vol):</i>	NA	37.5
<i>Dry bulk density (g/cm³):</i>	NA	1.46
<i>Wet bulk density (g/cm³):</i>	NA	1.84
<i>Calculated Porosity (% vol):</i>	NA	44.7
<i>Percent Saturation:</i>	NA	83.8

Laboratory analysis by: D. O'Dowd D. O'Dowd
Data entered by: C. Krous C. Krous
Checked by: J. Hines J. Hines

Comments:

* Weight including tares
NA = Not analyzed
--- = This sample was not remolded



**Data for Initial Moisture Content,
Bulk Density, Porosity, and Percent Saturation**

Job Name: EA Engineering, Science, and Technology, Inc.
Job Number: DB18.1333.00
Sample Number: WT-S9-171018-464
Project Name: 62735DM02 KAFB-Bulk Fuels Fac Vadose Zone
PO Number: 18189

	<u>As Received</u>	<u>Remolded</u>
<i>Test Date:</i>	22-Oct-18	22-Oct-18
<i>Field weight* of sample (g):</i>	1600.01	2452.33
<i>Tare weight, ring (g):</i>	0.00	630.94
<i>Tare weight, pan/plate (g):</i>	268.45	0.00
<i>Tare weight, other (g):</i>	0.00	0.00
<i>Dry weight of sample (g):</i>	1282.75	1750.11
<i>Sample volume (cm³):</i>	NA	957.70
<i>Assumed particle density (g/cm³):</i>	2.65	2.65
<i>Gravimetric Moisture Content (% g/g):</i>	3.8	4.1
<i>Volumetric Moisture Content (% vol):</i>	NA	7.4
<i>Dry bulk density (g/cm³):</i>	NA	1.83
<i>Wet bulk density (g/cm³):</i>	NA	1.90
<i>Calculated Porosity (% vol):</i>	NA	31.0
<i>Percent Saturation:</i>	NA	24.0

<i>Laboratory analysis by:</i> D. O'Dowd	D. O'Dowd
<i>Data entered by:</i> C. Krous	C. Krous
<i>Checked by:</i> J. Hines	J. Hines

Comments:

* Weight including tares
NA = Not analyzed
--- = This sample was not remolded



**Data for Initial Moisture Content,
Bulk Density, Porosity, and Percent Saturation**

Job Name: EA Engineering, Science, and Technology, Inc.
Job Number: DB18.1333.00
Sample Number: WT-S5-231018-490
Project Name: 62735DM02 KAFB-Bulk Fuels Fac Vadose Zone
PO Number: 18189

	<u>As Received</u>	<u>Remolded</u>
<i>Test Date:</i>	25-Oct-18	25-Oct-18
<i>Field weight* of sample (g):</i>	2671.30	2671.30
<i>Tare weight, ring (g):</i>	628.78	628.78
<i>Tare weight, pan/plate (g):</i>	0.00	0.00
<i>Tare weight, other (g):</i>	0.00	0.00
<i>Dry weight of sample (g):</i>	1805.54	1805.54
<i>Sample volume (cm³):</i>	NA	952.69
<i>Assumed particle density (g/cm³):</i>	2.65	2.65
<i>Gravimetric Moisture Content (% g/g):</i>	13.1	13.1
<i>Volumetric Moisture Content (% vol):</i>	NA	24.9
<i>Dry bulk density (g/cm³):</i>	NA	1.90
<i>Wet bulk density (g/cm³):</i>	NA	2.14
<i>Calculated Porosity (% vol):</i>	NA	28.5
<i>Percent Saturation:</i>	NA	87.3

<i>Laboratory analysis by:</i> D. O'Dowd	D. O'Dowd
<i>Data entered by:</i> C. Krous	C. Krous
<i>Checked by:</i> J. Hines	J. Hines

Comments:

* Weight including tares
NA = Not analyzed
--- = This sample was not remolded

Thermal Properties



Summary of Thermal Properties

Sample	Reading	Gravimetric Moisture Content (g/g, %)	Volumetric Moisture Content ¹ (vol/vol, %)	Dry Bulk Density ¹ (g/cm ³)	Temp °C	K W/(m·K)	ρ °C·cm/W	C MJ/(m ³ ·K)	D mm ² /s
WT-S9-171018-459	Initial	25.60	37.49	1.46	22.55	1.429	70.0	3.952	0.362
WT-S9-171018-459	0	31.82	45.30	1.42	21.52	1.429	70.0	3.805	0.376
WT-S9-171018-459	Air Dry #1	27.44	39.45	1.44	18.86	1.370	73.0	3.373	0.406
WT-S9-171018-459	Air Dry #2	23.24	33.93	1.46	18.24	1.360	73.5	3.241	0.420
WT-S9-171018-459	Air Dry #3	18.58	29.18	1.57	19.07	1.293	77.3	3.166	0.408
WT-S9-171018-459	Air Dry #4	13.08	21.03	1.61	19.84	1.060	94.3	2.670	0.397
WT-S9-171018-459	Air Dry #5	11.67	18.78	1.61	20.24	0.996	100.4	2.607	0.382
WT-S9-171018-459	Air Dry #6	10.45	16.82	1.61	20.60	0.941	106.2	2.502	0.376
WT-S9-171018-459	Oven Dry	0	0	1.61	21.22	0.547	182.7	1.554	0.352
WT-S9-171018-464	Initial	4.07	7.44	1.83	22.20	0.681	146.8	1.559	0.437
WT-S9-171018-464	0	16.45	30.06	1.83	22.59	1.686	59.3	3.038	0.555
WT-S9-171018-464	Air Dry #1	14.53	27.13	1.87	21.31	1.676	59.7	2.631	0.637
WT-S9-171018-464	Air Dry #2	12.19	23.21	1.90	20.15	1.587	63.0	2.096	0.757
WT-S9-171018-464	Air Dry #3	8.23	16.09	1.95	17.87	1.167	85.7	1.793	0.651
WT-S9-171018-464	Air Dry #4	6.18	12.07	1.95	18.79	1.001	99.9	2.082	0.481
WT-S9-171018-464	Air Dry #5	4.57	8.94	1.95	18.76	0.810	123.5	1.619	0.500
WT-S9-171018-464	Air Dry #6	3.76	7.35	1.95	20.63	0.656	152.3	1.886	0.348
WT-S9-171018-464	Oven Dry	0	0	1.95	21.47	0.262	381.1	1.471	0.178

¹ Adjusted for volume changes during testing, if applicable.



Summary of Thermal Properties (Continued)

Sample	Reading	Gravimetric Moisture Content (g/g, %)	Volumetric Moisture Content ¹ (vol/vol, %)	Dry Bulk Density ¹ (g/cm ³)	Temp °C	K W/(m·K)	ρ °C·cm/W	C MJ/(m ³ ·K)	D mm ² /s
WT-S5-231018-490	Initial	13.13	24.87	1.90	21.72	1.907	52.4	4.024	0.474
WT-S5-231018-490	0	14.44	27.36	1.90	20.78	1.825	54.8	2.988	0.611
WT-S5-231018-490	Air Dry #1	11.75	22.88	1.95	18.34	1.554	64.3	2.778	0.559
WT-S5-231018-490	Air Dry #2	9.08	17.73	1.95	18.66	1.432	69.8	2.582	0.555
WT-S5-231018-490	Air Dry #3	6.74	13.17	1.95	18.80	0.922	108.5	2.023	0.456
WT-S5-231018-490	Air Dry #4	6.22	12.16	1.95	18.89	0.785	127.4	1.756	0.447
WT-S5-231018-490	Air Dry #5	5.72	11.18	1.95	18.83	0.658	152.1	2.106	0.312
WT-S5-231018-490	Oven Dry	0	0	1.95	21.76	0.282	355.0	0.921	0.306

¹ Adjusted for volume changes during testing, if applicable.



Daniel B. Stephens & Associates, Inc.

Thermal Properties Results Sheet for Sample: WT-S9-171018-464

Job Name: EA Engineering, Inc.

Job Number: DB18.1333.00

Sample Number: WT-S9-171018-464

Project Name: 62735DM02 KAFB-Bulk Fuels Fac Vadose Zone

PO Number: 18189

Instrument Description: Decagon KD2 Pro

Probe: ☐ KS-1, 6 cm length, 1.3 mm diameter, single needle

☐ TR-1, 10 cm length, 2.4 mm diameter, single needle

☒ SH-1, 3 cm length, 1.3 mm diameter, dual needle, 6 mm spacing

Test Start Date: 10/22/18

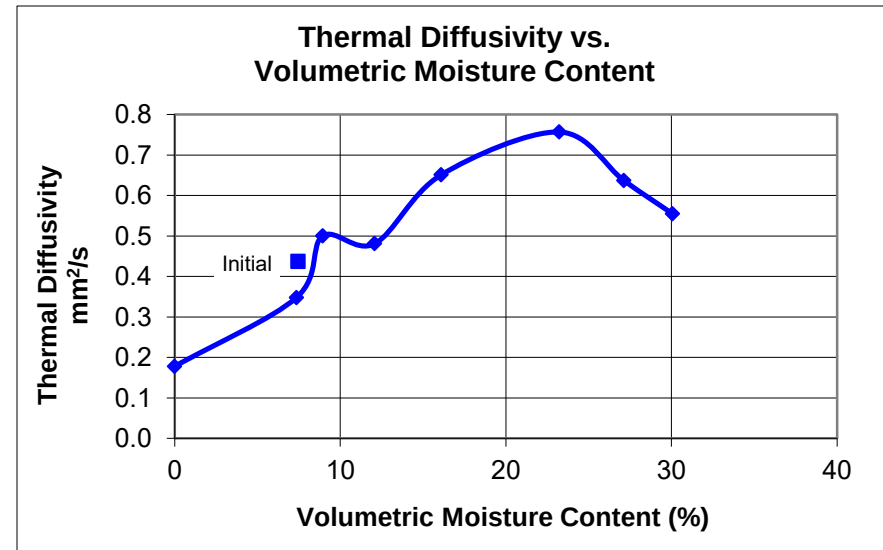
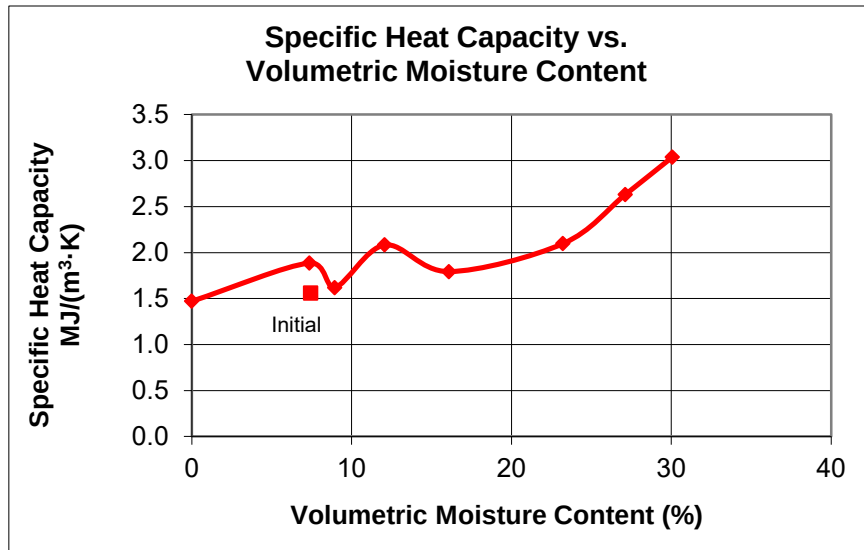
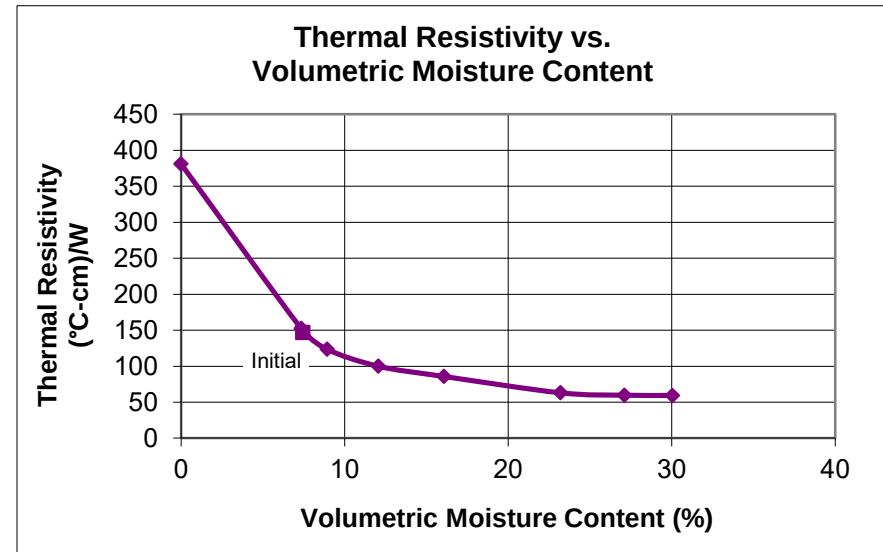
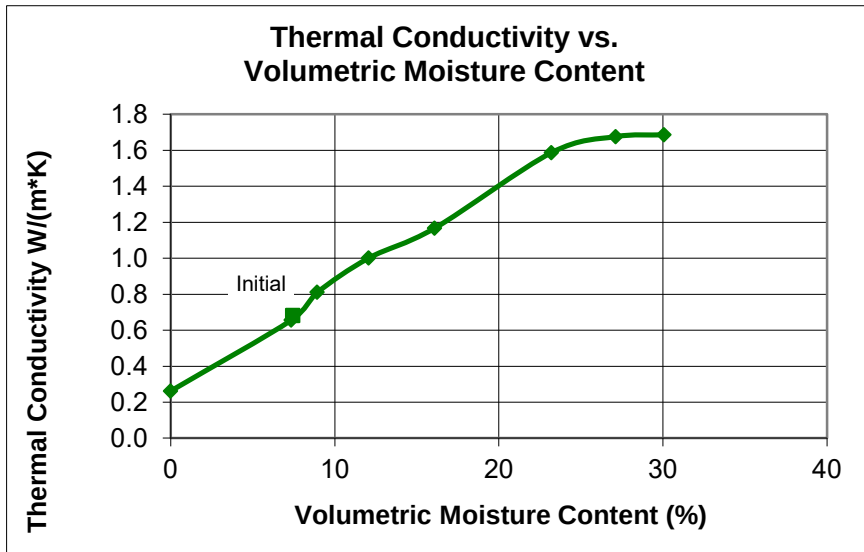
Reading	Water Potential (-cm water)	Gravimetric Moisture Content (g/g, %)	Volumetric Moisture Content ¹ (vol/vol, %)	Dry Bulk Density ¹ (g/cm ³)	Test Temperature (°C)	K Thermal Conductivity W/(m·K)	ρ Thermal Resistivity °C·cm/W	C Specific Heat Capacity MJ/(m ³ ·K)	D Thermal Diffusivity (mm ² /s)
Initial	---	4.07	7.44	1.83	22.20	0.681	146.8	1.559	0.437
Saturated	0	16.45	30.06	1.83	22.59	1.686	59.3	3.038	0.555
Air Dry #1	---	14.53	27.13	1.87	21.31	1.676	59.7	2.631	0.637
Air Dry #2	---	12.19	23.21	1.90	20.15	1.587	63.0	2.096	0.757
Air Dry #3	---	8.23	16.09	1.95	17.87	1.167	85.7	1.793	0.651
Air Dry #4	---	6.18	12.07	1.95	18.79	1.001	99.9	2.082	0.481
Air Dry #5	---	4.57	8.94	1.95	18.76	0.810	123.5	1.619	0.500
Air Dry #6	---	3.76	7.35	1.95	20.63	0.656	152.3	1.886	0.348
Oven Dry	---	0	0	1.95	21.47	0.262	381.1	1.471	0.178

--- = Value not measured.

¹ Adjusted for volume changes during testing, if applicable.



Thermal Properties Results Sheet for Sample: WT-S9-171018-464
Scatter Plots





Thermal Properties Data

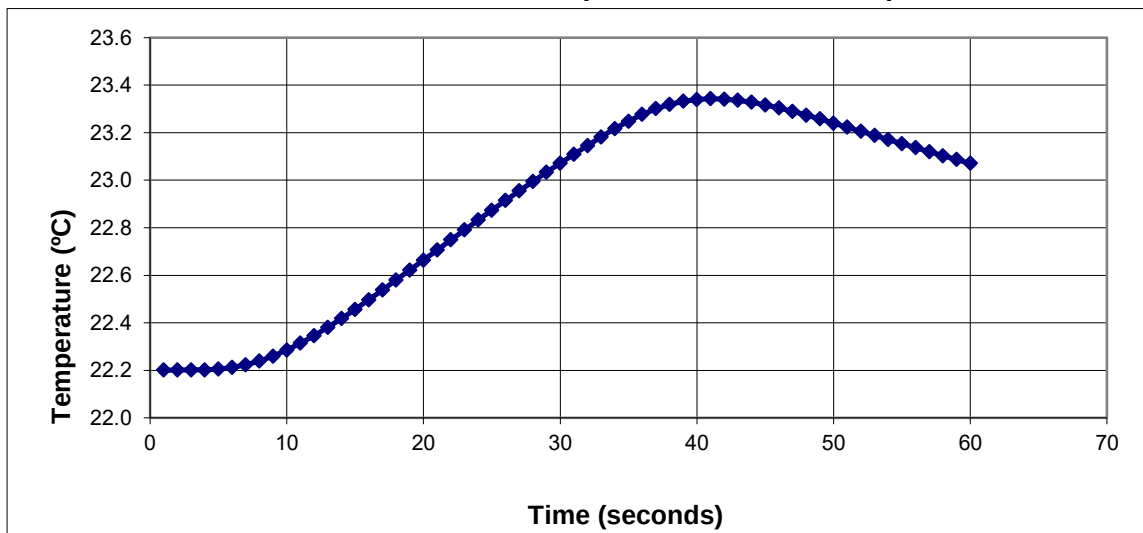
Sample Number: WT-S9-171018-464
Potential (-cm water): Initial

Test Date/Time: 10/22/2018 2:26 PM	K (W/(m·K)): 0.681
Sensor: SH-1	ρ (°C·cm/W): 146.8
Test Temp.(°C): 22.2	C (MJ/(m ³ ·K)): 1.559
KD2 Pro Sample ID: 464-AR	D (mm ² /s): 0.437
Power (W/m): 21.950	Err: 0.0032
Current (amps): 0.145	

Raw Data

Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)
1 22.201	16 22.497	31 23.109	46 23.304
2 22.201	17 22.538	32 23.146	47 23.290
3 22.202	18 22.580	33 23.182	48 23.274
4 22.202	19 22.622	34 23.217	49 23.258
5 22.205	20 22.664	35 23.248	50 23.240
6 22.212	21 22.706	36 23.277	51 23.223
7 22.223	22 22.749	37 23.301	52 23.206
8 22.239	23 22.791	38 23.319	53 23.189
9 22.260	24 22.833	39 23.332	54 23.171
10 22.285	25 22.874	40 23.340	55 23.154
11 22.314	26 22.915	41 23.343	56 23.137
12 22.346	27 22.955	42 23.341	57 23.120
13 22.381	28 22.994	43 23.336	58 23.103
14 22.418	29 23.034	44 23.328	59 23.087
15 22.456	30 23.072	45 23.317	60 23.071

WT-S9-171018-464,Potential: Initial - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

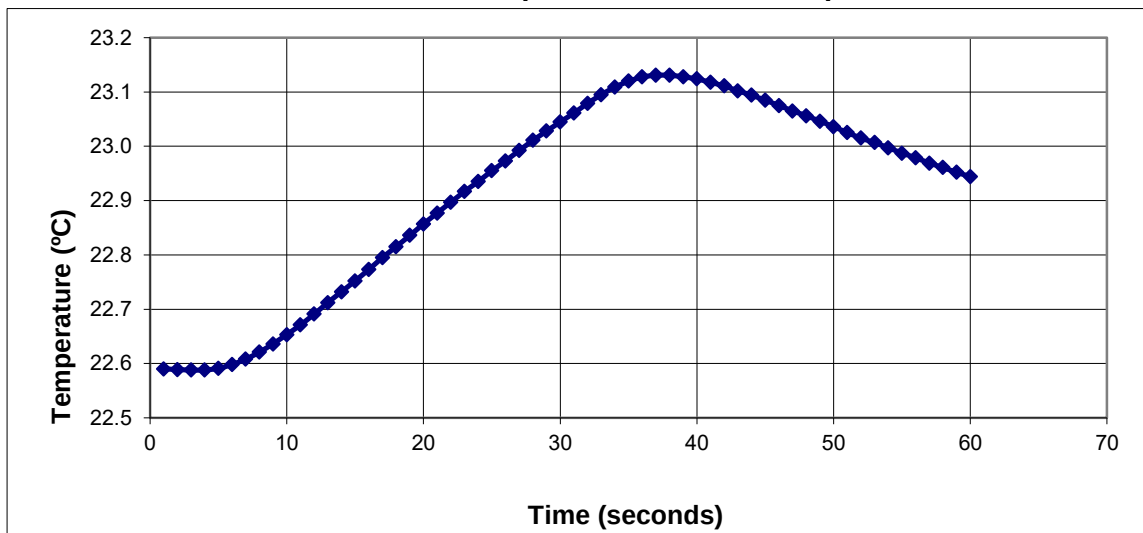
Sample Number: WT-S9-171018-464
Potential (-cm water): 0

Test Date/Time:	10/23/2018 9:50 AM	K (W/(m·K)):	1.686
Sensor:	SH-1	ρ (°C·cm/W):	59.3
Test Temp.(°C):	22.6	C (MJ/(m ³ ·K)):	3.038
KD2 Pro Sample ID:	464-SA	D (mm ² /s):	0.555
Power (W/m):	21.190	Err:	0.0012
Current (amps):	0.143		

Raw Data

Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)	
1	22.590	16	22.773	31	23.061	46	23.075
2	22.589	17	22.795	32	23.079	47	23.065
3	22.588	18	22.815	33	23.095	48	23.056
4	22.588	19	22.836	34	23.109	49	23.046
5	22.591	20	22.857	35	23.120	50	23.036
6	22.598	21	22.877	36	23.128	51	23.025
7	22.608	22	22.897	37	23.131	52	23.015
8	22.621	23	22.917	38	23.131	53	23.007
9	22.636	24	22.935	39	23.128	54	22.997
10	22.653	25	22.955	40	23.124	55	22.987
11	22.671	26	22.973	41	23.118	56	22.979
12	22.691	27	22.992	42	23.111	57	22.969
13	22.712	28	23.011	43	23.102	58	22.961
14	22.732	29	23.028	44	23.094	59	22.952
15	22.752	30	23.045	45	23.085	60	22.944

WT-S9-171018-464, Potential: 0 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

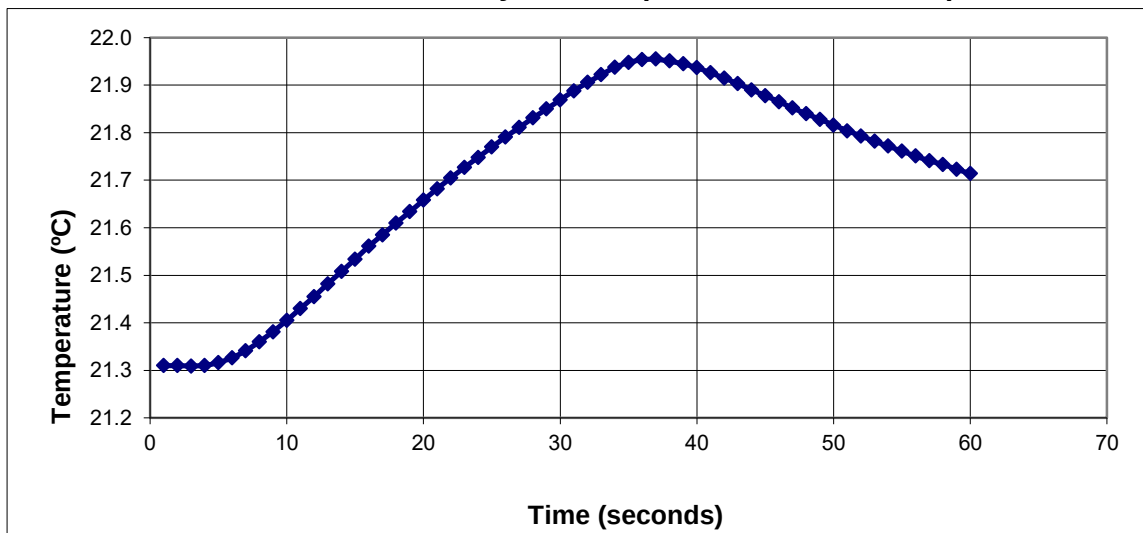
Sample Number: WT-S9-171018-464
Potential (-cm water): Air Dry #1

Test Date/Time: 10/24/2018 8:04 AM	K (W/(m·K)): 1.676
Sensor: SH-1	ρ (°C·cm/W): 59.7
Test Temp.(°C): 21.3	C (MJ/(m ³ ·K)): 2.631
KD2 Pro Sample ID: 464-AD1	D (mm ² /s): 0.637
Power (W/m): 21.690	Err: 0.0033
Current (amps): 0.144	

Raw Data

Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)	
1	21.310	16	21.561	31	21.888	46	21.865
2	21.310	17	21.585	32	21.906	47	21.852
3	21.309	18	21.610	33	21.922	48	21.840
4	21.310	19	21.634	34	21.938	49	21.828
5	21.316	20	21.658	35	21.948	50	21.816
6	21.326	21	21.682	36	21.954	51	21.804
7	21.341	22	21.705	37	21.955	52	21.793
8	21.360	23	21.727	38	21.951	53	21.782
9	21.381	24	21.748	39	21.945	54	21.772
10	21.405	25	21.770	40	21.937	55	21.761
11	21.430	26	21.791	41	21.926	56	21.751
12	21.455	27	21.811	42	21.915	57	21.741
13	21.482	28	21.831	43	21.903	58	21.733
14	21.508	29	21.850	44	21.890	59	21.723
15	21.534	30	21.869	45	21.878	60	21.714

WT-S9-171018-464,Potential: Air Dry #1 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

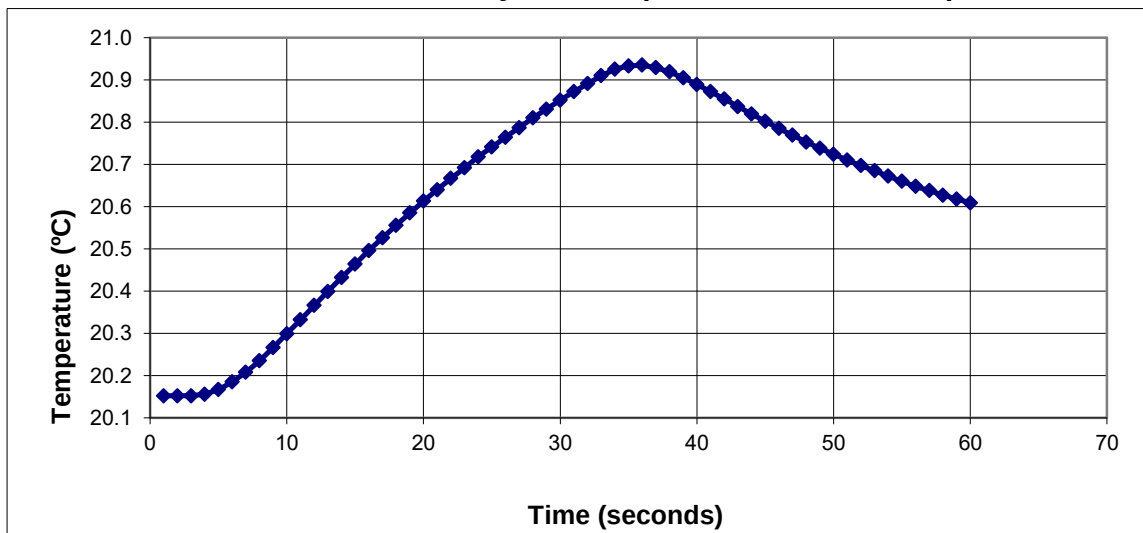
Sample Number: WT-S9-171018-464
Potential (-cm water): Air Dry #2

Test Date/Time:	10/25/2018 8:31 AM	K (W/(m·K)):	1.587
Sensor:	SH-1	ρ (°C·cm/W):	63.0
Test Temp.(°C):	20.2	C (MJ/(m ³ ·K)):	2.096
KD2 Pro Sample ID:	464-AD2	D (mm ² /s):	0.757
Power (W/m):	21.330	Err:	0.0058
Current (amps):	0.143		

Raw Data

Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)	
1	20.152	16	20.496	31	20.872	46	20.785
2	20.152	17	20.526	32	20.891	47	20.769
3	20.152	18	20.556	33	20.910	48	20.753
4	20.156	19	20.585	34	20.925	49	20.738
5	20.167	20	20.613	35	20.933	50	20.724
6	20.185	21	20.640	36	20.935	51	20.710
7	20.208	22	20.667	37	20.929	52	20.697
8	20.235	23	20.692	38	20.919	53	20.685
9	20.266	24	20.718	39	20.905	54	20.672
10	20.299	25	20.741	40	20.889	55	20.660
11	20.332	26	20.764	41	20.872	56	20.648
12	20.366	27	20.787	42	20.855	57	20.638
13	20.399	28	20.810	43	20.837	58	20.627
14	20.432	29	20.831	44	20.819	59	20.618
15	20.464	30	20.852	45	20.802	60	20.609

WT-S9-171018-464,Potential: Air Dry #2 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

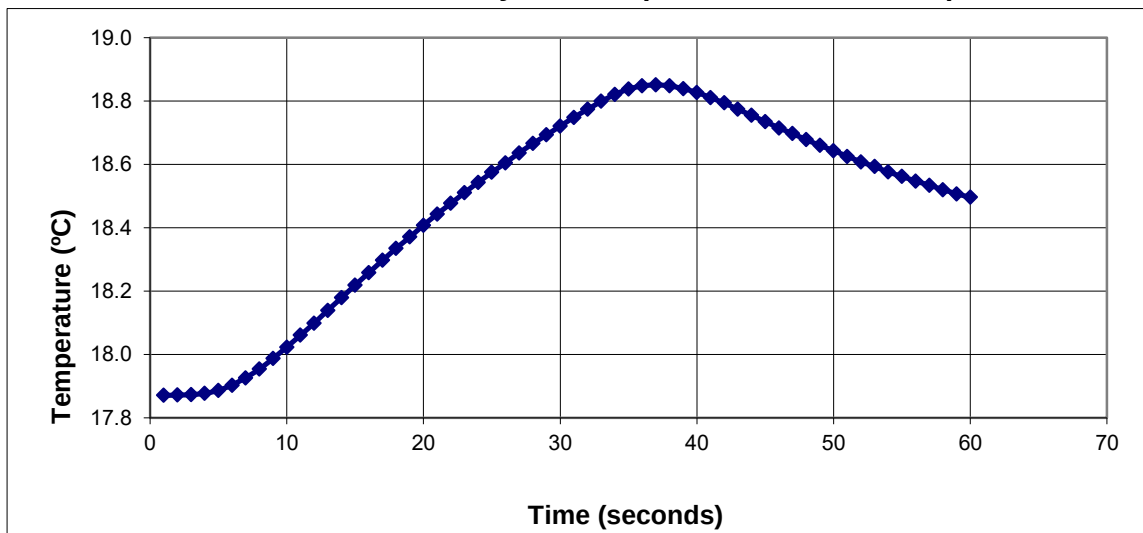
Sample Number: WT-S9-171018-464
Potential (-cm water): Air Dry #3

Test Date/Time: 10/26/2018 2:47 PM	K (W/(m·K)): 1.167
Sensor: SH-1	ρ (°C·cm/W): 85.7
Test Temp.(°C): 17.9	C (MJ/(m ³ ·K)): 1.793
KD2 Pro Sample ID: 464-AD3	D (mm ² /s): 0.651
Power (W/m): 22.190	Err: 0.0043
Current (amps): 0.146	

Raw Data

Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)
1 17.871	16 18.258	31 18.748	46 18.715
2 17.872	17 18.297	32 18.774	47 18.697
3 17.873	18 18.335	33 18.799	48 18.678
4 17.877	19 18.371	34 18.821	49 18.660
5 17.886	20 18.408	35 18.838	50 18.643
6 17.903	21 18.443	36 18.848	51 18.625
7 17.926	22 18.477	37 18.851	52 18.608
8 17.954	23 18.511	38 18.848	53 18.593
9 17.987	24 18.543	39 18.839	54 18.576
10 18.023	25 18.575	40 18.827	55 18.562
11 18.061	26 18.605	41 18.811	56 18.547
12 18.099	27 18.636	42 18.794	57 18.534
13 18.139	28 18.666	43 18.774	58 18.520
14 18.179	29 18.693	44 18.755	59 18.507
15 18.219	30 18.721	45 18.735	60 18.496

WT-S9-171018-464,Potential: Air Dry #3 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

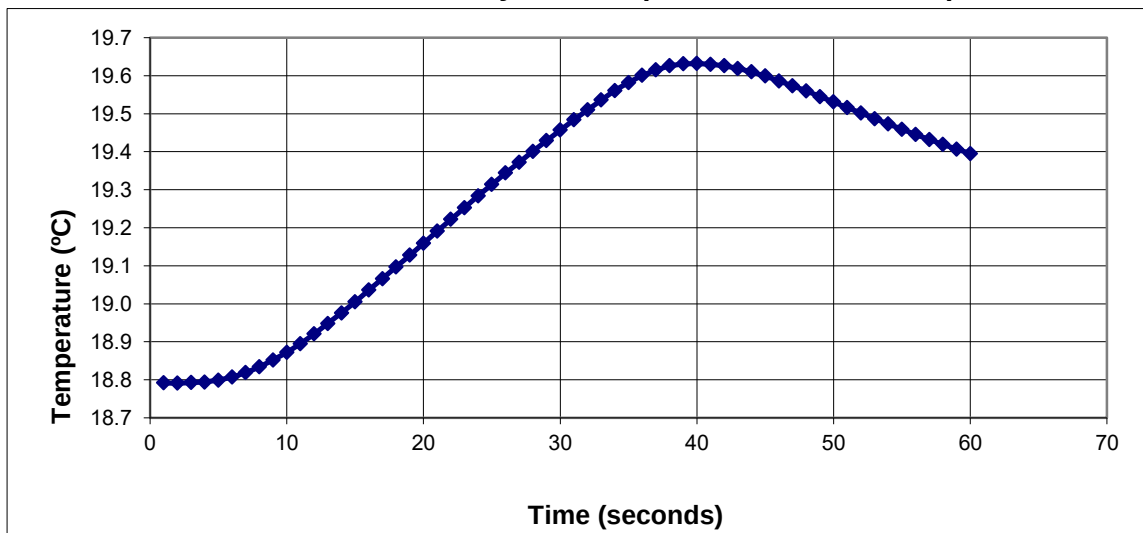
Sample Number: WT-S9-171018-464
Potential (-cm water): Air Dry #4

Test Date/Time: 10/29/2018 2:57 PM	K (W/(m·K)): 1.001
Sensor: SH-1	ρ (°C·cm/W): 99.9
Test Temp.(°C): 18.8	C (MJ/(m ³ ·K)): 2.082
KD2 Pro Sample ID: 464-AD4	D (mm ² /s): 0.481
Power (W/m): 21.780	Err: 0.0044
Current (amps): 0.145	

Raw Data

Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)
1 18.792	16 19.036	31 19.484	46 19.586
2 18.791	17 19.066	32 19.510	47 19.573
3 18.793	18 19.097	33 19.536	48 19.560
4 18.794	19 19.128	34 19.561	49 19.545
5 18.799	20 19.159	35 19.582	50 19.531
6 18.807	21 19.191	36 19.601	51 19.516
7 18.819	22 19.222	37 19.615	52 19.502
8 18.834	23 19.253	38 19.626	53 19.487
9 18.852	24 19.284	39 19.631	54 19.473
10 18.872	25 19.314	40 19.632	55 19.459
11 18.895	26 19.344	41 19.630	56 19.445
12 18.921	27 19.372	42 19.626	57 19.432
13 18.948	28 19.401	43 19.619	58 19.419
14 18.976	29 19.429	44 19.610	59 19.407
15 19.005	30 19.457	45 19.599	60 19.395

WT-S9-171018-464,Potential: Air Dry #4 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

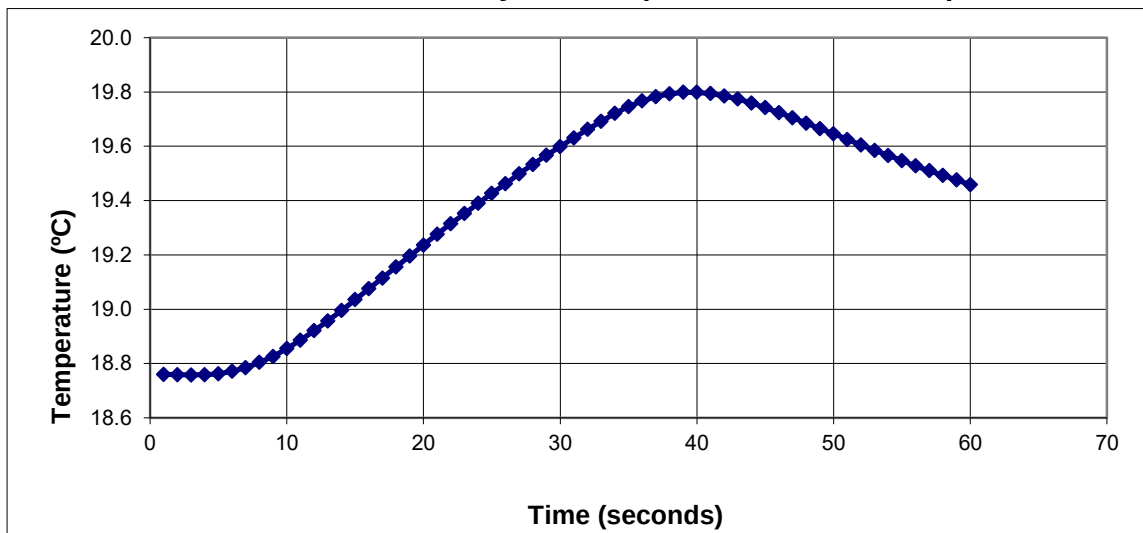
Sample Number: WT-S9-171018-464
Potential (-cm water): Air Dry #5

Test Date/Time: 10/30/2018 8:30 AM	K (W/(m·K)): 0.810
Sensor: SH-1	ρ (°C·cm/W): 123.5
Test Temp.(°C): 18.8	C (MJ/(m ³ ·K)): 1.619
KD2 Pro Sample ID: 464-AD5	D (mm ² /s): 0.500
Power (W/m): 21.940	Err: 0.0041
Current (amps): 0.145	

Raw Data

Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)
1 18.760	16 19.075	31 19.631	46 19.724
2 18.759	17 19.115	32 19.662	47 19.705
3 18.758	18 19.156	33 19.692	48 19.685
4 18.759	19 19.196	34 19.721	49 19.665
5 18.762	20 19.236	35 19.746	50 19.646
6 18.772	21 19.276	36 19.767	51 19.625
7 18.784	22 19.315	37 19.783	52 19.605
8 18.804	23 19.352	38 19.793	53 19.585
9 18.826	24 19.390	39 19.799	54 19.566
10 18.855	25 19.427	40 19.799	55 19.547
11 18.886	26 19.462	41 19.794	56 19.528
12 18.921	27 19.499	42 19.785	57 19.510
13 18.957	28 19.533	43 19.774	58 19.493
14 18.995	29 19.567	44 19.759	59 19.476
15 19.035	30 19.599	45 19.742	60 19.459

WT-S9-171018-464,Potential: Air Dry #5 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

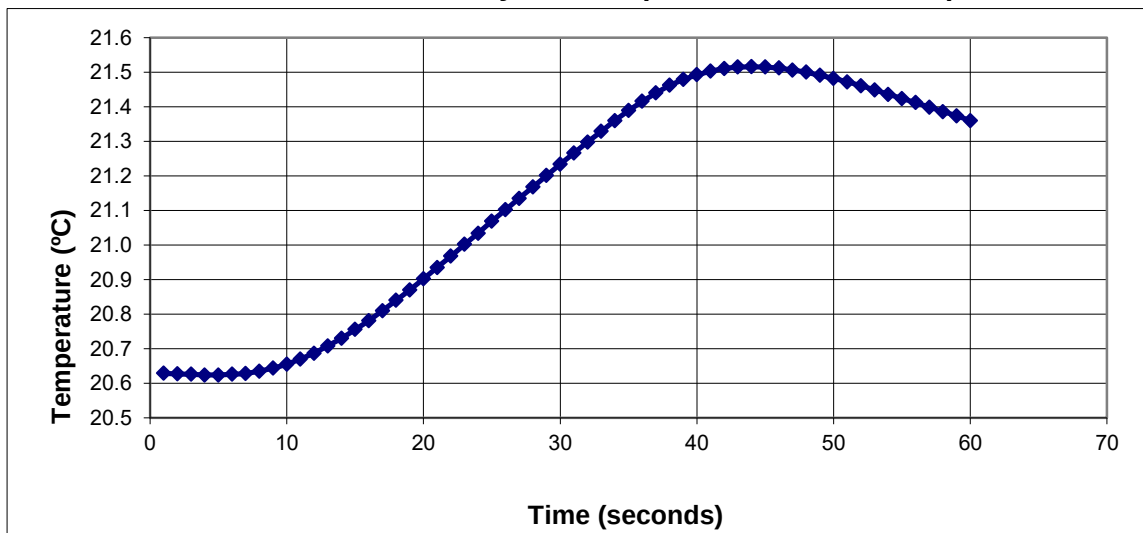
Sample Number: WT-S9-171018-464
Potential (-cm water): Air Dry #6

Test Date/Time:	10/31/2018 9:25 AM	K (W/(m·K)):	0.656
Sensor:	SH-1	ρ (°C·cm/W):	152.3
Test Temp.(°C):	20.6	C (MJ/(m ³ ·K)):	1.886
KD2 Pro Sample ID:	464-AD6	D (mm ² /s):	0.348
Power (W/m):	21.580	Err:	0.0023
Current (amps):	0.144		

Raw Data

Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)	
1	20.629	16	20.781	31	21.266	46	21.512
2	20.627	17	20.810	32	21.298	47	21.506
3	20.626	18	20.840	33	21.329	48	21.500
4	20.624	19	20.870	34	21.360	49	21.491
5	20.624	20	20.902	35	21.389	50	21.482
6	20.626	21	20.935	36	21.416	51	21.472
7	20.628	22	20.968	37	21.440	52	21.461
8	20.635	23	21.002	38	21.462	53	21.449
9	20.644	24	21.034	39	21.479	54	21.436
10	20.655	25	21.069	40	21.493	55	21.424
11	20.670	26	21.102	41	21.503	56	21.412
12	20.687	27	21.135	42	21.511	57	21.399
13	20.708	28	21.168	43	21.515	58	21.386
14	20.730	29	21.201	44	21.516	59	21.374
15	20.756	30	21.234	45	21.515	60	21.360

WT-S9-171018-464,Potential: Air Dry #6 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

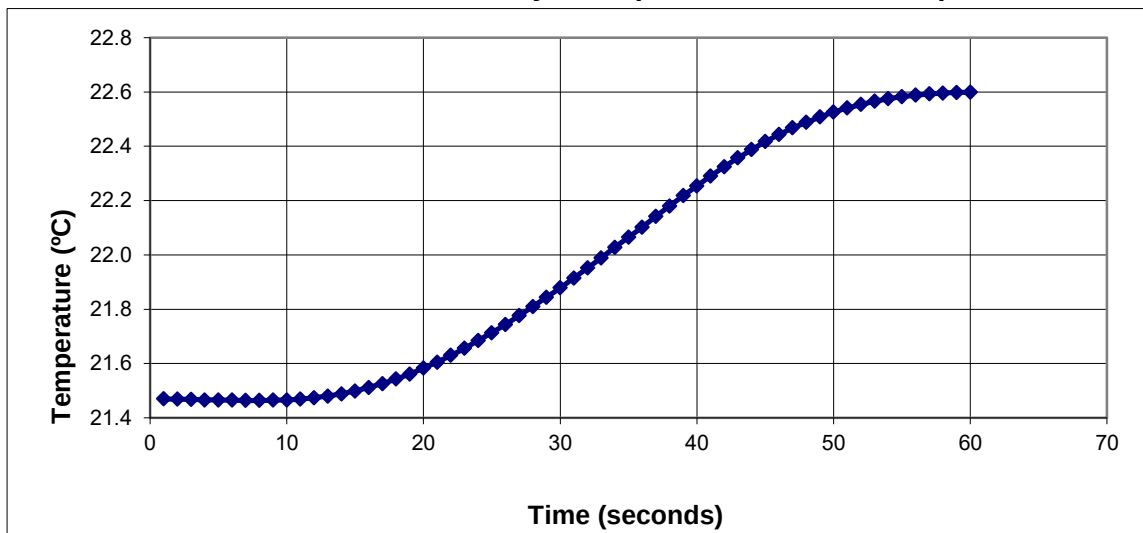
Sample Number: WT-S9-171018-464
Potential (-cm water): Oven Dry

Test Date/Time:	11/6/2018 3:50 PM	K (W/(m·K)):	0.262
Sensor:	SH-1	ρ (°C·cm/W):	381.1
Test Temp.(°C):	21.5	C (MJ/(m ³ ·K)):	1.471
KD2 Pro Sample ID:	464-OD	D (mm ² /s):	0.178
Power (W/m):	20.990	Err:	0.0009
Current (amps):	0.142		

Raw Data

Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)	
1	21.470	16	21.512	31	21.915	46	22.443
2	21.469	17	21.526	32	21.952	47	22.468
3	21.468	18	21.543	33	21.989	48	22.488
4	21.466	19	21.561	34	22.028	49	22.508
5	21.466	20	21.583	35	22.065	50	22.526
6	21.465	21	21.605	36	22.102	51	22.541
7	21.464	22	21.630	37	22.142	52	22.554
8	21.464	23	21.656	38	22.180	53	22.566
9	21.465	24	21.685	39	22.218	54	22.575
10	21.466	25	21.713	40	22.254	55	22.583
11	21.469	26	21.744	41	22.290	56	22.588
12	21.474	27	21.776	42	22.325	57	22.593
13	21.480	28	21.810	43	22.357	58	22.596
14	21.488	29	21.844	44	22.388	59	22.598
15	21.499	30	21.879	45	22.417	60	22.599

WT-S9-171018-464,Potential: Oven Dry - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Daniel B. Stephens & Associates, Inc.

Thermal Properties Results Sheet for Sample: WT-S9-171018-459

Job Name: EA Engineering, Inc.

Job Number: DB18.1333.00

Sample Number: WT-S9-171018-459

Project Name: 62735DM02 KAFB-Bulk Fuels Fac Vadose Zone

PO Number: 18189

Instrument Description: Decagon KD2 Pro

Probe: ☐ KS-1, 6 cm length, 1.3 mm diameter, single needle

☐ TR-1, 10 cm length, 2.4 mm diameter, single needle

☒ SH-1, 3 cm length, 1.3 mm diameter, dual needle, 6 mm spacing

Test Start Date: 10/22/18

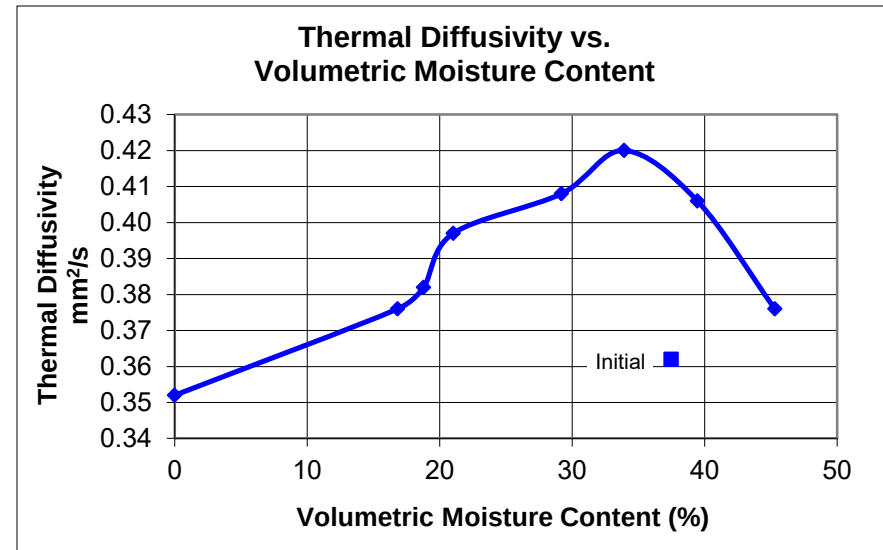
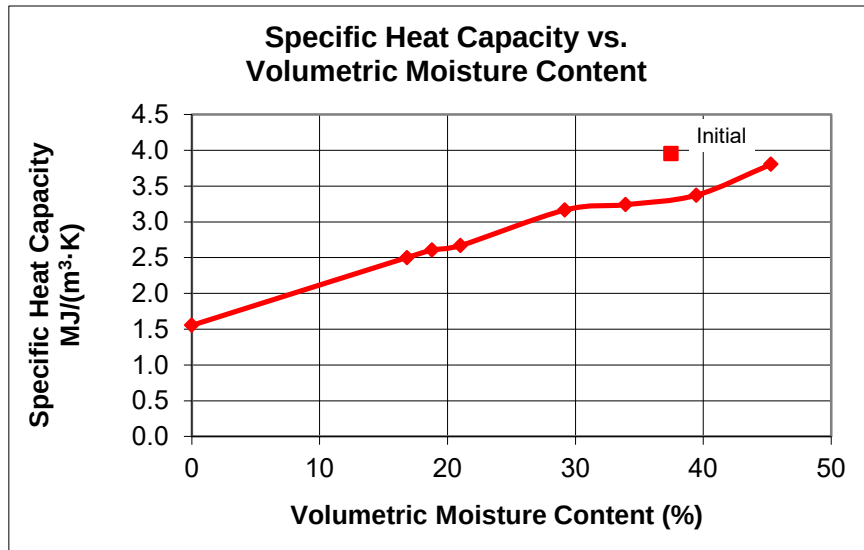
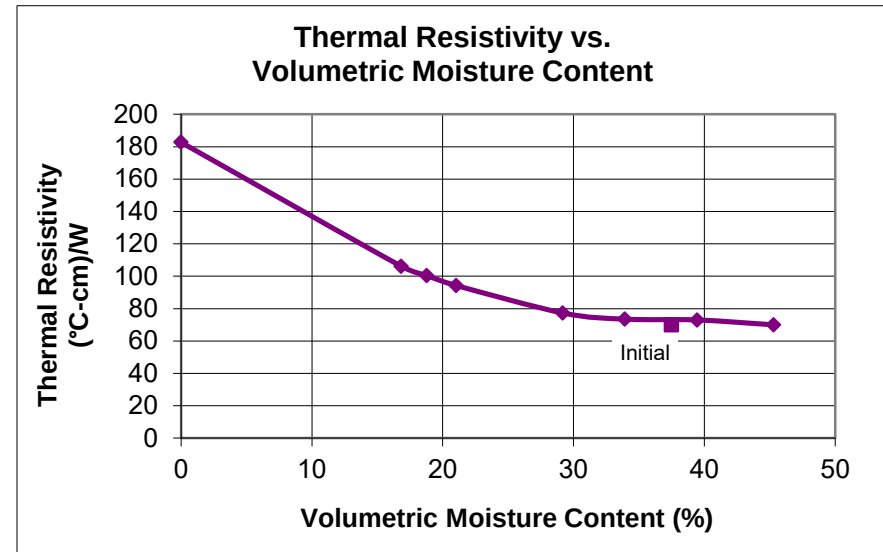
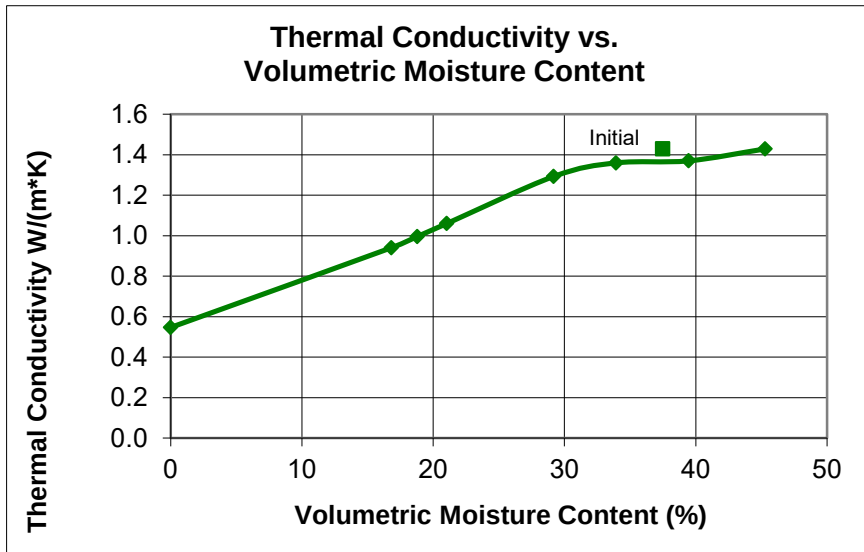
Reading	Water Potential (-cm water)	Gravimetric Moisture Content (g/g, %)	Volumetric Moisture Content ¹ (vol/vol, %)	Dry Bulk Density ¹ (g/cm ³)	Test Temperature (°C)	K Thermal Conductivity W/(m·K)	ρ Thermal Resistivity °C·cm/W	C Specific Heat Capacity MJ/(m ³ ·K)	D Thermal Diffusivity (mm ² /s)
Initial	---	25.60	37.49	1.46	22.55	1.429	70.0	3.952	0.362
Saturated	0	31.82	45.30	1.42	21.52	1.429	70.0	3.805	0.376
Air Dry #1	---	27.44	39.45	1.44	18.86	1.370	73.0	3.373	0.406
Air Dry #2	---	23.24	33.93	1.46	18.24	1.360	73.5	3.241	0.420
Air Dry #3	---	18.58	29.18	1.57	19.07	1.293	77.3	3.166	0.408
Air Dry #4	---	13.08	21.03	1.61	19.84	1.060	94.3	2.670	0.397
Air Dry #5	---	11.67	18.78	1.61	20.24	0.996	100.4	2.607	0.382
Air Dry #6	---	10.45	16.82	1.61	20.60	0.941	106.2	2.502	0.376
Oven Dry	---	0	0	1.61	21.22	0.547	182.7	1.554	0.352

--- = Value not measured.

¹ Adjusted for volume changes during testing, if applicable.



Thermal Properties Results Sheet for Sample: WT-S9-171018-459
Scatter Plots





Thermal Properties Data

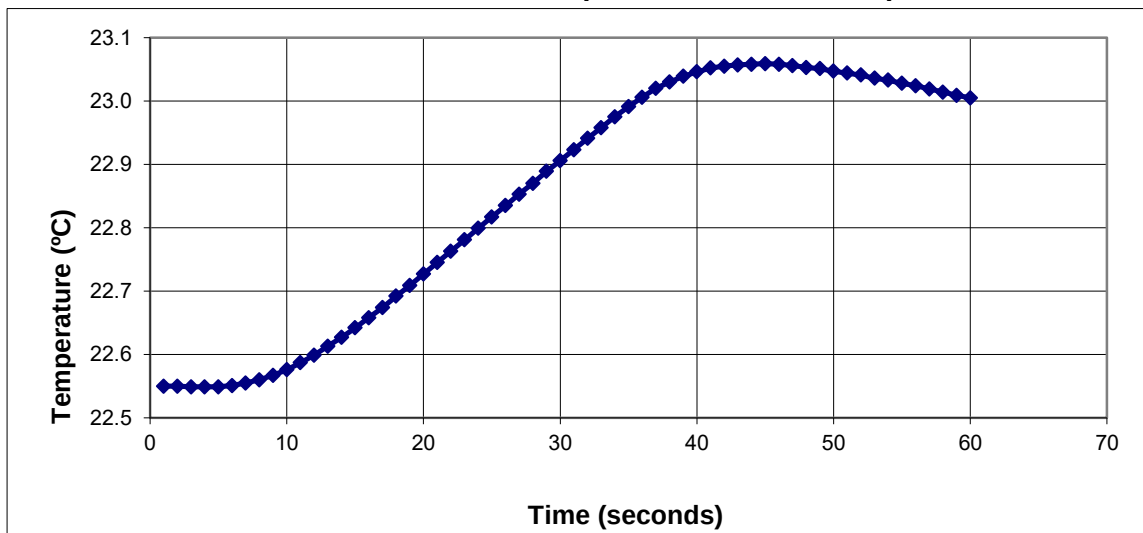
Sample Number: WT-S9-171018-459
Potential (-cm water): Initial

Test Date/Time: 10/22/2018 2:44 PM	K (W/(m·K)): 1.429
Sensor: SH-1	ρ (°C·cm/W): 70.0
Test Temp.(°C): 22.6	C (MJ/(m ³ ·K)): 3.952
KD2 Pro Sample ID: 459-AR	D (mm ² /s): 0.362
Power (W/m): 22.510	Err: 0.0023
Current (amps): 0.147	

Raw Data

Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)
1 22.550	16 22.658	31 22.923	46 23.058
2 22.550	17 22.674	32 22.941	47 23.056
3 22.549	18 22.692	33 22.958	48 23.053
4 22.549	19 22.709	34 22.975	49 23.051
5 22.549	20 22.727	35 22.991	50 23.047
6 22.551	21 22.745	36 23.006	51 23.044
7 22.555	22 22.763	37 23.020	52 23.041
8 22.560	23 22.781	38 23.030	53 23.036
9 22.567	24 22.799	39 23.039	54 23.033
10 22.576	25 22.817	40 23.046	55 23.028
11 22.587	26 22.835	41 23.052	56 23.024
12 22.599	27 22.853	42 23.055	57 23.019
13 22.613	28 22.870	43 23.057	58 23.014
14 22.627	29 22.889	44 23.058	59 23.009
15 22.642	30 22.906	45 23.059	60 23.005

WT-S9-171018-459,Potential: Initial - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

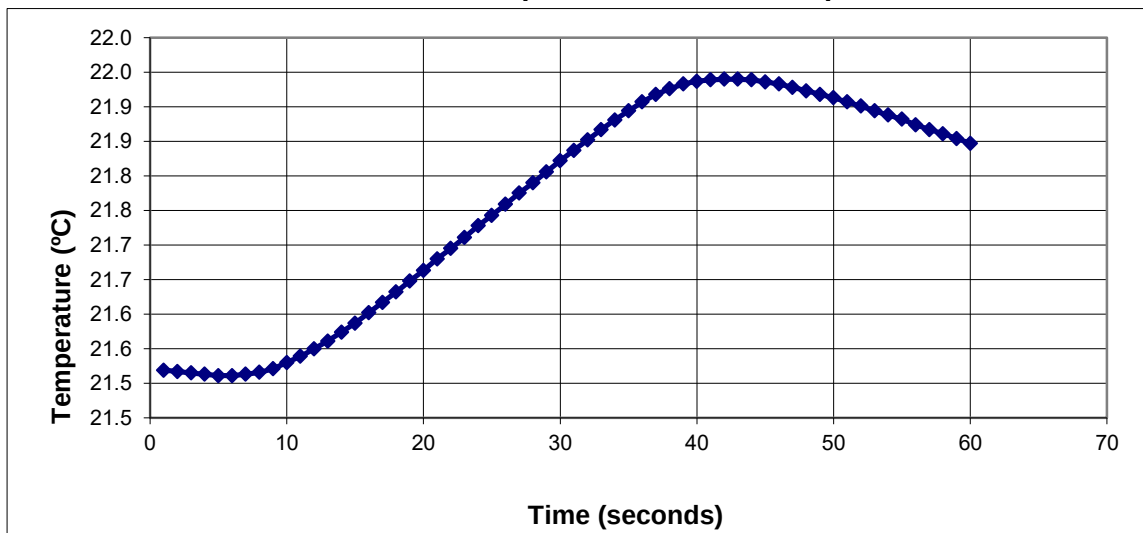
Sample Number: WT-S9-171018-459
Potential (-cm water): 0

Test Date/Time: 10/23/2018 10:14 AM	K (W/(m·K)): 1.429
Sensor: SH-1	ρ (°C·cm/W): 70.0
Test Temp.(°C): 21.5	C (MJ/(m³·K)): 3.805
KD2 Pro Sample ID: 459-SA	D (mm²/s): 0.376
Power (W/m): 21.470	Err: 0.0009
Current (amps): 0.144	

Raw Data

Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)	
1	21.519	16	21.602	31	21.837	46	21.933
2	21.517	17	21.617	32	21.852	47	21.928
3	21.515	18	21.632	33	21.867	48	21.923
4	21.513	19	21.648	34	21.881	49	21.918
5	21.511	20	21.663	35	21.894	50	21.913
6	21.511	21	21.680	36	21.907	51	21.907
7	21.513	22	21.695	37	21.918	52	21.901
8	21.516	23	21.711	38	21.926	53	21.894
9	21.521	24	21.728	39	21.933	54	21.888
10	21.530	25	21.743	40	21.937	55	21.882
11	21.539	26	21.759	41	21.939	56	21.874
12	21.550	27	21.775	42	21.940	57	21.867
13	21.561	28	21.790	43	21.940	58	21.861
14	21.574	29	21.806	44	21.939	59	21.854
15	21.587	30	21.822	45	21.936	60	21.847

WT-S9-171018-459,Potential: 0 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

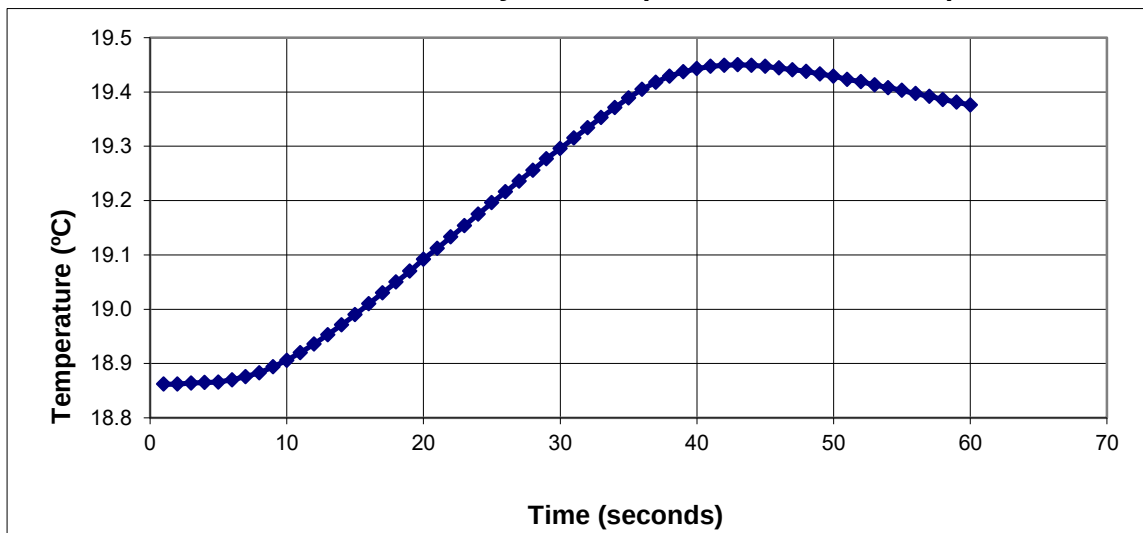
Sample Number: WT-S9-171018-459
Potential (-cm water): Air Dry #1

Test Date/Time: 10/24/2018 7:53 AM	K (W/(m·K)): 1.370
Sensor: SH-1	ρ (°C·cm/W): 73.0
Test Temp.(°C): 18.9	C (MJ/(m ³ ·K)): 3.373
KD2 Pro Sample ID: 459-AD1	D (mm ² /s): 0.406
Power (W/m): 21.390	Err: 0.0012
Current (amps): 0.143	

Raw Data

Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)	
1	18.862	16	19.010	31	19.315	46	19.444
2	18.862	17	19.030	32	19.334	47	19.441
3	18.864	18	19.050	33	19.353	48	19.438
4	18.865	19	19.070	34	19.371	49	19.433
5	18.866	20	19.092	35	19.389	50	19.429
6	18.870	21	19.112	36	19.405	51	19.423
7	18.876	22	19.133	37	19.418	52	19.419
8	18.883	23	19.154	38	19.429	53	19.413
9	18.894	24	19.175	39	19.437	54	19.408
10	18.906	25	19.196	40	19.443	55	19.403
11	18.920	26	19.216	41	19.447	56	19.397
12	18.936	27	19.236	42	19.449	57	19.392
13	18.953	28	19.256	43	19.450	58	19.386
14	18.971	29	19.277	44	19.449	59	19.381
15	18.990	30	19.296	45	19.447	60	19.376

WT-S9-171018-459,Potential: Air Dry #1 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

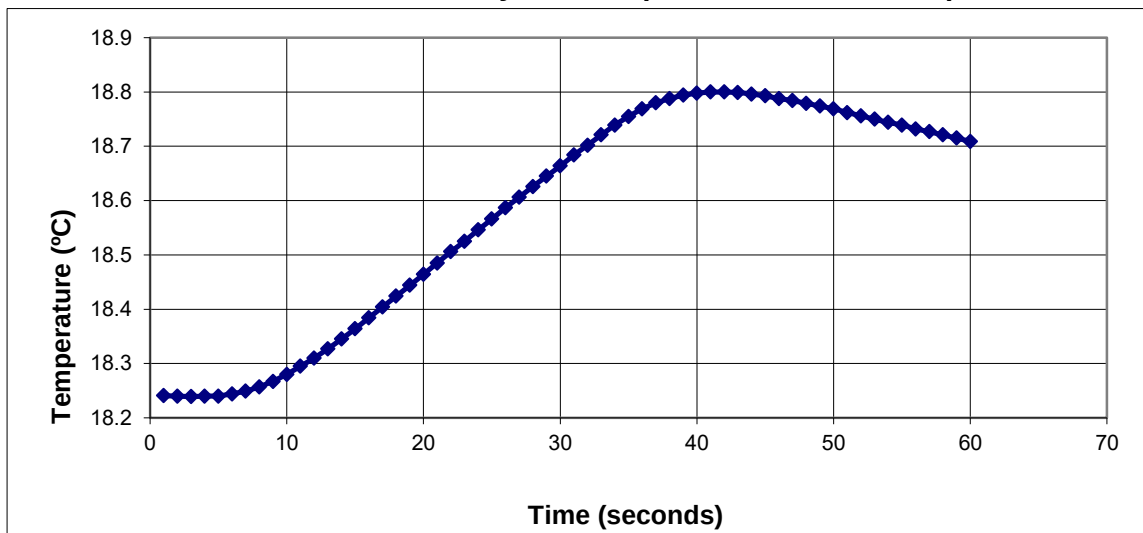
Sample Number: WT-S9-171018-459
Potential (-cm water): Air Dry #2

Test Date/Time:	10/25/2018 8:25 AM	K (W/(m·K)):	1.360
Sensor:	SH-1	ρ (°C·cm/W):	73.5
Test Temp.(°C):	18.2	C (MJ/(m ³ ·K)):	3.241
KD2 Pro Sample ID:	459-AD2	D (mm ² /s):	0.420
Power (W/m):	20.930	Err:	0.0012
Current (amps):	0.142		

Raw Data

Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)	
1	18.241	16	18.384	31	18.684	46	18.788
2	18.240	17	18.404	32	18.702	47	18.784
3	18.239	18	18.424	33	18.721	48	18.779
4	18.240	19	18.444	34	18.739	49	18.774
5	18.240	20	18.464	35	18.755	50	18.769
6	18.244	21	18.485	36	18.769	51	18.762
7	18.249	22	18.506	37	18.780	52	18.756
8	18.257	23	18.525	38	18.788	53	18.750
9	18.267	24	18.546	39	18.794	54	18.744
10	18.280	25	18.566	40	18.798	55	18.739
11	18.295	26	18.587	41	18.800	56	18.732
12	18.310	27	18.606	42	18.800	57	18.727
13	18.327	28	18.626	43	18.799	58	18.721
14	18.345	29	18.645	44	18.796	59	18.715
15	18.364	30	18.664	45	18.793	60	18.709

WT-S9-171018-459,Potential: Air Dry #2 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

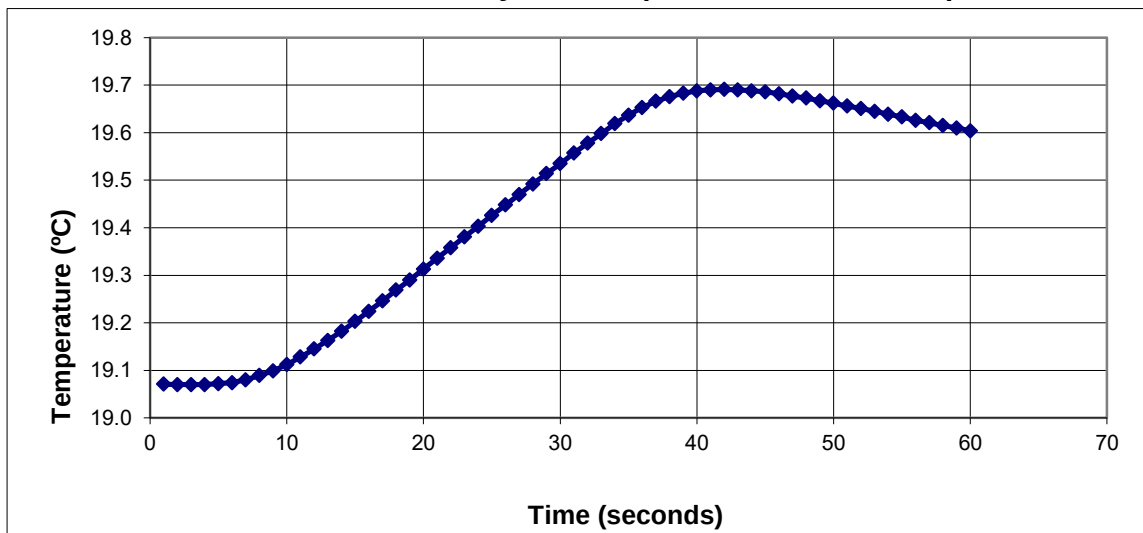
Sample Number: WT-S9-171018-459
Potential (-cm water): Air Dry #3

Test Date/Time:	10/26/2018 2:36 PM	K (W/(m·K)):	1.293
Sensor:	SH-1	ρ (°C·cm/W):	77.3
Test Temp.(°C):	19.1	C (MJ/(m ³ ·K)):	3.166
KD2 Pro Sample ID:	459-AD3	D (mm ² /s):	0.408
Power (W/m):	22.320	Err:	0.0016
Current (amps):	0.147		

Raw Data

Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)	
1	19.071	16	19.224	31	19.557	46	19.682
2	19.070	17	19.246	32	19.578	47	19.677
3	19.070	18	19.269	33	19.598	48	19.673
4	19.070	19	19.290	34	19.619	49	19.667
5	19.072	20	19.313	35	19.637	50	19.662
6	19.074	21	19.336	36	19.653	51	19.656
7	19.080	22	19.358	37	19.666	52	19.651
8	19.089	23	19.381	38	19.676	53	19.645
9	19.099	24	19.403	39	19.683	54	19.639
10	19.112	25	19.426	40	19.688	55	19.633
11	19.128	26	19.448	41	19.690	56	19.626
12	19.145	27	19.470	42	19.691	57	19.621
13	19.163	28	19.492	43	19.690	58	19.615
14	19.182	29	19.514	44	19.688	59	19.610
15	19.203	30	19.535	45	19.686	60	19.604

WT-S9-171018-459,Potential: Air Dry #3 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

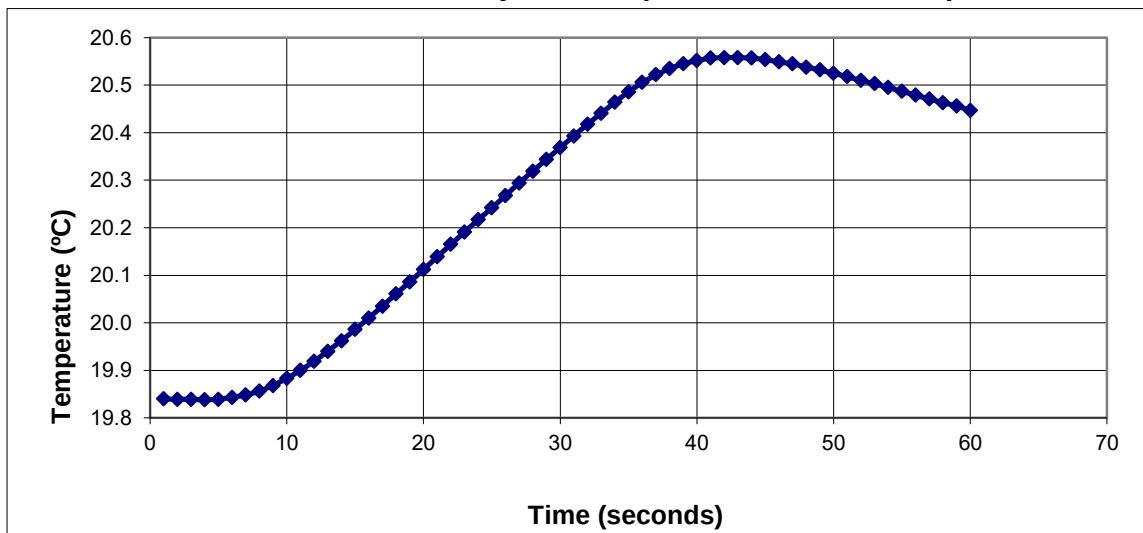
Sample Number: WT-S9-171018-459
Potential (-cm water): Air Dry #4

Test Date/Time:	10/29/2018 7:58 AM	K (W/(m·K)):	1.060
Sensor:	SH-1	ρ (°C·cm/W):	94.3
Test Temp.(°C):	19.8	C (MJ/(m ³ ·K)):	2.670
KD2 Pro Sample ID:	459-AD4	D (mm ² /s):	0.397
Power (W/m):	22.340	Err:	0.0009
Current (amps):	0.147		

Raw Data

Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)	
1	19.840	16	20.010	31	20.393	46	20.549
2	19.839	17	20.035	32	20.418	47	20.545
3	19.839	18	20.061	33	20.441	48	20.538
4	19.838	19	20.086	34	20.464	49	20.532
5	19.839	20	20.112	35	20.486	50	20.525
6	19.843	21	20.139	36	20.506	51	20.518
7	19.848	22	20.165	37	20.522	52	20.510
8	19.856	23	20.191	38	20.535	53	20.503
9	19.868	24	20.217	39	20.545	54	20.495
10	19.883	25	20.242	40	20.552	55	20.487
11	19.900	26	20.268	41	20.557	56	20.479
12	19.919	27	20.294	42	20.558	57	20.471
13	19.940	28	20.319	43	20.558	58	20.463
14	19.962	29	20.344	44	20.557	59	20.456
15	19.986	30	20.369	45	20.554	60	20.447

WT-S9-171018-459,Potential: Air Dry #4 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

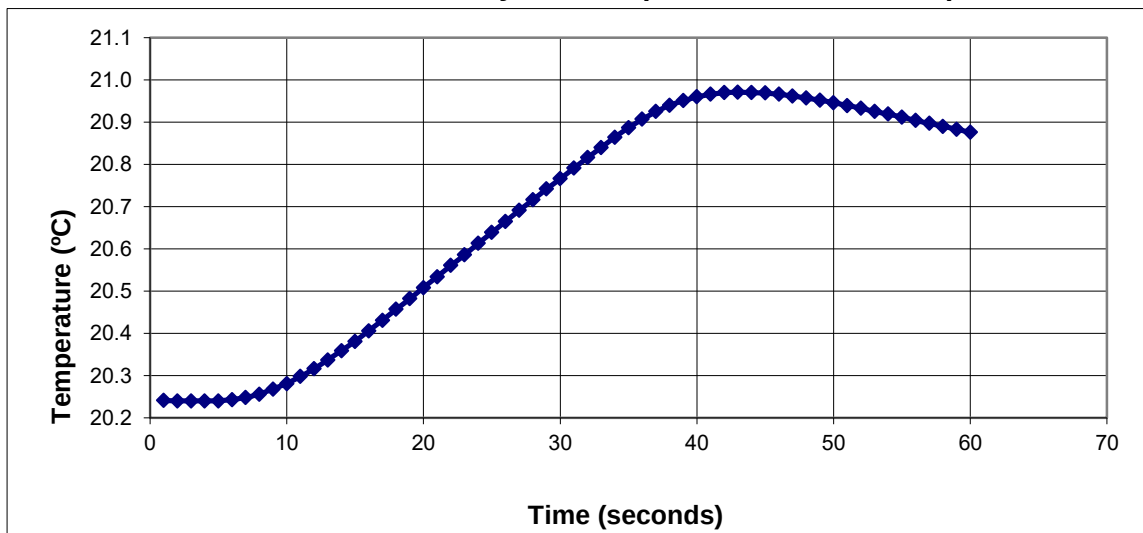
Sample Number: WT-S9-171018-459
Potential (-cm water): Air Dry #5

Test Date/Time: 10/30/2018 8:13 AM	K (W/(m·K)): 0.996
Sensor: SH-1	ρ (°C·cm/W): 100.4
Test Temp.(°C): 20.2	C (MJ/(m ³ ·K)): 2.607
KD2 Pro Sample ID: 459-AD5	D (mm ² /s): 0.382
Power (W/m): 21.700	Err: 0.0009
Current (amps): 0.145	

Raw Data

Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)
1 20.241	16 20.406	31 20.791	46 20.966
2 20.240	17 20.431	32 20.816	47 20.962
3 20.240	18 20.457	33 20.840	48 20.957
4 20.240	19 20.482	34 20.864	49 20.952
5 20.240	20 20.508	35 20.887	50 20.946
6 20.243	21 20.534	36 20.907	51 20.939
7 20.248	22 20.561	37 20.925	52 20.933
8 20.256	23 20.586	38 20.940	53 20.925
9 20.268	24 20.613	39 20.951	54 20.919
10 20.281	25 20.639	40 20.960	55 20.912
11 20.298	26 20.665	41 20.966	56 20.904
12 20.316	27 20.691	42 20.970	57 20.897
13 20.337	28 20.716	43 20.971	58 20.890
14 20.359	29 20.742	44 20.970	59 20.883
15 20.381	30 20.766	45 20.969	60 20.876

WT-S9-171018-459,Potential: Air Dry #5 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

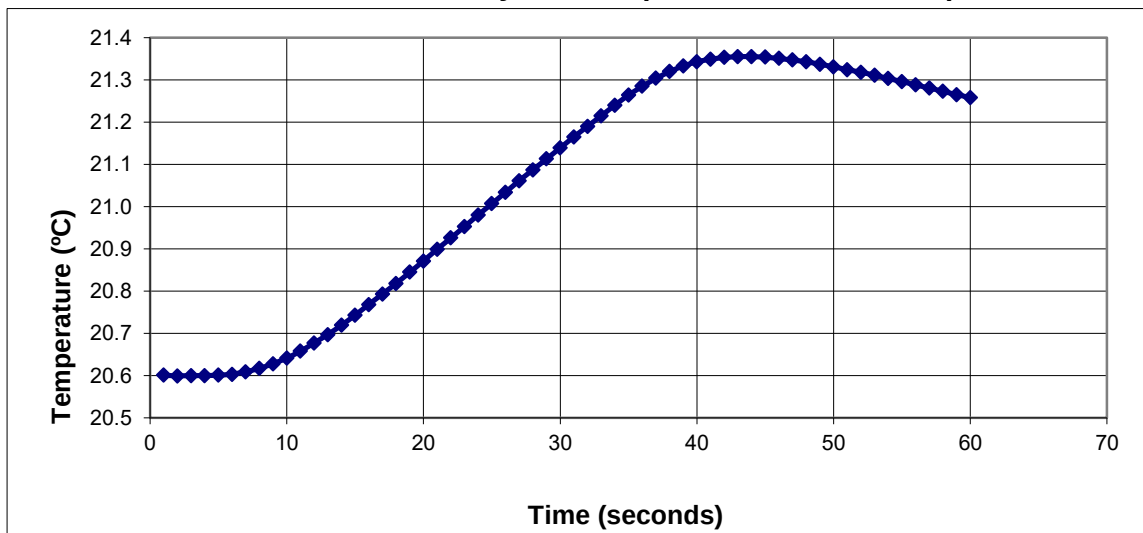
Sample Number: WT-S9-171018-459
Potential (-cm water): Air Dry #6

Test Date/Time:	10/31/2018 9:08 AM	K (W/(m·K)):	0.941
Sensor:	SH-1	ρ (°C·cm/W):	106.2
Test Temp.(°C):	20.6	C (MJ/(m ³ ·K)):	2.502
KD2 Pro Sample ID:	459-AD6	D (mm ² /s):	0.376
Power (W/m):	21.590	Err:	0.0009
Current (amps):	0.144		

Raw Data

Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)	
1	20.601	16	20.768	31	21.165	46	21.351
2	20.599	17	20.793	32	21.190	47	21.347
3	20.600	18	20.818	33	21.215	48	21.343
4	20.600	19	20.845	34	21.240	49	21.337
5	20.601	20	20.871	35	21.264	50	21.331
6	20.603	21	20.899	36	21.285	51	21.324
7	20.609	22	20.926	37	21.304	52	21.318
8	20.617	23	20.953	38	21.320	53	21.311
9	20.628	24	20.980	39	21.333	54	21.303
10	20.641	25	21.007	40	21.343	55	21.296
11	20.658	26	21.034	41	21.349	56	21.288
12	20.677	27	21.061	42	21.353	57	21.281
13	20.697	28	21.087	43	21.355	58	21.273
14	20.719	29	21.113	44	21.355	59	21.265
15	20.743	30	21.139	45	21.354	60	21.258

WT-S9-171018-459,Potential: Air Dry #6 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

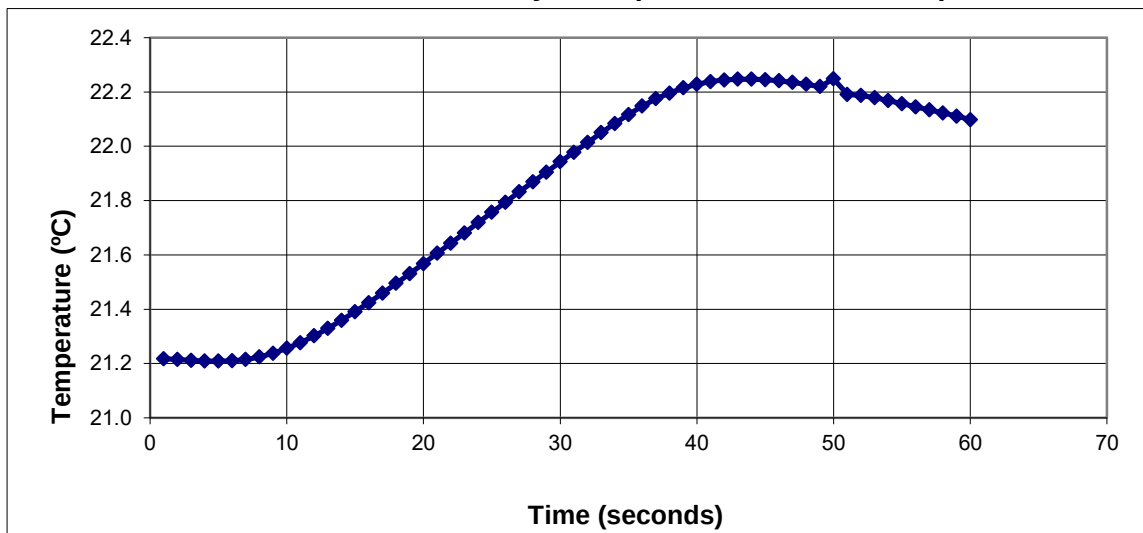
Sample Number: WT-S9-171018-459
Potential (-cm water): Oven Dry

Test Date/Time:	11/6/2018 3:57 PM	K (W/(m·K)):	0.547
Sensor:	SH-1	ρ (°C·cm/W):	182.7
Test Temp.(°C):	21.2	C (MJ/(m ³ ·K)):	1.554
KD2 Pro Sample ID:	459-OD	D (mm ² /s):	0.352
Power (W/m):	20.990	Err:	0.0023
Current (amps):	0.142		

Raw Data

Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)	
1	21.218	16	21.424	31	21.978	46	22.241
2	21.215	17	21.459	32	22.014	47	22.235
3	21.212	18	21.495	33	22.050	48	22.228
4	21.209	19	21.531	34	22.084	49	22.220
5	21.209	20	21.568	35	22.117	50	22.248
6	21.210	21	21.606	36	22.148	51	22.191
7	21.215	22	21.643	37	22.175	52	22.187
8	21.224	23	21.681	38	22.196	53	22.179
9	21.238	24	21.719	39	22.215	54	22.168
10	21.256	25	21.757	40	22.228	55	22.157
11	21.277	26	21.794	41	22.238	56	22.145
12	21.302	27	21.832	42	22.244	57	22.134
13	21.329	28	21.869	43	22.247	58	22.123
14	21.359	29	21.905	44	22.247	59	22.111
15	21.391	30	21.943	45	22.245	60	22.098

WT-S9-171018-459,Potential: Oven Dry - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Results Sheet for Sample: WT-S5-231018-490

Job Name: EA Engineering, Inc.

Job Number: DB18.1333.00

Sample Number: WT-S5-231018-490

Project Name: 62735DM02 KAFB-Bulk Fuels Fac Vadose Zone

PO Number: 18189

Instrument Description: Decagon KD2 Pro

Probe: ☐ KS-1, 6 cm length, 1.3 mm diameter, single needle

☐ TR-1, 10 cm length, 2.4 mm diameter, single needle

☒ SH-1, 3 cm length, 1.3 mm diameter, dual needle, 6 mm spacing

Test Start Date: 10/25/18

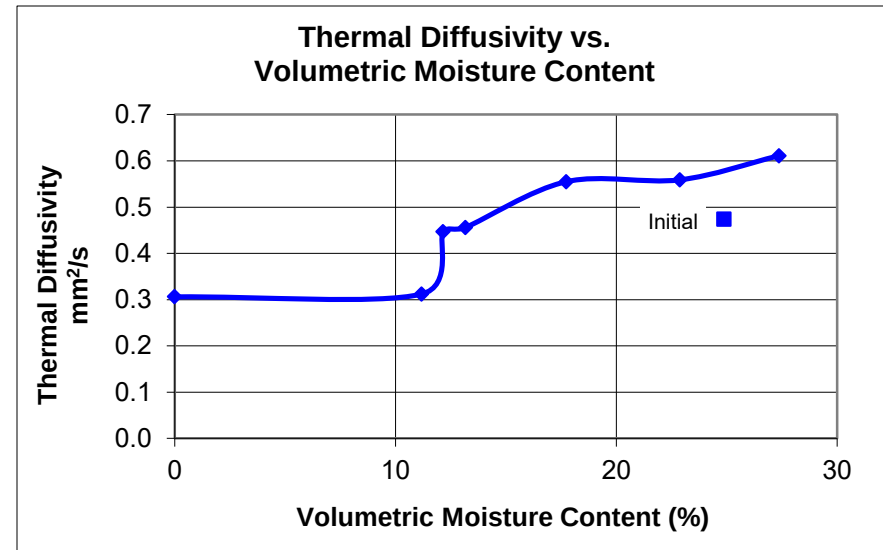
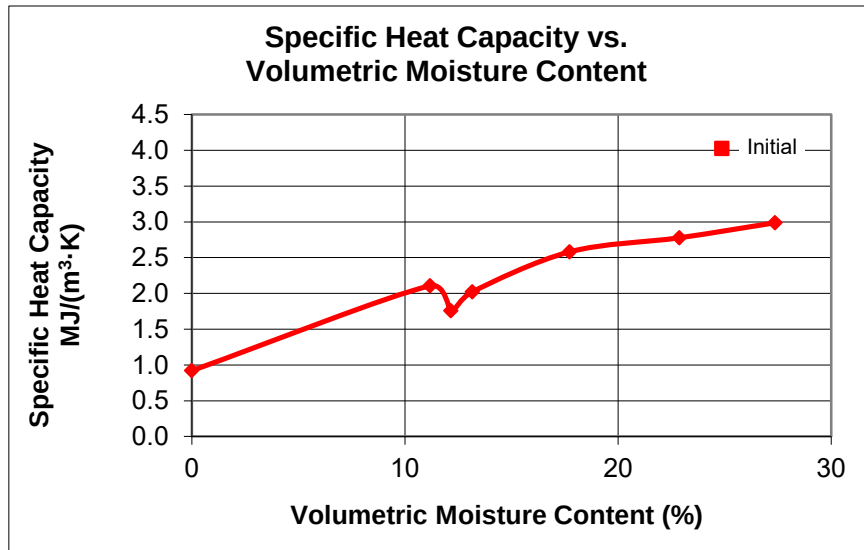
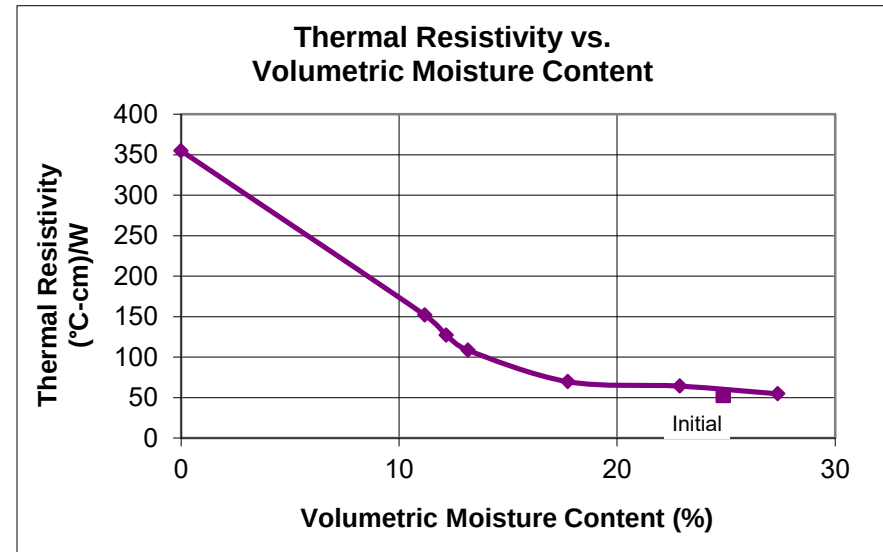
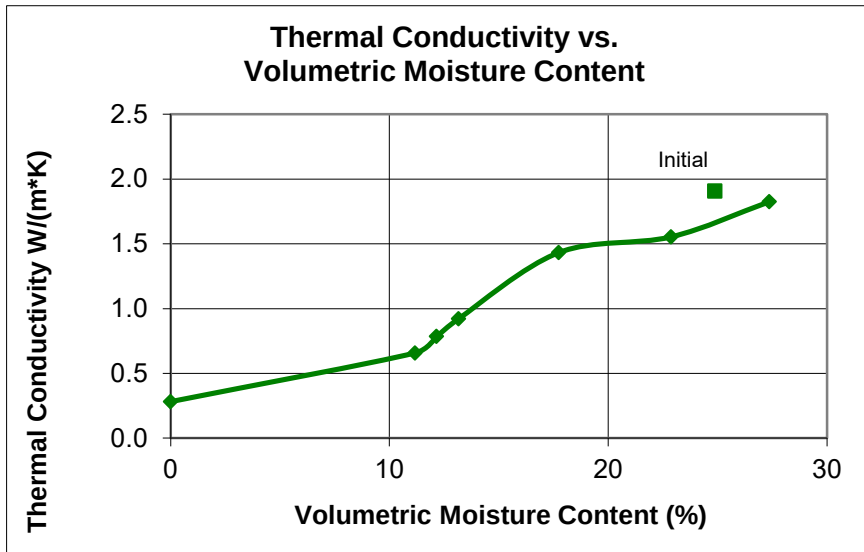
Reading	Water Potential (-cm water)	Gravimetric Moisture Content (g/g, %)	Volumetric Moisture Content ¹ (vol/vol, %)	Dry Bulk Density ¹ (g/cm ³)	Test Temperature (°C)	K Thermal Conductivity W/(m·K)	ρ Thermal Resistivity °C·cm/W	C Specific Heat Capacity MJ/(m ³ ·K)	D Thermal Diffusivity (mm ² /s)
Initial	---	13.13	24.87	1.90	21.72	1.907	52.4	4.024	0.474
Saturated	0	14.44	27.36	1.90	20.78	1.825	54.8	2.988	0.611
Air Dry #1	---	11.75	22.88	1.95	18.34	1.554	64.3	2.778	0.559
Air Dry #2	---	9.08	17.73	1.95	18.66	1.432	69.8	2.582	0.555
Air Dry #3	---	6.74	13.17	1.95	18.80	0.922	108.5	2.023	0.456
Air Dry #4	---	6.22	12.16	1.95	18.89	0.785	127.4	1.756	0.447
Air Dry #5	---	5.72	11.18	1.95	18.83	0.658	152.1	2.106	0.312
Oven Dry	0	0	0	1.95	21.76	0.282	355.0	0.921	0.306

--- = Value not measured.

¹ Adjusted for volume changes during testing, if applicable.



Thermal Properties Results Sheet for Sample: WT-S5-231018-490
Scatter Plots





Thermal Properties Data

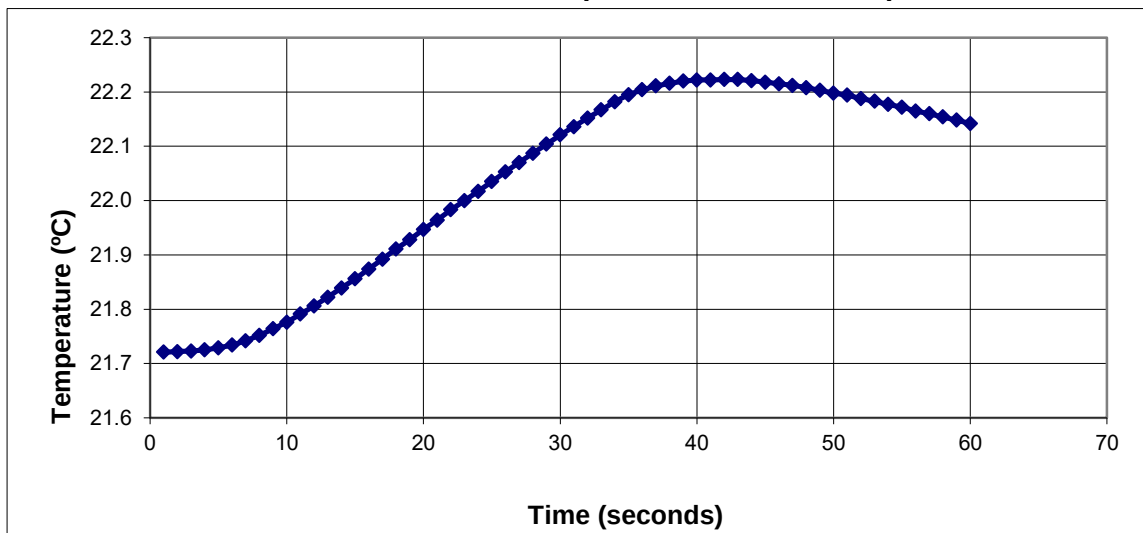
Sample Number: WT-S5-231018-490
Potential (-cm water): Initial

Test Date/Time: 10/25/2018 1:13 PM	K (W/(m·K)): 1.907
Sensor: SH-1	ρ (°C·cm/W): 52.4
Test Temp.(°C): 21.7	C (MJ/(m ³ ·K)): 4.024
KD2 Pro Sample ID: 490-AR	D (mm ² /s): 0.474
Power (W/m): 21.590	Err: 0.0025
Current (amps): 0.144	

Raw Data

Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)	
1	21.721	16	21.874	31	22.136	46	22.215
2	21.722	17	21.892	32	22.152	47	22.212
3	21.723	18	21.911	33	22.167	48	22.208
4	21.725	19	21.928	34	22.182	49	22.203
5	21.729	20	21.947	35	22.195	50	22.198
6	21.734	21	21.964	36	22.204	51	22.194
7	21.742	22	21.983	37	22.211	52	22.188
8	21.752	23	22.000	38	22.216	53	22.183
9	21.764	24	22.017	39	22.220	54	22.177
10	21.776	25	22.035	40	22.222	55	22.172
11	21.791	26	22.053	41	22.222	56	22.165
12	21.806	27	22.070	42	22.223	57	22.160
13	21.822	28	22.087	43	22.223	58	22.154
14	21.839	29	22.104	44	22.221	59	22.148
15	21.856	30	22.121	45	22.218	60	22.142

WT-S5-231018-490,Potential: Initial - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

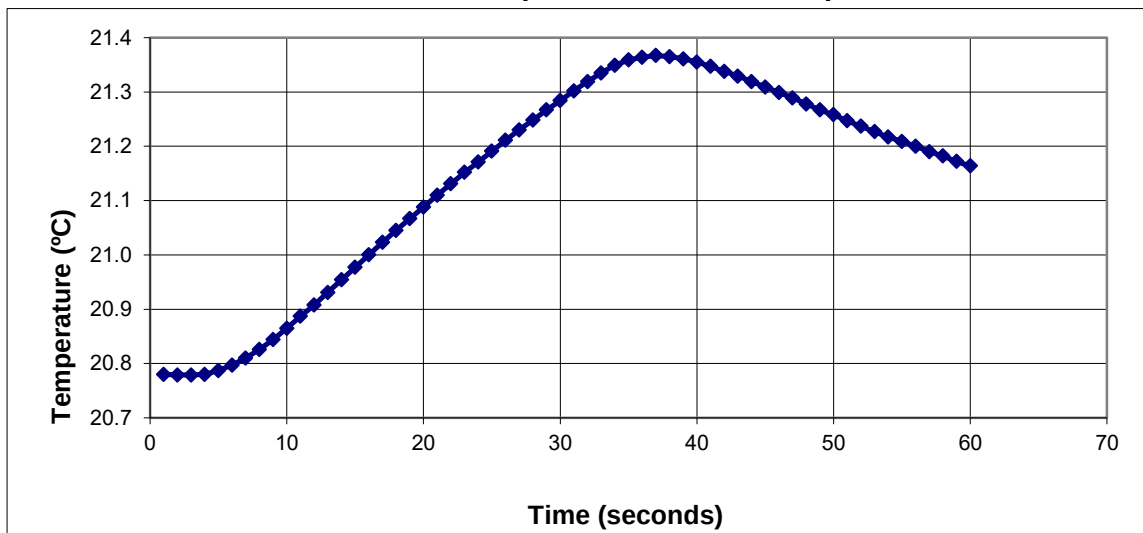
Sample Number: WT-S5-231018-490
Potential (-cm water): 0

Test Date/Time: 10/26/2018 10:35 AM	K (W/(m·K)): 1.825
Sensor: SH-1	ρ (°C·cm/W): 54.8
Test Temp.(°C): 20.8	C (MJ/(m ³ ·K)): 2.988
KD2 Pro Sample ID: 490-SA	D (mm ² /s): 0.611
Power (W/m): 21.890	Err: 0.0024
Current (amps): 0.145	

Raw Data

Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)	
1	20.780	16	21.000	31	21.302	46	21.299
2	20.779	17	21.023	32	21.319	47	21.289
3	20.779	18	21.045	33	21.335	48	21.278
4	20.780	19	21.067	34	21.349	49	21.267
5	20.787	20	21.088	35	21.359	50	21.258
6	20.797	21	21.110	36	21.364	51	21.247
7	20.810	22	21.131	37	21.367	52	21.237
8	20.826	23	21.152	38	21.365	53	21.227
9	20.844	24	21.171	39	21.361	54	21.217
10	20.865	25	21.191	40	21.355	55	21.209
11	20.887	26	21.211	41	21.347	56	21.200
12	20.908	27	21.230	42	21.338	57	21.190
13	20.931	28	21.248	43	21.329	58	21.182
14	20.954	29	21.267	44	21.319	59	21.172
15	20.977	30	21.284	45	21.309	60	21.164

WT-S5-231018-490,Potential: 0 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

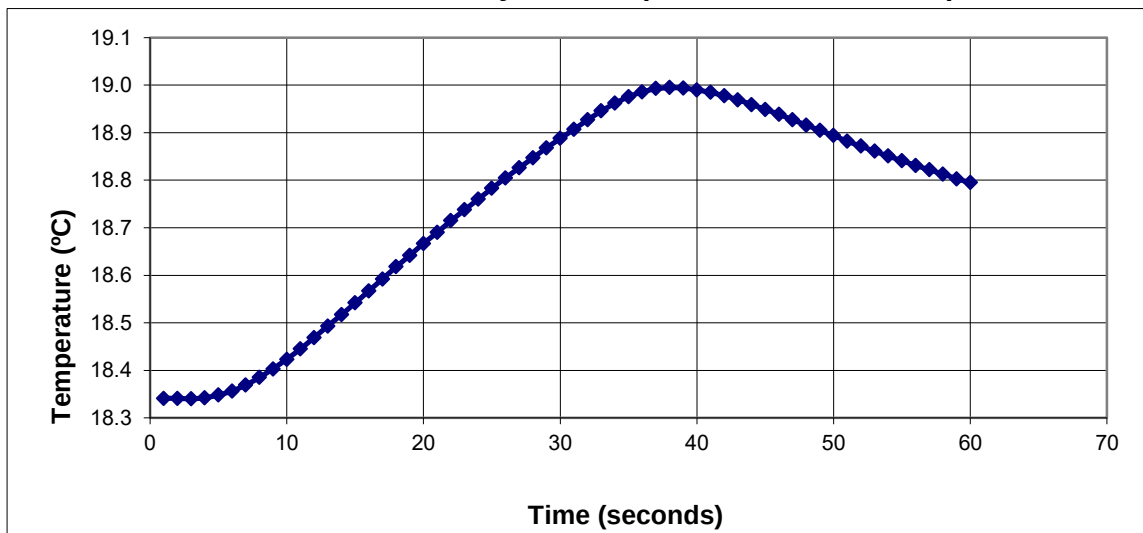
Sample Number: WT-S5-231018-490
Potential (-cm water): Air Dry #1

Test Date/Time: 10/29/2018 10:11 AM	K (W/(m·K)): 1.554
Sensor: SH-1	ρ (°C·cm/W): 64.3
Test Temp.(°C): 18.3	C (MJ/(m ³ ·K)): 2.778
KD2 Pro Sample ID: 490-AD1	D (mm ² /s): 0.559
Power (W/m): 22.280	Err: 0.0027
Current (amps): 0.146	

Raw Data

Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)	
1	18.341	16	18.567	31	18.907	46	18.939
2	18.341	17	18.592	32	18.927	47	18.927
3	18.340	18	18.618	33	18.946	48	18.916
4	18.342	19	18.642	34	18.962	49	18.905
5	18.348	20	18.667	35	18.976	50	18.894
6	18.356	21	18.690	36	18.986	51	18.882
7	18.369	22	18.715	37	18.993	52	18.872
8	18.385	23	18.738	38	18.995	53	18.861
9	18.403	24	18.760	39	18.994	54	18.851
10	18.423	25	18.783	40	18.990	55	18.841
11	18.445	26	18.805	41	18.985	56	18.831
12	18.469	27	18.826	42	18.978	57	18.822
13	18.493	28	18.847	43	18.969	58	18.813
14	18.517	29	18.868	44	18.959	59	18.803
15	18.542	30	18.888	45	18.949	60	18.795

WT-S5-231018-490,Potential: Air Dry #1 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

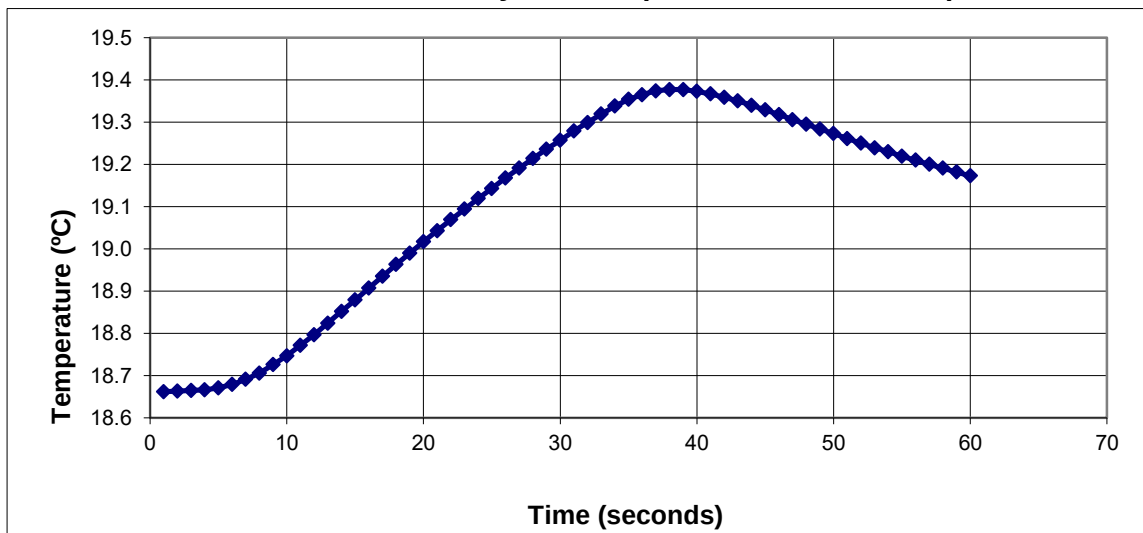
Sample Number: WT-S5-231018-490
Potential (-cm water): Air Dry #2

Test Date/Time: 10/29/2018 3:38 PM	K (W/(m·K)): 1.432
Sensor: SH-1	ρ (°C·cm/W): 69.8
Test Temp.(°C): 18.7	C (MJ/(m ³ ·K)): 2.582
KD2 Pro Sample ID: 490-AD2	D (mm ² /s): 0.555
Power (W/m): 22.230	Err: 0.0035
Current (amps): 0.146	

Raw Data

Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)
1 18.662	16 18.907	31 19.279	46 19.318
2 18.663	17 18.935	32 19.299	47 19.306
3 18.665	18 18.963	33 19.319	48 19.295
4 18.666	19 18.990	34 19.338	49 19.284
5 18.671	20 19.017	35 19.354	50 19.273
6 18.679	21 19.043	36 19.365	51 19.261
7 18.691	22 19.069	37 19.374	52 19.250
8 18.706	23 19.094	38 19.377	53 19.239
9 18.726	24 19.119	39 19.377	54 19.230
10 18.747	25 19.143	40 19.373	55 19.219
11 18.772	26 19.168	41 19.367	56 19.210
12 18.797	27 19.191	42 19.359	57 19.200
13 18.824	28 19.214	43 19.350	58 19.191
14 18.852	29 19.236	44 19.340	59 19.182
15 18.879	30 19.257	45 19.329	60 19.173

WT-S5-231018-490,Potential: Air Dry #2 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

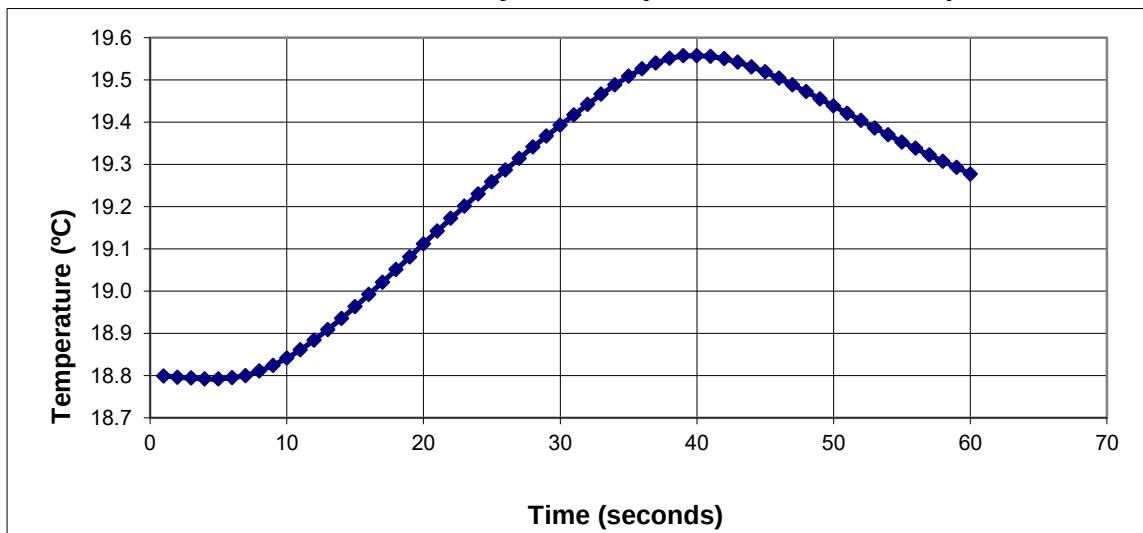
Sample Number: WT-S5-231018-490
Potential (-cm water): Air Dry #3

Test Date/Time:	10/30/2018 8:24 AM	K (W/(m·K)):	0.922
Sensor:	SH-1	ρ (°C·cm/W):	108.5
Test Temp.(°C):	18.8	C (MJ/(m ³ ·K)):	2.023
KD2 Pro Sample ID:	490-AD3	D (mm ² /s):	0.456
Power (W/m):	21.890	Err:	0.0037
Current (amps):	0.145		

Raw Data

Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)	
1	18.799	16	18.992	31	19.417	46	19.504
2	18.796	17	19.021	32	19.442	47	19.488
3	18.794	18	19.051	33	19.466	48	19.472
4	18.792	19	19.081	34	19.488	49	19.455
5	18.792	20	19.112	35	19.509	50	19.438
6	18.795	21	19.142	36	19.526	51	19.421
7	18.800	22	19.172	37	19.540	52	19.404
8	18.811	23	19.201	38	19.551	53	19.386
9	18.824	24	19.230	39	19.557	54	19.370
10	18.841	25	19.259	40	19.557	55	19.353
11	18.861	26	19.287	41	19.556	56	19.338
12	18.884	27	19.314	42	19.550	57	19.322
13	18.909	28	19.341	43	19.542	58	19.307
14	18.935	29	19.367	44	19.531	59	19.293
15	18.963	30	19.393	45	19.519	60	19.277

WT-S5-231018-490,Potential: Air Dry #3 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

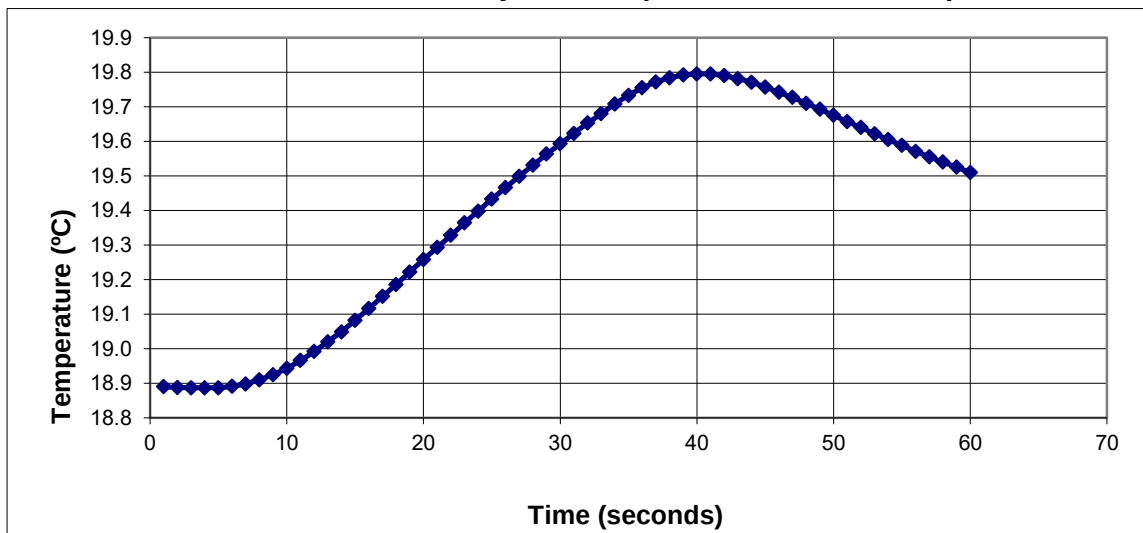
Sample Number: WT-S5-231018-490
Potential (-cm water): Air Dry #4

Test Date/Time: 10/31/2018 9:31 AM	K (W/(m·K)): 0.785
Sensor: SH-1	ρ (°C·cm/W): 127.4
Test Temp.(°C): 18.9	C (MJ/(m ³ ·K)): 1.756
KD2 Pro Sample ID: 490-AD4	D (mm ² /s): 0.447
Power (W/m): 21.650	Err: 0.0034
Current (amps): 0.144	

Raw Data

Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)	Second Temp.(°C)
1 18.890	16 19.116	31 19.623	46 19.742
2 18.888	17 19.151	32 19.653	47 19.727
3 18.887	18 19.186	33 19.680	48 19.710
4 18.887	19 19.222	34 19.708	49 19.693
5 18.887	20 19.258	35 19.733	50 19.675
6 18.891	21 19.293	36 19.755	51 19.657
7 18.898	22 19.328	37 19.772	52 19.640
8 18.910	23 19.364	38 19.784	53 19.622
9 18.925	24 19.398	39 19.792	54 19.605
10 18.943	25 19.433	40 19.795	55 19.588
11 18.966	26 19.466	41 19.795	56 19.571
12 18.992	27 19.499	42 19.790	57 19.555
13 19.020	28 19.531	43 19.781	58 19.540
14 19.049	29 19.563	44 19.771	59 19.525
15 19.082	30 19.593	45 19.757	60 19.510

WT-S5-231018-490,Potential: Air Dry #4 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

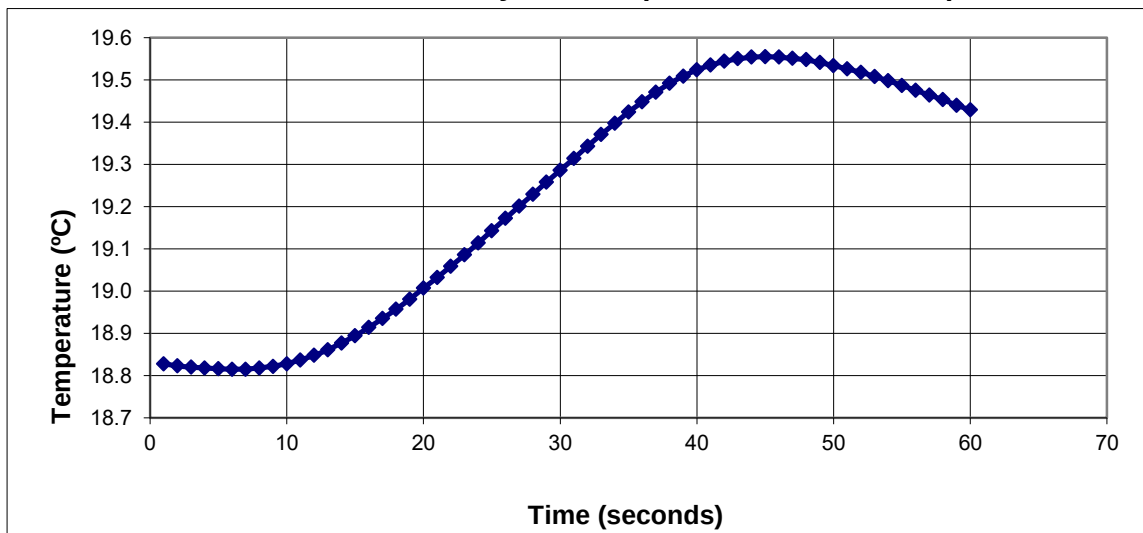
Sample Number: WT-S5-231018-490
Potential (-cm water): Air Dry #5

Test Date/Time: 11/1/2018 8:07 AM	K (W/(m·K)): 0.658
Sensor: SH-1	ρ (°C·cm/W): 152.1
Test Temp.(°C): 18.8	C (MJ/(m ³ ·K)): 2.106
KD2 Pro Sample ID: 490-AD5	D (mm ² /s): 0.312
Power (W/m): 21.570	Err: 0.0023
Current (amps): 0.144	

Raw Data

Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)	
1	18.828	16	18.914	31	19.314	46	19.554
2	18.823	17	18.935	32	19.343	47	19.551
3	18.820	18	18.957	33	19.371	48	19.548
4	18.818	19	18.981	34	19.397	49	19.541
5	18.816	20	19.007	35	19.424	50	19.534
6	18.815	21	19.032	36	19.448	51	19.526
7	18.815	22	19.059	37	19.471	52	19.518
8	18.818	23	19.086	38	19.492	53	19.508
9	18.822	24	19.114	39	19.509	54	19.498
10	18.828	25	19.143	40	19.524	55	19.487
11	18.837	26	19.172	41	19.535	56	19.475
12	18.848	27	19.201	42	19.544	57	19.464
13	18.861	28	19.229	43	19.550	58	19.453
14	18.877	29	19.258	44	19.554	59	19.440
15	18.894	30	19.286	45	19.555	60	19.429

WT-S5-231018-490,Potential: Air Dry #5 - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines



Thermal Properties Data

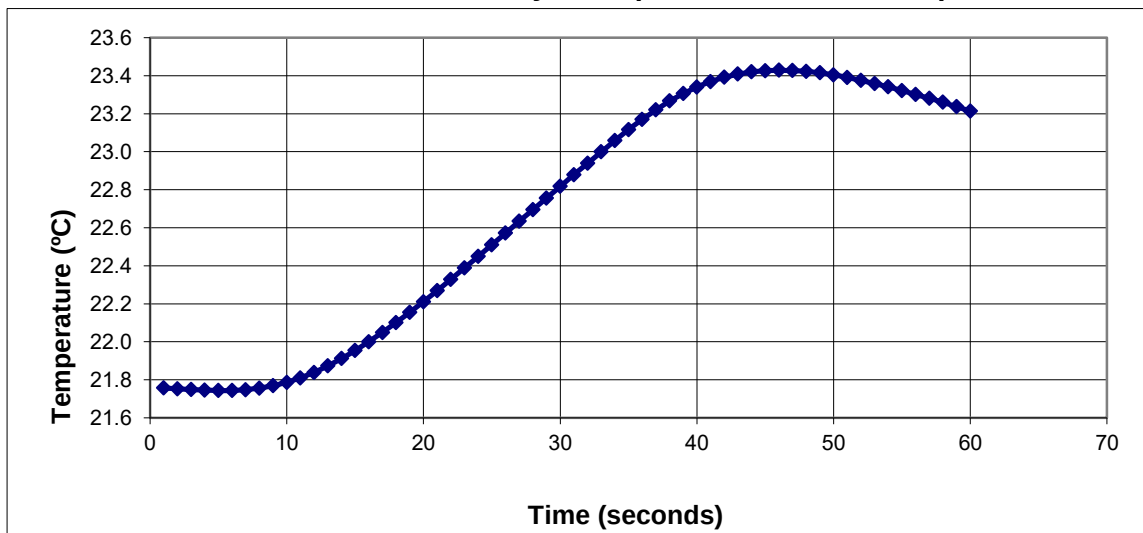
Sample Number: WT-S5-231018-490
Potential (-cm water): Oven Dry

Test Date/Time: 11/6/2018 3:43 PM	K (W/(m·K)): 0.282
Sensor: SH-1	ρ (°C·cm/W): 355.0
Test Temp.(°C): 21.8	C (MJ/(m ³ ·K)): 0.921
KD2 Pro Sample ID: 490-OD	D (mm ² /s): 0.306
Power (W/m): 20.950	Err: 0.0018
Current (amps): 0.142	

Raw Data

Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)		Second Temp.(°C)	
1	21.757	16	22.000	31	22.879	46	23.429
2	21.753	17	22.049	32	22.939	47	23.427
3	21.749	18	22.101	33	22.999	48	23.423
4	21.746	19	22.155	34	23.058	49	23.415
5	21.744	20	22.211	35	23.116	50	23.404
6	21.744	21	22.269	36	23.170	51	23.391
7	21.748	22	22.329	37	23.220	52	23.375
8	21.756	23	22.389	38	23.267	53	23.359
9	21.769	24	22.449	39	23.306	54	23.341
10	21.787	25	22.510	40	23.340	55	23.322
11	21.810	26	22.572	41	23.369	56	23.301
12	21.839	27	22.634	42	23.392	57	23.281
13	21.873	28	22.695	43	23.408	58	23.260
14	21.912	29	22.756	44	23.420	59	23.237
15	21.954	30	22.818	45	23.426	60	23.214

WT-S5-231018-490,Potential: Oven Dry - Temperature vs. Time Graph



Laboratory analysis by: D. O'Dowd
Data entered by: C. Krous
Checked by: J. Hines

Laboratory Tests and Methods



Tests and Methods

Dry Bulk Density:	ASTM D7263
Moisture Content:	ASTM D7263, ASTM D2216
Calculated Porosity:	ASTM D7263
Thermal Properties:	ASTM D5334



Daniel B. Stephens & Associates, Inc.

Daniel B. Stephens & Associates, Inc.
Hydraulic Properties Soils Laboratory
4400 Alameda Blvd. NE, Suite C
ALBUQUERQUE, NM 87113
Ph: (505) 889-7752 (800) 933-3105
FAX (505) 889-0258

Name/Address/Phone for report:

Bonnie Backisch
EA Engineering, Science, + Tech
320 Gold Ave, SW, Ste 1300
Albuquerque, NM 87102
bbackisch@eaerst.com 505-234-1105

cc: Name/Address/Phone for invoice: (or same as report)

Same as Report.
p.moss@eaerst.com

Project Name or P.O. # PO 18189

Kirkland Vadosezone

Project Contact & Phone:

161018

DATE SAMPLE IDENTIFICATION

DATE	SAMPLE IDENTIFICATION	# of containers	Type of Container	As Received	Initial Soil Properties	Saturated Hydraulic Conductivity	Hydraulic Properties Package (HPP) with	Hydraulic Properties Package (HPP) with	Particle Size Analysis	Proctor Compaction	Specific Gravity, Fines	Specific Gravity, Coarse	Atterberg Limits	Thermal Properties, 1-point	Thermal Properties, 5-6-point	REMARKS
				Gravimetric Moisture Content	(density, moisture, porosity)	(rigid wall or flexible wall)	Flexible Wall Ksat	Rigid Wall Ksat	(wet sieve & hydrometer)	(standard or modified)						
10/16/18	W-59-171018-435	1	Bag	X												
10/17/18	WT-59-171018-459	1	Bag	X										X		
10/17/18	WT-59-171018-464	1	Bag	X										X		

RELINQUISHED BY: (SIGNATURE)

DATE

10/19/18

RECEIVED BY: (SIGNATURE)

DATE

10/19/18

Comments:

APPENDIX H

Data Quality Evaluation Report – Source Zone Characterization Samples (October 2018 – February 2019)

LIST OF ACRONYMS AND ABBREVIATIONS

%	percent
AFB	Air Force Base
ASTM	ASTM International
BFF	Bulk Fuels Facility
CCV	continuing calibration verification
DL	detection limit
DoD	Department of Defense
DRO	diesel range organic
EDB	ethylene dibromide
EPA	U.S. Environmental Protection Agency
GRO	gasoline range organic
IS	internal standard
LCS	laboratory control sample
LCSD	laboratory control sample duplicate
LOD	limit of detection
LOQ	limit of quantification
MS	matrix spike
MSD	matrix spike duplicate
QAPjP	Quality Assurance Project Plan
QC	quality control
QSM	Quality Systems Manual
RB	equipment rinse blank
RL	reporting limit
RPD	relative percent different
SDG	sample delivery group
SW	Solid Waste
SWMU	Solid Waste Management Unit
TA	TestAmerica Laboratories, Inc.
TPH	total petroleum hydrocarbon
USACE	U.S. Army Corps of Engineers
VOC	volatile organic compound

DATA QUALITY EVALUATION REPORT

Source Zone Characterization Samples

(October 2018 – February 2019)

1. LABORATORY DATA QUALITY SUMMARY

This Data Quality Evaluation Report describes the findings of the data validation performed for the source zone characterization chemical analytical data collected in support of the Work Plan for Vadose Zone Coring, Vapor Monitoring, and Water Supply Sampling Revision 2, Bulk Fuels Facility (BFF), Solid Waste Management (SWMU) ST-106/SS-111, Kirtland Air Force Base (AFB), New Mexico (U.S. Army Corps of Engineers [USACE], 2017a). Sampling and analysis for the source zone characterization coring activities were conducted in accordance with the procedures and overall quality control (QC) and quality assurance protocols presented in the Work Plan and Quality Assurance Project Plan (QAPjP), Appendix A (USACE, 2017b).

Seventy-three soil samples for volatile organic compound (VOC) and ethylene dibromide (EDB) analysis and 87 samples for total petroleum hydrocarbon (TPH) analysis were collected from the continuous coring of 11 monitoring wells installed during the period of October 6, 2018 through February 5, 2019. The wells were installed to obtain data to support vadose zone treatability studies to address data gaps in the horizontal and vertical extent of light non-aqueous phase liquids in the vadose zone and groundwater associated with Kirtland AFB BFF, SWMU ST-106/SS-111.

The following field QC samples were associated with the vadose zone soil coring and sampling: 10 field duplicates, 18 equipment rinse blanks (RBs), 46 trip blanks, and 1 source water blank for equipment rinse water. Samples were shipped to TestAmerica Laboratories, Inc. (TA), Arvada, Colorado for analysis. TA maintains a current Department of Defense (DoD) Environmental Laboratory Accreditation Program certification to perform the analyses required for this project. Sample analyses were performed in accordance with the following guidance documents:

- DoD Quality Systems Manual (QSM), Version 5.1.1 (DoD, 2018).
- U.S. Environmental Protection Agency (EPA) Test Methods for Evaluating Solid Waste (SW-846), Third Edition and Updates, 1986.
- ASTM International (ASTM) Standard Test Method for Laboratory Determination of Water (Moisture) Content of Soil and Rock by Mass, 2005.

Source zone soil samples were analyzed for the following list of parameters and methods as indicated in Appendix H – Table 1.

- **VOCs**—Method SW8260B.
- **EDB**—Method SW8011.
- **TPH-Gasoline Range Organics (GRO), Diesel Range Organics (DRO) and Motor Oil**—Method SW8015D modified.
- **Percent Moisture**—Method ASTM D2216-90.

Chemical analytical data for the source zone soil samples were reported by TA in 52 sample delivery groups (SDGs). Appendix H– Table 1 summarizes samples collected from coring and the associated field QC samples, collection date, laboratory SDG, and analytical parameters.

A third-party subcontractor, Environmental Data Services, Virginia Beach, Virginia, conducted EPA Stage 3 data validation on 100 percent (%) of the source zone coring sample data.

Analytical data validation was performed using the quality criteria specified in the following analytical guidelines and methods as applicable:

- Work Plan (USACE, 2017a) and QAPjP (USACE, 2017b)
- DoD QSM for Environmental Laboratories, Version 5.1.1 (DoD, 2018)
- EPA Test Methods for Evaluating Solids Waste, Physical/Chemical Methods (SW 846, Third Edition and updates) (1986)
- EPA Contract Laboratory Program, National Functional Guidelines for Superfund Organic Methods Data Review (EPA, 2014).

The following QC criteria were included in the EPA Stage 3 validation per the QAPjP, as applicable to the analytical method:

- Sample preservation and extraction and analysis holding times
- Laboratory method blank contamination
- Surrogate spike and internal standard (IS) recoveries (organic analyses)
- Laboratory control sample (LCS) and LCS duplicate (LCSD) recoveries
- Matrix spike (MS) and matrix spike duplicate (MSD) sample recoveries
- Relative percent difference (RPD)
- Initial and continuing calibrations
- Second column confirmation (for EDB only)
- Trip, RB, and source water blank results
- Field duplicate sample precision.

Analytical data were reviewed to evaluate precision, accuracy (bias), representativeness, comparability, completeness, and sensitivity as defined below:

- *Precision* is expressed as the RPD between the results of replicate sample analyses: sample duplicates, LCSDs, and MSDs. When analyte RPDs exceed the acceptance criteria, the data are qualified accordingly.
- *Accuracy (bias)* is demonstrated by recovery of target analytes from fortified blank and sample matrices, LCS/LCSD, and MS/MSD, respectively. For organic methods, bias is also demonstrated through recovery of surrogates from each field and QC sample. A comparison was made from the recovery of target analytes from fortified samples to the acceptance criteria defined in the QAPjP (USACE, 2017b) and DoD QSM (2018). When the acceptance criteria are not available in the QAPjP or DoD QSM, results are compared with the laboratory in-house control limits. When these criteria are not met, the data are qualified accordingly. Bias may be indicated as high or low.

- *Representativeness* of the samples submitted for analysis is ensured by adherence to standard sampling techniques and standard analytical method protocols.
- *Comparability* of sample results is ensured through the use of approved sampling and analysis methods and comparison of sample results to historical sample data.
- *Completeness* of data is evaluated based on contractual, analytical, and technical criteria for data analysis. Technical completeness of data is used to assess overall project data completeness and is expressed as a percentage of the ratio of the number of usable data results to the total number of analytical data results. Only rejected data (R-qualified) are considered not usable to achieve project objectives.
- *Sensitivity* is determined by the ability to achieve the established method-specific reporting limits (RLs) in accordance with DoD QSM requirements and includes establishing the detection limit (DL), limit of detection (LOD), and limit of quantitation (LOQ). For this project, the laboratory reported positive results to the DL, and results between the DL and LOQ are flagged with a J-qualifier and reported as estimated data. Sensitivity will be evaluated based on comparison of the sample RLs to the project screening levels.

The following sections present the EPA Stage 3 data validation findings. The discussion summarizes data quality exceedances and their potential impact on the quality and usability of analytical results.

Appendix H – Table 2 presents the definitions of data qualification and the reasons why these codes were applied to the analytical results. Appendix H – Table 3 summarizes the qualified analytical data for the source zone coring sample results based on data validation findings.

1.1 DATA QUALITY FINDINGS

1.1.1 Sample Preservation and Sample Extraction and Analysis Holding Times (Reason Code HT)

The sample coolers and samples contained within the coolers were received intact at the laboratory at ≤ 6 degrees Celsius, per EPA guidelines. Samples were preserved appropriately per the requirements of EPA methods with exceptions noted below. Sample holding times were evaluated by comparing (1) the sample collection date to the sample extraction date, and (2) the extraction date to the analysis date to determine if the method-specified holding times were exceeded. Sample receipt discrepancies, preservation, and holding time exceedances include the following:

- Terra Core® samplers for 13 soil samples for low-level VOC analysis were not able to be frozen within 48 hours due to sampling and shipping logistics but were kept on ice below 6 degrees Celsius and were immediately frozen upon receipt at the laboratory.
- One vial for sample TB-S5-231018-02 was received broken and could not be analyzed for VOCs.
- One methanol vial for sample V-V2-141218-215 for VOC analysis was received broken at the laboratory and the VOCs were analyzed from the associated 4-ounce glass jar.
- VOC samples including trip blank and RB samples were analyzed outside of but within 1 week of the 14-day hold time.

- GRO field and trip blank samples were analyzed outside of but within 1 week of the 14-day hold time.
- DRO field and RB samples were extracted outside of but within 1 week of the 14-day hold time.
- One EDB field duplicate sample was analyzed 7 days outside of the 14-day hold time.

Qualified sample data based on holding time exceedance are presented on Appendix H – Table 3.

1.1.2 Laboratory Method Blanks (Reason Code MB)

The field sample results were evaluated with respect to the laboratory method blank prepared and analyzed for each analytical batch for each analytical method. Based on the DoD QSM requirements (2018), laboratory method blank concentrations are considered acceptable when contaminant levels in the blank are less than one-half the LOQ for target analytes and less than the LOQ for common laboratory contaminants. Detections of VOCs (acetone, methylene chloride, chloroform, toluene, and naphthalene), GRO, and DRO were reported in method blank samples. Associated sample detections were “U” qualified, signifying non-detect sample results if the analyte result in the sample was less than five times the amount reported in the blank; or less than 10 times the amount for common laboratory contaminants (acetone, methylene chloride). Qualified sample data based on method blank detections are presented on Appendix H – Table 3.

1.1.3 Initial and Continuing Calibration Blanks (Reason Code CB/CCB)

Initial and continuing calibration blank criteria were reviewed to ensure that the instruments were free of contamination prior to sample analysis.

Based on the DoD QSM requirements (2018), calibration blank concentrations are considered acceptable when contaminant levels in the blank are less than one-half the LOQ for target analytes and less than the LOQ for common laboratory contaminants. Initial and continuing calibration blank data were within control criteria for the source zone sample analyses and no data were qualified based on calibration blank criteria.

1.1.4 Surrogate and Internal Standard Recoveries (Reason Code SURR/IS)

Surrogate and internal compounds are added to field and laboratory QC samples for organic analysis to evaluate the matrix effect and method performance on an individual sample basis. IS recoveries for the source zone sample data were within method control criteria. Surrogate recoveries for VOC, TPH-GRO, and TPH-DRO analyses resulted in data qualification. Qualified sample data based on surrogate recoveries are presented on Appendix H – Table 3.

1.1.5 Laboratory Control Sample/Laboratory Control Sample Duplicate Recoveries and Precision (Reason Codes LCS/RPD)

The LCS is an aliquot of an analyte-free matrix spiked with target analytes that are prepared with each analytical batch for each analytical method. The recovery of target analytes from the LCS analysis is a measurement of method performance in an interference-free sample matrix. LCS recoveries for the source zone coring analyses were within method control limits or did not result in data qualification with one exception. One LCS recovery for TPH motor oil exceeded the control limit and resulted in one RB

result “J” qualified for sample RB-S3-271118-01 (Appendix H – Table 3). All other LCS recoveries for the source zone analyses were within method control criteria.

1.1.6 Matrix Spike/Matrix Spike Duplicate Recoveries and Precision (Reason Codes MS/MSD and RPD)

The MS and MSD samples are a portion of a field sample spiked with target analytes that are prepared with each analytical batch and method as appropriate. The MS/MSD results are used to evaluate any bias introduced to the method due to matrix interference, and to measure bias and precision for each analytical batch.

In accordance with the QAPjP requirements, the additional volume for MS/MSD samples was collected at a rate of one per 20 soil samples. During the source zone coring sampling event, five MS/MSD samples were collected associated with the 73 VOC samples and 97 TPH samples and analyzed for the same analytical parameters as identified on Appendix H – Table 1. MS/MSD sample collection was achieved at the one per 20 samples requirement for the sampling event. Additional MS/MSD samples were analyzed by the laboratory as necessary to meet method and analytical batch requirements.

Sample-specific MS/MSD analyte recoveries exceeded control criteria for VOC analytes, EDB, and TPH-GRO in soil samples. As a result, the analytes in the associated samples were qualified “J” or “UJ” signifying estimated detect and non-detect data. All other MS/MSD recoveries were within method control criteria. Qualified source zone sample data based on MS/MSD sample recoveries are presented on Appendix H – Table 3.

1.1.7 Initial and Continuing Calibration Verification (Reason Code CCV)

Instrument calibration is performed for all analyses in accordance with method requirements. The linear analytical range is established for each method by analysis of calibration standards prepared at increasing concentrations that cover the expected sample concentration range. The acceptability of the initial calibration is determined by calculation of a percent relative standard deviation or coefficient. Routinely during sample analysis, the stability of the analytical system is monitored by analysis of continuing calibration standards at concentrations near the mid-point of the instrument calibration range. The percent difference values between the relative response factor in the initial calibration and the relative response factor in the continuing calibration are reviewed to ensure instrument calibration criteria are within method control limits. Initial and continuing calibration verification met the method-specific control criteria for the source zone analytical data with exception of the VOC analytes n-butylbenzene in three SDGs and trans-1,3-dichloropropene in one SDG. Qualified source zone sample data based on continuing calibration verification (CCV) criteria are presented on Appendix H – Table 3.

1.1.8 Sample Confirmation (Reason Code RPD)

As required by DoD and EPA analytical method guidance, sample detections for EDB require confirmation using a second column analysis. EDB samples were analyzed by EPA Method SW8011, confirmed by a second column analysis, and reported from the primary column. Any detection of EDB on the second column is considered confirmation unless it appears to be associated with matrix interferences. There were no data reporting issues associated with EDB detections and second column confirmation for the source zone sampling event.

1.1.9 Trip Blanks (Reason Code TB)

Trip blanks were prepared by the laboratory and stored with the containers to be used for collection of samples to be analyzed for VOCs, EDB, and TPH-GRO. As samples are collected for volatile-type analyses, trip blanks are placed in the cooler with the sample containers; therefore, they are exposed to any potential contamination along with the field samples. In accordance with the QAPjP requirements, trip blank samples are to be included at a rate of one per cooler when collecting samples for volatile organic-type analyses. A total of 46 trip blank samples were shipped during the source zone coring sampling event accompanying samples for VOC, EDB, and TPH-GRO analyses. Detections of acetone, methylene chloride, and TPH-GRO were reported in trip blanks resulting in “U” qualified (non-detect) sample results. Qualified source zone sample data based on trip blank detections are presented on Appendix H – Table 3. Appendix H – Table 4 summarizes the results for trip blank samples for the source zone sampling event.

1.1.10 Equipment Rinse Blanks (Reason Code RB)

RBs are collected in the field to assess potential contamination from sampling equipment. Results for the RB samples are used to evaluate the efficiency of equipment decontamination procedures in the field.

Eighteen RB samples were collected during the source zone coring sampling event. The RB samples were prepared by rinsing the decontaminated sampling equipment with laboratory-grade water provided by Environmental Sampling Supply, and then collecting the final rinse water into appropriate sample containers for analysis. The RB and source water samples associated with the source zone coring sampling event were analyzed for VOCs, EDB, and TPH. Low-level detections of TPH and VOCs were reported in RBs and the source water blank but only TPH-GRO detections resulted in qualified sample data. The associated detections of TPH-GRO in three samples were “U” qualified, signifying non-detect data. The RB sample results were acceptable to demonstrate the equipment decontamination procedures for the source zone coring sampling event achieved project objectives. Qualified source zone sample data based on RB detections are presented on Appendix H – Table 3. Appendix H – Table 4 presents the results of the RB and source water blank samples.

1.1.11 Field Duplicate Samples (Reason Code FD)

In accordance with the project QAPjP requirements (USACE, 2017b), field duplicate samples are collected at a frequency of 10% of the total number of soil samples for the source zone coring event. Ten field duplicate samples were collected for a total of 73 VOC soil samples and 87 TPH soil samples. The 10 % frequency for field duplicate samples was achieved for source zone sampling at 14% and 12%, respectively.

Field duplicate RPD is evaluated by calculating the RPD between the parent sample and the duplicate sample. The RPD was calculated using the following equation:

$$RPD = |(S-D)/[(S+D)/2]| \times 100$$

where

- S = Sample result.
- D = Duplicate result.

Acceptable precision control criteria are established at less than or equal to 50% for soil samples as specified in the QAPjP. The RPD was calculated between pairs of field duplicate samples when both results are reported at or above the LOQ. Field duplicate sample results that exceeded the RPD and resulted in qualified data include VOCs, EDB, and TPH-GRO. The analytes exceeding RPD were “J” qualified, signifying estimated data. Qualified source zone sample data based on field duplicates are presented on Appendix H – Table 3. All other field duplicate samples achieved the acceptable precision control criterion. The results for the field duplicate samples collected during the source zone coring event are presented on Appendix H – Table 5. The field duplicate results demonstrate acceptable overall field sampling procedures and analytical method precision.

1.1.12 Professional Judgement

Professional judgement may be applied by a third-party data validation subcontractor or the project chemist during the data review process to apply validation qualifiers based on site-specific knowledge, historical data, comparability of data, and analytical expertise. There were no exceptions to the validation qualifiers as applied to the data in accordance with the project QAPjP and other guidelines used in the validation review for the source zone coring data.

1.2 COMPLETENESS

The following sections present a discussion of contractual, analytical, and technical completeness for the source zone coring analytical data. Completeness calculations were performed for the samples that are used for project decisions. Completeness results are presented in the following sections.

1.2.1 Contractual Completeness

Contractual completeness is a quantitative determination of the number of unqualified results compared to the total number of sample results expressed as a percentage, based on data qualified for QC outliers related to analytical method performance. These include data qualified for calibration or method blank contamination, missed holding times, LCS recovery, and/or precision. The contractual completeness goal is 95%. Contractual completeness was calculated as follows:

$$\text{Percent Contractual Completeness} = \frac{\text{Number of Unqualified Results}}{\text{Total Number of Results}} \times 100$$

Source zone sampling data were qualified based on calibration verification, method blank detections, and holding time exceedance (234 qualified analytes out of the 5,852 total analytes). Contractual completeness for the source zone coring data is 96%. The 95% contractual completeness objective was achieved for all of the methods for the source zone coring sampling event.

1.2.2 Analytical Completeness

Analytical completeness is a quantitative measure of the number of unqualified data results compared to the total number of results expressed as a percentage, based on the target analytes qualified for exceedances of QC requirements for calibration, LCS, MS/MSD, surrogate, method precision, and laboratory method blank contamination results. The analytical completeness goal is 90% for the project. Analytical completeness was calculated as follows:

$$\text{Percent Analytical Completeness} = \frac{\text{Number of Unqualified Results}}{\text{Total Number of Results}} \times 100$$

A total of 595 soil sample results were qualified “U” and “J” based on method blank, trip blank, RB, CCV, surrogate recovery, MS/MSD, field duplicate RPD, and hold time, out of a total number of 5,852 analyte results, for an analytical completeness of 90%. The analytical completeness objective was achieved for the source zone coring sampling event.

Estimated data (“J” qualified) are usable to achieve project objectives. No analytical data for the source zone coring samples were rejected (“R” qualified) and qualified unusable based on analytical completeness.

1.2.3 Technical Completeness

Technical completeness is a quantitative measure of the data usability based on the number of rejected data compared to the total number of sample results. The technical completeness goal for the source zone coring event is equal to or greater than 95%. The technical completeness calculation considers all data that are not rejected (“R” qualified) usable data to achieve project objectives. The technical completeness was calculated as follows:

$$\text{Percent Technical Completeness} = \frac{\text{Number of Usable Results}}{\text{Total Number of Results}} \times 100$$

The project data quality objectives were achieved for all methods for the sampling event. No data were rejected, and the technical completeness for the soil data is 100% for all parameters and is provided in Appendix H – Table 6.

1.2.4 Data Analysis Completeness

As a part of the data review process, chain-of-custody forms and project data deliverables are reviewed against the Work Plan Table 3-3 – Proposed Sampling Plan for Continuous Coring to ensure compliance with the sampling plan, and that analytical results were reported for all planned methods and samples with the laboratory receipt exceptions noted in Section 1.1.1.

Data deliverable completeness for the source zone coring data deliverables was determined to be 100% complete as received from the laboratory. Level 4 analytical data packages are provided in Appendix G-1.

1.3 REPRESENTATIVENESS AND COMPARABILITY

Source zone characterization sampling was conducted in accordance with the sampling and analysis protocols and standard operating procedures documented in the Work Plan (USACE, 2017a). Approved procedures were used to collect, preserve, document, and ship samples to the TA laboratory, thus ensuring the samples collected for the coring sampling event were representative of the project site and conditions.

The project laboratory (TA) maintains current DoD Environmental Laboratory Accreditation Program certification and adhered to the analytical methods documented in the project QAPjP and DoD QSM requirements to prepare and analyze samples and report the data. This certification ensures the

comparability of the analytical results between different samples. For the source zone coring data, an EPA Stage 3 data review was performed on 100% of the analytical data to verify that the laboratory complied with the DoD QSM, project QAPjP, and method requirements. QC results that exceeded method control criteria resulted in data qualification as presented in the previous sections. Based on a review of the completed sample collection logs, chain-of-custody forms, sample receipt forms, and laboratory data packages, the analytical data reported for the source zone coring event achieved the project data representativeness and comparability requirements.

1.4 SENSITIVITY

Data sensitivity for source zone coring was achieved by complying with the analytical method guidelines and RLs specified in the project QAPjP and in accordance with DoD QSM requirements. Elevated sample RLs are associated with elevated concentrations of target analytes or sample matrix interference in the soil samples, resulting in sample dilution. For the analytical results, detections of target compounds reported below the LOQ are “J” flagged as estimated values. Non-detect analytes are reported at the LOD per the DoD QSM requirements unless as noted above.

1.5 CONCLUSIONS

The analytical data reported for the source zone coring sampling event have been reviewed for precision, accuracy (bias), representativeness, comparability, completeness, and sensitivity. Data quality exceedances that resulted in data qualification include: (1) method blank contamination for VOCs and TPH-GRO/DRO; (2) RB contamination for TPH-GRO; (3) trip blank contamination for VOCs and TPH-GRO; (4) MS/MSD recovery exceedances for VOCs, EDB, and TPH-GRO; (5) surrogate recovery exceedance for VOCs and TPH-GRO/DRO; (6) CCV criteria exceedance for two VOCs analytes; (7) holding time exceedances for VOCs, TPH-GRO/DRO, and EDB; and (8) field duplicate RPD exceedance for VOCs, EDB, and TPH-GRO. Analytical data are qualified as estimated detect and non-detect data results. Estimated sample data are usable to achieve project objectives. The 95% technical completeness goal was achieved for all analytical methods for the source zone coring sampling event. Data are usable to achieve the project data quality objectives as qualified based on validation.

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TABLES

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Table H-1
Source Zone Characterization Sample Collection Summary

Sample Location ID	Field Sample ID	Sample Date	Sample Delivery Group	Analytical Parameter ^a	Comments
KAFB-106247	P-247-050219-480	2/5/2019	280-119877-1	TPH	
KAFB-106247	P-247-050219-489	2/5/2019	280-119877-1	TPH	
KAFB-106247	P-247-050219-499	2/5/2019	280-119877-1	TPH	
KAFB-106247	V-247-050219-480	2/5/2019	280-119877-1	EDB, VOCs	
KAFB-106247	V-247-050219-489	2/5/2019	280-119877-1	EDB, VOCs	
KAFB-106247	V-247-050219-499	2/5/2019	280-119877-1	EDB, VOCs	
KAFB-106247	P-247-310119-164	1/31/2019	280-119743-1	TPH	
KAFB-106247	P-247-310119-208	1/31/2019	280-119743-1	TPH	
KAFB-106247	V-247-310119-164	1/31/2019	280-119743-1	EDB, VOCs	
KAFB-106247	V-247-310119-208	1/31/2019	280-119743-1	EDB, VOCs	
KAFB-106247	P-247-300119-120	1/30/2019	280-119694-1	TPH	
KAFB-106247	V-247-300119-120	1/30/2019	280-119694-1	EDB, VOCs	
KAFB-106S1	P-S1-271118-140	11/27/2018	280-117531-1	TPH	
KAFB-106S1	P-S1-291118-260	11/29/2018	280-117531-1	TPH	
KAFB-106S1	P-S1-291118-279	11/29/2018	280-117531-1	TPH	
KAFB-106S1	V-S1-271118-140	11/27/2018	280-117531-1	EDB, VOCs	
KAFB-106S1	V-S1-291118-260	11/29/2018	280-117531-1	EDB, VOCs	
KAFB-106S1	V-S1-291118-279	11/29/2018	280-117531-1	EDB, VOCs	
KAFB-106S1	P-S1-041218-459	12/4/2018	280-117710-1	TPH	
KAFB-106S1	P-S1-041218-461	12/4/2018	280-117710-1	TPH	
KAFB-106S1	V-S1-041218-459	12/4/2018	280-117710-1	EDB, VOCs	
KAFB-106S1	V-S1-041218-461	12/4/2018	280-117710-1	EDB, VOCs	
KAFB-106S1	P-S1-011218-400	12/1/2018	280-117650-1	TPH	
KAFB-106S1	P-S1-031218-414	12/3/2018	280-117650-1	TPH	
KAFB-106S1	V-S1-011218-400	12/1/2018	280-117650-1	EDB, VOCs	
KAFB-106S1	V-S1-031218-414	12/3/2018	280-117650-1	EDB, VOCs	
KAFB-106S1	P-S1-051218-510	12/5/2018	280-117794-1	TPH	Field Duplicate
KAFB-106S1	V-S1-051218-510	12/5/2018	280-117794-1	EDB, VOCs	Field Duplicate
KAFB-106S1	P-S1-271118-020	11/27/2018	280-117391-1	TPH	
KAFB-106S1	P-S1-051218-475	12/5/2018	280-117883-1	TPH	
KAFB-106S1	P-S1-051218-480	12/5/2018	280-117883-1	TPH	
KAFB-106S1	P-S1-051218-489	12/5/2018	280-117883-1	TPH	
KAFB-106S1	V-S1-051218-475	12/5/2018	280-117883-1	EDB, VOCs	
KAFB-106S1	V-S1-051218-480	12/5/2018	280-117883-1	EDB, VOCs	
KAFB-106S1	V-S1-051218-489	12/5/2018	280-117883-1	EDB, VOCs	
KAFB-106S2	P-S2-061118-050	11/6/2018	280-116903-1	TPH	
KAFB-106S2	P-S2-061118-105	11/6/2018	280-116903-1	TPH	
KAFB-106S2	P-S2-081118-278	11/8/2018	280-116903-1	TPH	
KAFB-106S2	P-S2-131118-338	11/13/2018	280-117064-1	TPH	
KAFB-106S2	P-S2-141118-404	11/14/2018	280-117064-1	TPH	
KAFB-106S2	P-S2-151118-419	11/15/2018	280-117064-1	TPH	
KAFB-106S2	V-S2-131118-338	11/13/2018	280-117064-1	EDB, VOCs	
KAFB-106S2	V-S2-141118-404	11/14/2018	280-117064-1	EDB, VOCs	
KAFB-106S2	V-S2-151118-419	11/15/2018	280-117064-1	EDB, VOCs	
KAFB-106S2	P-S2-161118-510	11/16/2018	280-117130-1	TPH	Field Duplicate, MS/MSD
KAFB-106S2	V-S2-161118-510	11/16/2018	280-117130-1	EDB, VOCs	Field Duplicate, MS/MSD
KAFB-106S2	P-S2-161118-474	11/16/2018	280-117130-2	TPH	

Table H-1
Source Zone Characterization Sample Collection Summary

Sample Location ID	Field Sample ID	Sample Date	Sample Delivery Group	Analytical Parameter ^a	Comments
KAFB-106S2	P-S2-161118-499	11/16/2018	280-117130-2	TPH	
KAFB-106S2	V-S2-161118-474	11/16/2018	280-117130-2	EDB, VOCs	
KAFB-106S2	V-S2-161118-499	11/16/2018	280-117130-2	EDB, VOCs	
KAFB-106S3	P-S3-191018-020	10/19/2018	280-115886-1	TPH	
KAFB-106S3	P-S3-191018-160	10/19/2018	280-115886-1	TPH	
KAFB-106S3	P-S3-051118-240	11/4/2018	280-116615-3	TPH	
KAFB-106S3	P-S3-051118-360	11/5/2018	280-116615-3	TPH	
KAFB-106S3	P-S3-191118-424	11/19/2018	280-117172-1	TPH	
KAFB-106S3	V-S3-191118-424	11/19/2018	280-117172-1	EDB, VOCs	
KAFB-106S3	P-S3-201118-459	11/20/2018	280-117258-1	TPH	Field Duplicate, MS/MSD
KAFB-106S3	V-S3-201118-459	11/20/2018	280-117258-1	EDB, VOCs	Field Duplicate, MS/MSD
KAFB-106S3	P-S3-201118-477	11/20/2018	280-117265-1	TPH	
KAFB-106S3	P-S3-201118-489	11/20/2018	280-117265-1	TPH	
KAFB-106S3	V-S3-201118-477	11/20/2018	280-117265-1	EDB, VOCs	
KAFB-106S3	V-S3-201118-489	11/20/2018	280-117265-1	EDB, VOCs	
KAFB-106S3	P-S3-211118-492	11/21/2018	280-117359-1	TPH	
KAFB-106S3	P-S3-211118-512	11/21/2018	280-117359-1	TPH	
KAFB-106S3	V-S3-211118-492	11/21/2018	280-117359-1	EDB, VOCs	
KAFB-106S3	V-S3-211118-512	11/21/2018	280-117359-1	EDB, VOCs	
KAFB-106S4	P-S4-171018-040	10/17/2018	280-115762-1	TPH	
KAFB-106S4	P-S4-171018-110	10/17/2018	280-115762-1	TPH	
KAFB-106S4	P-S4-171018-300	10/17/2018	280-115762-1	TPH	
KAFB-106S4	P-S4-011118-366	11/1/2018	280-116535-1	TPH	
KAFB-106S4	P-S4-021118-416	11/2/2018	280-116535-1	TPH	
KAFB-106S4	V-S4-021118-416	11/2/2018	280-116535-1	EDB, VOCs	
KAFB-106S4	P-S4-041118-467	11/4/2018	280-116615-1	TPH	
KAFB-106S4	P-S4-051118-494	11/5/2018	280-116615-1	TPH	
KAFB-106S4	V-S4-041118-467	11/4/2018	280-116615-1	EDB, VOCs	
KAFB-106S4	P-S4-051118-504	11/5/2018	280-116615-2	TPH	Field Duplicate, MS/MSD
KAFB-106S4	V-S4-051118-504	11/5/2018	280-116615-2	EDB, VOCs	Field Duplicate, MS/MSD
KAFB-106S4	V-S4-051118-494	11/5/2018	280-116615-1	EDB, VOCs	
KAFB-106S5	P-S5-091018-210	10/9/2018	280-115297-1	TPH	
KAFB-106S5	P-S5-091018-360	10/9/2018	280-115297-1	TPH	
KAFB-106S5	P-S5-211018-417	10/21/2018	280-115956-1	TPH	
KAFB-106S5	V-S5-211018-417	10/21/2018	280-115956-1	EDB, VOCs	
KAFB-106S5	P-S5-231018-467	10/23/2018	280-116044-1	TPH	Field Duplicate, MS/MSD
KAFB-106S5	P-S5-231018-491	10/23/2018	280-116044-1	TPH	
KAFB-106S5	P-S5-231018-506	10/23/2018	280-116044-1	TPH	
KAFB-106S5	V-S5-231018-467	10/23/2018	280-116044-1	EDB, VOCs	Field Duplicate, MS/MSD
KAFB-106S5	V-S5-231018-491	10/23/2018	280-116044-1	EDB, VOCs	
KAFB-106S5	V-S5-231018-506	10/23/2018	280-116044-1	EDB, VOCs	
KAFB-106S7	P-S7-220119-469	1/22/2019	280-119401-1	TPH	
KAFB-106S7	P-S7-220119-485	1/22/2019	280-119401-1	TPH	

Table H-1
Source Zone Characterization Sample Collection Summary

Sample Location ID	Field Sample ID	Sample Date	Sample Delivery Group	Analytical Parameter^a	Comments
KAFB-106S7	P-S7-220119-495	1/22/2019	280-119401-1	TPH	
KAFB-106S7	P-S7-220119-506	1/22/2019	280-119401-1	TPH	Field Duplicate
KAFB-106S7	V-S7-220119-469	1/22/2019	280-119401-1	EDB, VOCs	
KAFB-106S7	V-S7-220119-485	1/22/2019	280-119401-1	EDB, VOCs	
KAFB-106S7	V-S7-220119-495	1/22/2019	280-119401-1	EDB, VOCs	
KAFB-106S7	V-S7-220119-506	1/22/2019	280-119401-1	EDB, VOCs	Field Duplicate
KAFB-106S7	P-S7-210119-420	1/21/2019	280-119377-1	TPH	
KAFB-106S7	V-S7-210119-420	1/21/2019	280-119377-1	EDB, VOCs	
KAFB-106S8	P-S8-160119-419	1/16/2019	280-119255-1	TPH	
KAFB-106S8	V-S8-160119-419	1/16/2019	280-119255-1	EDB, VOCs	
KAFB-106S8	P-S8-180119-475	1/18/2019	280-119347-1	TPH	
KAFB-106S8	P-S8-180119-499	1/18/2019	280-119347-1	TPH	
KAFB-106S8	P-S8-180119-514	1/18/2019	280-119347-1	TPH	Field Duplicate
KAFB-106S8	V-S8-180119-475	1/18/2019	280-119347-1	EDB, VOCs	
KAFB-106S8	V-S8-180119-499	1/18/2019	280-119347-1	EDB, VOCs	
KAFB-106S8	V-S8-180119-514	1/18/2019	280-119347-1	EDB, VOCs	Field Duplicate
KAFB-106S9	P-S9-061018-102	10/6/2018	280-115294-1	TPH	
KAFB-106S9	P-S9-071018-174	10/7/2018	280-115294-1	TPH	
KAFB-106S9	V-S9-061018-102	10/6/2018	280-115294-1	EDB, VOCs	
KAFB-106S9	V-S9-071018-174	10/7/2018	280-115294-1	EDB, VOCs	
KAFB-106S9	P-S9-111018-342	10/11/2018	280-115488-1	TPH	
KAFB-106S9	V-S9-111018-342	10/11/2018	280-115488-1	EDB, VOCs	
KAFB-106S9	P-S9-161018-415	10/16/2018	280-115708-1	TPH	
KAFB-106S9	V-S9-161018-415	10/16/2018	280-115708-1	EDB, VOCs	
KAFB-106S9	P-S9-171018-470	10/17/2018	280-115813-1	TPH	
KAFB-106S9	P-S9-171018-475	10/17/2018	280-115813-1	TPH	
KAFB-106S9	V-S9-171018-470	10/17/2018	280-115813-1	EDB, VOCs	
KAFB-106S9	V-S9-171018-475	10/17/2018	280-115813-1	EDB, VOCs	
KAFB-106S9	P-S9-301018-496	10/30/2018	280-116409-1	TPH	
KAFB-106S9	P-S9-301018-501	10/30/2018	280-116409-1	TPH	
KAFB-106S9	V-S9-301018-496	10/30/2018	280-116409-1	EDB, VOCs	
KAFB-106S9	V-S9-301018-501	10/30/2018	280-116409-1	EDB, VOCs	
KAFB-106S9	P-S9-091018-252	10/9/2018	280-115432-1	TPH	
KAFB-106S9	V-S9-091018-252	10/9/2018	280-115432-1	EDB, VOCs	
KAFB-106S9	P-S9-191018-490	10/19/2018	280-115896-1	TPH	Field Duplicate, MS/MSD
KAFB-106S9	V-S9-191018-490	10/19/2018	280-115896-1	EDB, VOCs	Field Duplicate, MS/MSD
KAFB-106V1	P-V1-161218-061	12/16/2018	280-118336-1	TPH	
KAFB-106V1	P-V1-161218-101	12/16/2018	280-118336-1	TPH	
KAFB-106V1	P-V1-161218-115	12/16/2018	280-118336-1	TPH	
KAFB-106V1	P-V1-171218-147	12/17/2018	280-118336-1	TPH	
KAFB-106V1	V-V1-161218-061	12/16/2018	280-118336-1	EDB, VOCs	
KAFB-106V1	V-V1-161218-101	12/16/2018	280-118336-1	EDB, VOCs	
KAFB-106V1	V-V1-161218-115	12/16/2018	280-118336-1	EDB, VOCs	
KAFB-106V1	V-V1-171218-147	12/17/2018	280-118336-1	EDB, VOCs	
KAFB-106V1	P-V1-171218-161	12/17/2018	280-118390-1	TPH	
KAFB-106V1	P-V1-181218-216	12/18/2018	280-118390-1	TPH	
KAFB-106V1	V-V1-171218-161	12/17/2018	280-118390-1	EDB, VOCs	

Table H-1
Source Zone Characterization Sample Collection Summary

Sample Location ID	Field Sample ID	Sample Date	Sample Delivery Group	Analytical Parameter ^a	Comments
KAFB-106V1	V-V1-181218-216	12/18/2018	280-118390-1	EDB, VOCs	
KAFB-106V1	P-V1-191218-254	12/19/2018	280-118473-1	TPH	
KAFB-106V1	P-V1-191218-271	12/19/2018	280-118473-1	TPH	
KAFB-106V1	P-V1-191218-285	12/19/2018	280-118473-1	TPH	Field Duplicate
KAFB-106V1	V-V1-191218-254	12/19/2018	280-118473-1	EDB, VOCs	
KAFB-106V1	V-V1-191218-271	12/19/2018	280-118473-1	EDB, VOCs	
KAFB-106V1	V-V1-191218-285	12/19/2018	280-118473-1	EDB, VOCs	Field Duplicate
KAFB-106V2	P-V2-121218-080	12/12/2018	280-118266-1	TPH	
KAFB-106V2	P-V2-121218-103	12/12/2018	280-118266-1	TPH	
KAFB-106V2	P-V2-131218-117	12/13/2018	280-118266-1	TPH	
KAFB-106V2	P-V2-131218-159	12/13/2018	280-118266-1	TPH	
KAFB-106V2	V-V2-121218-080	12/12/2018	280-118266-1	EDB, VOCs	
KAFB-106V2	V-V2-121218-103	12/12/2018	280-118266-1	EDB, VOCs	
KAFB-106V2	V-V2-131218-117	12/13/2018	280-118266-1	EDB, VOCs	
KAFB-106V2	V-V2-131218-159	12/13/2018	280-118266-1	EDB, VOCs	
KAFB-106V2	P-V2-141218-215	12/14/2018	280-118290-1	TPH	
KAFB-106V2	P-V2-141218-254	12/14/2018	280-118290-1	TPH	
KAFB-106V2	V-V2-141218-215	12/14/2018	280-118290-1	EDB, VOCs	
KAFB-106V2	V-V2-141218-254	12/14/2018	280-118290-1	EDB, VOCs	
KAFB-106V2	P-V2-141218-270	12/14/2018	280-118291-1	TPH	Field Duplicate
KAFB-106V2	V-V2-141218-270	12/14/2018	280-118291-1	EDB, VOCs	Field Duplicate
KAFB-106V2	P-V2-151218-287	12/15/2018	280-118337-1	TPH	
KAFB-106V2	V-V2-151218-287	12/15/2018	280-118337-1	EDB, VOCs	
Trip Blank	TB-S9-091018-01	10/9/2018	280-115294-1	TPH, VOCs	
Trip Blank	TB-S5-091018-01	10/9/2018	280-115297-1	TPH	
Trip Blank	TB-S9-111018-01	10/11/2018	280-115488-1	TPH, VOCs	
Trip Blank	TB-S9-161018-01	10/16/2018	280-115708-1	TPH, VOCs	
Trip Blank	TB-S4-171018-01	10/17/2018	280-115762-1	TPH	
Trip Blank	TB-S9-181018-01	10/18/2018	280-115813-1	TPH, VOCs	
Trip Blank	TB-S3-191018-01	10/19/2018	280-115886-1	TPH	
Trip Blank	TB-S4-021118-01	11/2/2018	280-116535-1	TPH, EDB, VOCs	
Trip Blank	TB-S4-041118-01	11/4/2018	280-116615-1	TPH, EDB, VOCs	
Trip Blank	TB-S4-051118-01	11/5/2018	280-116615-2	TPH, EDB, VOCs	
Trip Blank	TB-S3-051118-01	11/5/2018	280-116615-3	TPH	
Trip Blank	TB-S5-221018-01	10/22/2018	280-115956-1	TPH, VOCs	
Trip Blank	TB-S9-221018-01	10/22/2018	280-115951-1	TPH, EDB, VOCs	
Trip Blank	TB-S9-311018-01	10/31/2018	280-116409-1	TPH, EDB, VOCs	
Trip Blank	TB-S2-061118-01	11/6/2018	280-116903-1	TPH	
Trip Blank	TB-S9-101018-01	10/10/2018	280-115432-1	TPH, VOCs	
Trip Blank	TB-S5-231018-01	10/23/2018	280-116044-1	TPH, VOCs	
Trip Blank	TB-S5-231018-02	10/23/2018	280-116044-1	TPH	
Trip Blank	TB-S9-191018-01	10/19/2018	280-115896-1	TPH, VOCs	
Trip Blank	TB-S2-151118-01	11/15/2018	280-117064-1	TPH, EDB, VOCs	
Trip Blank	TB-S2-161118-02	11/16/2018	280-117130-1	TPH, EDB, VOCs	
Trip Blank	TB-S2-161118-01	11/16/2018	280-117130-2	TPH, EDB, VOCs	
Trip Blank	TB-S3-191118-01	11/19/2018	280-117172-1	TPH, EDB, VOCs	
Trip Blank	TB-S3-2018-01	11/20/2018	280-117258-1	TPH, EDB, VOCs	
Trip Blank	TB-S3-201118-02	11/20/2018	280-117265-1	TPH, EDB, VOCs	
Trip Blank	TB-S3-211118-01	11/21/2018	280-117359-1	TPH, EDB, VOCs	

Table H-1
Source Zone Characterization Sample Collection Summary

Sample Location ID	Field Sample ID	Sample Date	Sample Delivery Group	Analytical Parameter ^a	Comments
Trip Blank	TB-S1-291118-01	11/29/2018	280-117531-1	TPH, EDB, VOCs	
Trip Blank	TB-S1-041218-01	12/4/2018	280-117710-1	TPH, EDB, VOCs	
Trip Blank	TB-S1-031218-01	12/3/2018	280-117650-1	TPH, EDB, VOCs	
Trip Blank	TB-S1-051218-01	12/5/2018	280-117883-1	TPH, VOCs	
Trip Blank	TB-S1-051218-02	12/5/2018	280-117794-1	TPH, EDB, VOCs	
Trip Blank	TB-S1-271118-01	11/27/2018	280-117391-1	TPH	
Trip Blank	TB-V1-171218-01	12/17/2018	280-118336-1	TPH, EDB, VOCs	
Trip Blank	TB-V1-181218-01	12/18/2018	280-118390-1	TPH, EDB, VOCs	
Trip Blank	TB-V2-131218-01	12/13/2018	280-118266-1	TPH, EDB, VOCs	
Trip Blank	TB-V2-141218-01	12/14/2018	280-118290-1	TPH, EDB, VOCs	
Trip Blank	TB-V2-141218-02	12/14/2018	280-118291-1	TPH, VOCs	
Trip Blank	TB-V2-171218-01	12/17/2018	280-118337-1	TPH, EDB, VOCs	
Trip Blank	TB-V1-191218-01	12/19/2018	280-118473-1	TPH, EDB, VOCs	
Trip Blank	TB-S8-160119-01	1/16/2019	280-119255-1	TPH, EDB, VOCs	
Trip Blank	TB-S8-180119-01	1/18/2019	280-119347-1	TPH, EDB, VOCs	
Trip Blank	TB-247-050219-01	2/5/2019	280-119877-1	TPH, EDB, VOCs	
Trip Blank	TB-247-310119-01	1/31/2019	280-119743-1	TPH, EDB, VOCs	
Trip Blank	TB-247-300119-01	1/30/2019	280-119694-1	TPH, EDB, VOCs	
Trip Blank	TB-S7-220119-01	1/22/2019	280-119401-1	TPH, EDB, VOCs	
Trip Blank	TB-S7-210119-01	1/21/2019	280-119377-1	TPH, EDB, VOCs	
Rinsate Blank	RB-247-290119-01	1/29/2019	280-119694-1	VOCs	
Rinsate Blank	RB-247-310119-01	1/31/2019	280-119743-1	EDB, VOCs, TPH	
Rinsate Blank	RB-S2-181118-01	11/18/2018	280-117169-1	EDB, VOCs, TPH	
Rinsate Blank	RB-S2-181118-02	11/18/2018	280-117169-1	EDB, VOCs, TPH	
Rinsate Blank	RB-S3-271118-01	11/27/2018	280-117400-1	EDB, VOCs, TPH	
Rinsate Blank	RB-S3-271118-02	11/27/2018	280-117400-1	EDB, VOCs, TPH	
Rinsate Blank	RB-S4-051118-01	11/5/2018	280-116899-1	EDB, VOCs, TPH	
Rinsate Blank	RB-S4-051118-02	11/5/2018	280-116899-1	EDB, VOCs, TPH	
Rinsate Blank	RB-S5-251018-01	10/25/2018	280-116206-1	EDB, VOCs, TPH	
Rinsate Blank	RB-S5-251018-02	10/25/2018	280-116206-1	EDB, VOCs, TPH	
Rinsate Blank	RB-S7-200119-01	1/20/2019	280-119377-1	EDB, VOCs, TPH	
Rinsate Blank	RB-S9-011118-01	11/1/2018	280-116567-1	EDB, VOCs, TPH	
Rinsate Blank	RB-S9-011118-02	11/1/2018	280-116567-1	EDB, VOCs, TPH	
Rinsate Blank	RB-S9-201018-01	10/20/2018	280-115951-1	EDB, VOCs, TPH	
Rinsate Blank	RB-V2-111218-01	12/11/2018	280-118241-1	EDB, VOCs, TPH	
Rinsate Blank	RB-S8-060119-01	1/6/2019	280-119077-1	EDB, VOCs, TPH	
Rinsate Blank	RB-S9-201018-02	10/20/2018	280-115951-1	EDB, VOCs, TPH	
Rinsate Blank	RB-V1-151218-01	12/15/2018	280-118337-1	EDB, VOCs, TPH	
Source Water Blank	SWB-S9-201018-01	10/20/2018	280-115951-1	EDB, VOCs, TPH	

a. Analytical methods include: Method SW8260B for VOCs; Method SW8011 for EDB; and Method SW8015D modified for TPH-gasoline range organics, diesel range organics, and motor oil. Trip blanks were analyzed for VOCs, EDB, and TPH-gasoline range organics.

EDB = ethylene dibromide

ID = identification

MS = matrix spike

MSD = matrix spike duplicate

TPH = total petroleum hydrocarbons

VOC = volatile organic compound

Table H-2
Data Qualification Flags and Reason Codes

Data Qualifier Definitions for Data Validation

Qualifier	Definition
	No Qualifier indicates that the data are acceptable both qualitatively and quantitatively.
U	The analyte was analyzed for but was not detected above the detection limit. The value associated with the U-qualifier is the limit of detection.
J	The analyte was analyzed for and was positively identified, but the reported numerical value may not be consistent with the amount actually present in the environmental sample. Results are estimated, although the data are considered usable and may be used as appropriate to meet project objectives. Results are qualitatively acceptable and quantitatively uncertain.
J-	The analyte was positively identified; the associated numerical value is its approximate concentration with a low bias in the sample.
J+	The analyte was positively identified; the associated numerical value is its approximate concentration with a high bias in the sample.
UJ	The analyte was analyzed for but was not detected. The reported quantitation limit is approximate and may be inaccurate or imprecise.
R	The analyte was analyzed for, but the presence <u>or</u> absence of the analyte has not been verified. Re-sampling and re-analysis may be necessary to confirm or deny the presence of the analyte. Results are rejected, and data are <u>unusable</u> for any purposes.

Reason Codes for Data Validation

Reason Code	Description
CB/CCB	Calibration blank or continuing calibration blank outside of control limits
CCV	Calibration verification outside of control limits
EB	Equipment rinse blank contamination
FB	Field blank contamination
FD	Field duplicate sample results out of control criteria
HT	Holding time exceedance
ICS	Interference check sample
LCS	Laboratory control sample recovery out of control criteria
MB	Method blank contamination
MS/MSD	Matrix spike/ matrix spike duplicate recovery outside of control criteria
RPD	Relative percent difference outside of control limits
SD	Inductively coupled plasma serial dilution out of control criteria
SURR	Surrogate recovery outside of control limits
TB	Trip blank contamination

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
KAFB-106247	V-247-050219-480	280-119877-1	G	N	Chloroform	U	Method Blank
KAFB-106247	V-247-050219-489	280-119877-1	G	N	Chloroform	U	Method Blank
KAFB-106247	V-247-050219-499	280-119877-1	G	N	Chloroform	U	Method Blank
KAFB-106247	V-247-310119-164	280-119743-1	G	N	Acetone	U	Method Blank
KAFB-106247	V-247-310119-164	280-119743-1	G	N	Chloroform	U	Method Blank
KAFB-106247	V-247-310119-208	280-119743-1	G	N	Chloroform	U	Method Blank
KAFB-106247	V-247-300119-120	280-119694-1	G	N	Acetone	U	Method Blank
KAFB-106247	V-247-300119-120	280-119694-1	G	N	Chloroform	U	Method Blank
KAFB-106S1	P-S1-271118-140	280-117531-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S1	P-S1-291118-260	280-117531-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S1	P-S1-291118-279	280-117531-1	G	N	Motor Oil (C20-C38)	UJ	Surrogate Recovery
KAFB-106S1	P-S1-291118-279	280-117531-1	G	N	TPH-DRO (C10-C28)	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	Ethylbenzene	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	Styrene	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	cis-1,3-Dichloropropene	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	trans-1,3-Dichloropropene	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	n-Propylbenzene	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	n-Butylbenzene	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	4-Chlorotoluene	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	1,4-Dichlorobenzene	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	1,2-Dibromoethane	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	1,2-Dichloroethane	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	4-Methyl-2-pentanone	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	1,3,5-Trimethylbenzene	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	Bromobenzene	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	Toluene	J	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	Chlorobenzene	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	1,2,4-Trichlorobenzene	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	Dibromochloromethane	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	Tetrachloroethene	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	sec-Butylbenzene	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	1,3-Dichloropropane	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	cis-1,2-Dichloroethene	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	trans-1,2-Dichloroethene	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	Methyl tert-butyl ether	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	m- & p-Xylenes	J	Surrogate Recovery

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	1,2-Dichloroethene	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	1,3-Dichlorobenzene	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	Carbon tetrachloride	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	1,1-Dichloropropene	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	2-Hexanone	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	2,2-Dichloropropane	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	1,1,1,2-Tetrachloroethane	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	Acetone	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	Chloroform	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	Benzene	J	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	1,1,1-Trichloroethane	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	Bromomethane	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	Chloromethane	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	Dibromomethane	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	Bromochloromethane	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	Chloroethane	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	Vinyl Chloride	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	Methylene chloride	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	Carbon disulfide	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	Bromoform	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	Bromodichloromethane	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	1,1-Dichloroethane	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	1,1-Dichloroethene	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	Trichlorofluoromethane	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	Dichlorodifluoromethane	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	1,2-Dichloropropane	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	2-Butanone	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	1,1,2-Trichloroethane	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	Trichloroethene	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	1,1,2,2-Tetrachloroethane	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	1,2,3-Trichlorobenzene	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	Hexachloro-1,3-butadiene	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	Naphthalene	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	o-Xylene	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	2-Chlorotoluene	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	1,2-Dichlorobenzene	UJ	Surrogate Recovery

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	1,2,4-Trimethylbenzene	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	1,2-Dibromo-3-chloropropane	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	1,2,3-Trichloropropane	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	tert-Butylbenzene	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	Isopropylbenzene	UJ	Surrogate Recovery
KAFB-106S1	V-S1-291118-279	280-117531-1	G	N	4-Isopropyltoluene	UJ	Surrogate Recovery
KAFB-106S1	P-S1-011218-400	280-117650-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S1	P-S1-031218-414	280-117650-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S1	V-S1-031218-414	280-117650-1	G	N	n-Butylbenzene	UJ	CCV
KAFB-106S1	P-S1-051218-510	280-117794-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S1	P-S1-051258-510	280-117794-1	G	FD	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S1	P-S1-271118-020	280-117391-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S2	P-S2-061118-050	280-116903-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S2	P-S2-061118-105	280-116903-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S2	P-S2-081118-278	280-116903-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S2	P-S2-131118-338	280-117064-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S2	P-S2-141118-404	280-117064-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S2	P-S2-151118-419	280-117064-1	G	N	Motor Oil (C20-C38)	J	Surrogate Recovery
KAFB-106S2	P-S2-151118-419	280-117064-1	G	N	TPH-DRO (C10-C28)	UJ	Surrogate Recovery
KAFB-106S2	V-S2-131118-338	280-117064-1	G	N	Acetone	U	Method Blank
KAFB-106S2	V-S2-141118-404	280-117064-1	G	N	Acetone	U	Method Blank
KAFB-106S2	V-S2-151118-419	280-117064-1	G	N	Acetone	U	Method Blank
KAFB-106S2	P-S2-161118-510	280-117130-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S2	P-S2-161118-510	280-117130-1	G	N	TPH-GRO (C6-C10)	J	Field Duplicate
KAFB-106S2	P-S2-161158-510	280-117130-1	G	FD	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S2	P-S2-161158-510	280-117130-1	G	FD	TPH-GRO (C6-C10)	J	Field Duplicate
KAFB-106S2	V-S2-161118-510	280-117130-1	G	N	Ethylbenzene	J	Field Duplicate
KAFB-106S2	V-S2-161118-510	280-117130-1	G	N	n-Propylbenzene	J	MS/MSD - Percent recovery
KAFB-106S2	V-S2-161118-510	280-117130-1	G	N	n-Butylbenzene	J	MS/MSD - Percent recovery
KAFB-106S2	V-S2-161118-510	280-117130-1	G	N	1,3,5-Trimethylbenzene	J	Field Duplicate
KAFB-106S2	V-S2-161118-510	280-117130-1	G	N	Toluene	J	Field Duplicate
KAFB-106S2	V-S2-161118-510	280-117130-1	G	N	m- & p-Xylenes	J	Field Duplicate
KAFB-106S2	V-S2-161118-510	280-117130-1	G	N	Benzene	J	Field Duplicate
KAFB-106S2	V-S2-161118-510	280-117130-1	G	N	o-Xylene	J	MS/MSD - Percent recovery
KAFB-106S2	V-S2-161118-510	280-117130-1	G	N	1,2,4-Trimethylbenzene	J	Field Duplicate
KAFB-106S2	V-S2-161158-510	280-117130-1	G	FD	Ethylbenzene	J	Field Duplicate

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
KAFB-106S2	V-S2-161158-510	280-117130-1	G	FD	1,3,5-Trimethylbenzene	J	Field Duplicate
KAFB-106S2	V-S2-161158-510	280-117130-1	G	FD	Toluene	J	Field Duplicate
KAFB-106S2	V-S2-161158-510	280-117130-1	G	FD	m- & p-Xylenes	J	Field Duplicate
KAFB-106S2	V-S2-161158-510	280-117130-1	G	FD	Benzene	J	Field Duplicate
KAFB-106S2	V-S2-161158-510	280-117130-1	G	FD	1,2,4-Trimethylbenzene	J	Field Duplicate
KAFB-106S2	P-S2-161118-499	280-117130-2	G	N	TPH-GRO (C6-C10)	J	Surrogate Recovery
KAFB-106S2	V-S2-161118-474	280-117130-2	G	N	Acetone	U	Method Blank
KAFB-106S2	V-S2-161118-474	280-117130-2	G	N	1,2,4-Trimethylbenzene	U	Method Blank
KAFB-106S3	P-S3-191018-020	280-115886-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S3	P-S3-191018-020	280-115886-1	G	N	TPH-GRO (C6-C10)	U	Method Blank
KAFB-106S3	P-S3-191018-160	280-115886-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S3	P-S3-191018-160	280-115886-1	G	N	TPH-GRO (C6-C10)	U	Method Blank
KAFB-106S3	P-S3-051118-240	280-116615-3	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S3	P-S3-051118-360	280-116615-3	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S3	P-S3-191118-424	280-117172-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S3	P-S3-201118-459	280-117258-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S3	P-S3-201158-459	280-117258-1	G	FD	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S3	V-S3-201118-459	280-117258-1	G	N	1,2-Dibromoethane	UJ	MS/MSD - Percent recovery
KAFB-106S3	V-S3-201118-459	280-117258-1	G	N	2-Butanone	J	Field Duplicate
KAFB-106S3	V-S3-201158-459	280-117258-1	G	FD	2-Butanone	J	Field Duplicate
KAFB-106S3	P-S3-201118-477	280-117265-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S3	P-S3-201118-489	280-117265-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S3	V-S3-201118-477	280-117265-1	G	N	Methylene chloride	U	Trip Blank
KAFB-106S3	V-S3-201118-489	280-117265-1	G	N	Methylene chloride	U	Trip Blank
KAFB-106S3	P-S3-211118-512	280-117359-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S3	V-S3-211118-492	280-117359-1	G	N	trans-1,3-Dichloropropene	UJ	CCV
KAFB-106S3	V-S3-211118-512	280-117359-1	G	N	trans-1,3-Dichloropropene	UJ	CCV
KAFB-106S4	P-S4-171018-040	280-115762-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S4	P-S4-171018-040	280-115762-1	G	N	TPH-GRO (C6-C10)	U	Method Blank
KAFB-106S4	P-S4-171018-110	280-115762-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S4	P-S4-171018-110	280-115762-1	G	N	TPH-GRO (C6-C10)	U	Method Blank
KAFB-106S4	P-S4-171018-300	280-115762-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S4	P-S4-171018-300	280-115762-1	G	N	TPH-GRO (C6-C10)	U	Method Blank
KAFB-106S4	P-S4-011118-366	280-116535-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S4	P-S4-021118-416	280-116535-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S4	V-S4-021118-416	280-116535-1	G	N	n-Butylbenzene	UJ	CCV

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
KAFB-106S4	P-S4-041118-467	280-116615-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S4	P-S4-051118-504	280-116615-2	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S4	P-S4-051118-504	280-116615-2	G	N	TPH-GRO (C6-C10)	UJ	MS/MSD - Percent recovery
KAFB-106S4	P-S4-051158-504	280-116615-2	G	FD	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S4	P-S4-051158-504	280-116615-2	G	FD	TPH-GRO (C6-C10)	U	Equipment Blank
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	Ethylbenzene	J	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	Styrene	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	cis-1,3-Dichloropropene	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	trans-1,3-Dichloropropene	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	n-Propylbenzene	J	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	n-Butylbenzene	J	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	4-Chlorotoluene	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	1,4-Dichlorobenzene	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	1,2-Dibromoethane	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	1,2-Dichloroethane	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	4-Methyl-2-pentanone	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	1,3,5-Trimethylbenzene	J	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	Bromobenzene	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	Toluene	J	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	Chlorobenzene	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	1,2,4-Trichlorobenzene	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	Dibromochloromethane	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	Tetrachloroethene	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	sec-Butylbenzene	J	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	1,3-Dichloropropane	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	cis-1,2-Dichloroethene	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	trans-1,2-Dichloroethene	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	Methyl tert-butyl ether	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	m- & p-Xylenes	J	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	1,2-Dichloroethene	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	1,3-Dichlorobenzene	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	Carbon tetrachloride	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	1,1-Dichloropropene	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	2-Hexanone	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	2,2-Dichloropropane	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	1,1,1,2-Tetrachloroethane	UJ	MS/MSD - Percent recovery

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	Acetone	J	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	Chloroform	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	Benzene	J	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	1,1,1-Trichloroethane	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	Bromomethane	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	Chloromethane	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	Dibromomethane	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	Bromochloromethane	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	Chloroethane	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	Vinyl Chloride	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	Methylene chloride	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	Carbon disulfide	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	Bromoform	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	Bromodichloromethane	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	1,1-Dichloroethane	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	1,1-Dichloroethene	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	Trichlorofluoromethane	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	Dichlorodifluoromethane	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	1,2-Dichloropropane	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	2-Butanone	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	1,1,2-Trichloroethane	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	Trichloroethene	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	1,1,2,2-Tetrachloroethane	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	1,2,3-Trichlorobenzene	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	Hexachloro-1,3-butadiene	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	Naphthalene	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	o-Xylene	J	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	2-Chlorotoluene	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	1,2-Dichlorobenzene	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	1,2,4-Trimethylbenzene	J	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	1,2-Dibromo-3-chloropropane	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	1,2,3-Trichloropropane	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	tert-Butylbenzene	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	Isopropylbenzene	J	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051118-504	280-116615-2	G	N	4-Isopropyltoluene	UJ	MS/MSD - Percent recovery
KAFB-106S4	V-S4-051158-504	280-116615-2	G	FD	n-Butylbenzene	J	CCV

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
KAFB-106S5	P-S5-091018-210	280-115297-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S5	P-S5-091018-210	280-115297-1	G	N	TPH-GRO (C6-C10)	UJ	MS/MSD - Percent recovery
KAFB-106S5	P-S5-091018-360	280-115297-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S5	P-S5-091018-360	280-115297-1	G	N	TPH-GRO (C6-C10)	UJ	Surrogate Recovery
KAFB-106S5	P-S5-211018-417	280-115956-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S5	P-S5-211018-417	280-115956-1	G	N	TPH-GRO (C6-C10)	UJ	Hold Time
KAFB-106S5	V-S5-211018-417	280-115956-1	G	N	Methylene chloride	U	Method Blank
KAFB-106S5	P-S5-231018-506	280-116044-1	G	N	TPH-GRO (C6-C10)	U	Trip Blank
KAFB-106S5	P-S5-231058-467	280-116044-1	G	FD	TPH-GRO (C6-C10)	UJ	Surrogate Recovery
KAFB-106S5	V-S5-231018-467	280-116044-1	G	N	Acetone	U	Method Blank
KAFB-106S5	V-S5-231018-506	280-116044-1	G	N	Acetone	U	Method Blank
KAFB-106S5	V-S5-231058-467	280-116044-1	G	FD	Acetone	U	Method Blank
KAFB-106S7	P-S7-220119-469	280-119401-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S7	P-S7-220119-506	280-119401-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S7	P-S7-220159-506	280-119401-1	G	FD	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S7	V-S7-220119-469	280-119401-1	G	N	Toluene	U	Method Blank
KAFB-106S7	V-S7-220119-469	280-119401-1	G	N	Naphthalene	U	Method Blank
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	Styrene	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	cis-1,3-Dichloropropene	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	trans-1,3-Dichloropropene	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	4-Chlorotoluene	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	1,4-Dichlorobenzene	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	1,2-Dibromoethane	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	1,2-Dichloroethane	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	4-Methyl-2-pentanone	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	Bromobenzene	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	Chlorobenzene	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	1,2,4-Trichlorobenzene	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	Dibromochloromethane	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	Tetrachloroethene	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	1,3-Dichloropropane	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	cis-1,2-Dichloroethene	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	trans-1,2-Dichloroethene	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	Methyl tert-butyl ether	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	1,2-Dichloroethene	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	1,3-Dichlorobenzene	UJ	Surrogate Recovery

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	Carbon tetrachloride	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	1,1-Dichloropropene	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	2-Hexanone	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	2,2-Dichloropropane	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	1,1,1,2-Tetrachloroethane	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	Chloroform	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	1,1,1-Trichloroethane	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	Bromomethane	J	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	Chloromethane	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	Dibromomethane	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	Bromochloromethane	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	Chloroethane	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	Vinyl Chloride	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	Methylene chloride	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	Carbon disulfide	J	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	Bromoform	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	Bromodichloromethane	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	1,1-Dichloroethane	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	1,1-Dichloroethene	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	Trichlorofluoromethane	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	Dichlorodifluoromethane	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	1,2-Dichloropropane	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	2-Butanone	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	1,1,2-Trichloroethane	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	Trichloroethene	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	1,1,2,2-Tetrachloroethane	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	1,2,3-Trichlorobenzene	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	Hexachloro-1,3-butadiene	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	2-Chlorotoluene	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	1,2-Dichlorobenzene	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	1,2-Dibromo-3-chloropropane	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	1,2,3-Trichloropropane	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	tert-Butylbenzene	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	Ethylbenzene	J	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	n-Propylbenzene	J	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	n-Butylbenzene	J	Surrogate Recovery

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	1,3,5-Trimethylbenzene	J	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	Toluene	J	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	sec-Butylbenzene	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	m- & p-Xylenes	J	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	Acetone	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	Benzene	J	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	Naphthalene	UJ	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	o-Xylene	J	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	1,2,4-Trimethylbenzene	J	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	Isopropylbenzene	J	Surrogate Recovery
KAFB-106S7	V-S7-220119-495	280-119401-1	G	N	4-Isopropyltoluene	UJ	Surrogate Recovery
KAFB-106S7	P-S7-210119-420	280-119377-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S7	V-S7-210119-420	280-119377-1	G	N	Toluene	U	Method Blank
KAFB-106S8	P-S8-180119-475	280-119347-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S8	P-S8-180119-475	280-119347-1	G	N	TPH-GRO (C6-C10)	J	MS/MSD - Percent recovery
KAFB-106S8	P-S8-180119-499	280-119347-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S8	P-S8-180119-514	280-119347-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S8	P-S8-180159-514	280-119347-1	G	FD	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S8	V-S8-180119-475	280-119347-1	G	N	Ethylbenzene	J	Surrogate Recovery
KAFB-106S8	V-S8-180119-475	280-119347-1	G	N	n-Propylbenzene	J	Surrogate Recovery
KAFB-106S8	V-S8-180119-475	280-119347-1	G	N	n-Butylbenzene	J	Surrogate Recovery
KAFB-106S8	V-S8-180119-475	280-119347-1	G	N	4-Methyl-2-pentanone	J	Surrogate Recovery
KAFB-106S8	V-S8-180119-475	280-119347-1	G	N	1,3,5-Trimethylbenzene	J	Surrogate Recovery
KAFB-106S8	V-S8-180119-475	280-119347-1	G	N	Toluene	J	Surrogate Recovery
KAFB-106S8	V-S8-180119-475	280-119347-1	G	N	2-Hexanone	J	Surrogate Recovery
KAFB-106S8	V-S8-180119-475	280-119347-1	G	N	Benzene	J	Surrogate Recovery
KAFB-106S8	V-S8-180119-475	280-119347-1	G	N	tert-Butylbenzene	J	Surrogate Recovery
KAFB-106S8	V-S8-180119-475	280-119347-1	G	N	Isopropylbenzene	J	Surrogate Recovery
KAFB-106S8	V-S8-180119-475	280-119347-1	G	N	4-Isopropyltoluene	J	Surrogate Recovery
KAFB-106S8	V-S8-180119-499	280-119347-1	G	N	Toluene	U	Method Blank
KAFB-106S8	V-S8-180119-499	280-119347-1	G	N	Acetone	U	Trip Blank
KAFB-106S8	V-S8-180119-514	280-119347-1	G	N	Toluene	U	Method Blank
KAFB-106S8	V-S8-180119-514	280-119347-1	G	N	Acetone	U	Trip Blank
KAFB-106S8	V-S8-180159-514	280-119347-1	G	FD	Toluene	U	Method Blank
KAFB-106S8	V-S8-180159-514	280-119347-1	G	FD	Acetone	U	Trip Blank
KAFB-106S9	P-S9-061018-102	280-115294-1	G	N	TPH-GRO (C6-C10)	U	Method Blank

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
KAFB-106S9	P-S9-071018-174	280-115294-1	G	N	TPH-GRO (C6-C10)	U	Method Blank
KAFB-106S9	V-S9-061018-102	280-115294-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S9	V-S9-061018-102	280-115294-1	G	N	1,2-Dibromoethane	UJ	MS/MSD - Percent recovery
KAFB-106S9	V-S9-071018-174	280-115294-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S9	P-S9-111018-342	280-115488-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S9	P-S9-111018-342	280-115488-1	G	N	TPH-GRO (C6-C10)	U	Method Blank
KAFB-106S9	P-S9-161018-415	280-115708-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S9	P-S9-161018-415	280-115708-1	G	N	TPH-GRO (C6-C10)	U	Method Blank
KAFB-106S9	V-S9-171018-470	280-115813-1	G	N	1,3,5-Trimethylbenzene	J	Surrogate Recovery
KAFB-106S9	V-S9-171018-470	280-115813-1	G	N	m- & p-Xylenes	J	Surrogate Recovery
KAFB-106S9	V-S9-171018-470	280-115813-1	G	N	Acetone	J	Surrogate Recovery
KAFB-106S9	V-S9-171018-470	280-115813-1	G	N	Naphthalene	U	Method Blank
KAFB-106S9	V-S9-171018-470	280-115813-1	G	N	o-Xylene	J	Surrogate Recovery
KAFB-106S9	V-S9-171018-470	280-115813-1	G	N	1,2,4-Trimethylbenzene	J	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	Ethylbenzene	J	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	Styrene	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	cis-1,3-Dichloropropene	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	trans-1,3-Dichloropropene	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	n-Propylbenzene	J	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	n-Butylbenzene	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	4-Chlorotoluene	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	1,4-Dichlorobenzene	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	1,2-Dibromoethane	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	1,2-Dichloroethane	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	4-Methyl-2-pentanone	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	1,3,5-Trimethylbenzene	J	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	Bromobenzene	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	Toluene	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	Chlorobenzene	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	1,2,4-Trichlorobenzene	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	Dibromochloromethane	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	Tetrachloroethene	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	sec-Butylbenzene	J	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	1,3-Dichloropropane	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	cis-1,2-Dichloroethene	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	trans-1,2-Dichloroethene	UJ	Surrogate Recovery

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	Methyl tert-butyl ether	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	m- & p-Xylenes	J	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	1,2-Dichloroethene	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	1,3-Dichlorobenzene	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	Carbon tetrachloride	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	1,1-Dichloropropene	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	2-Hexanone	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	2,2-Dichloropropane	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	1,1,1,2-Tetrachloroethane	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	Acetone	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	Chloroform	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	Benzene	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	1,1,1-Trichloroethane	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	Bromomethane	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	Chloromethane	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	Dibromomethane	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	Bromochloromethane	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	Chloroethane	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	Vinyl Chloride	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	Methylene chloride	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	Carbon disulfide	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	Bromoform	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	Bromodichloromethane	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	1,1-Dichloroethane	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	1,1-Dichloroethene	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	Trichlorofluoromethane	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	Dichlorodifluoromethane	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	1,2-Dichloropropane	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	2-Butanone	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	1,1,2-Trichloroethane	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	Trichloroethene	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	1,1,2,2-Tetrachloroethane	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	1,2,3-Trichlorobenzene	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	Hexachloro-1,3-butadiene	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	Naphthalene	J	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	o-Xylene	J	Surrogate Recovery

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	2-Chlorotoluene	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	1,2-Dichlorobenzene	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	1,2,4-Trimethylbenzene	J	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	1,2-Dibromo-3-chloropropane	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	1,2,3-Trichloropropane	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	tert-Butylbenzene	UJ	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	Isopropylbenzene	J	Surrogate Recovery
KAFB-106S9	V-S9-171018-475	280-115813-1	G	N	4-Isopropyltoluene	UJ	Surrogate Recovery
KAFB-106S9	P-S9-301018-501	280-116409-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S9	V-S9-301018-496	280-116409-1	G	N	Carbon disulfide	J	Surrogate Recovery
KAFB-106S9	V-S9-301018-496	280-116409-1	G	N	Naphthalene	J	Surrogate Recovery
KAFB-106S9	V-S9-301018-501	280-116409-1	G	N	n-Propylbenzene	J	Surrogate Recovery
KAFB-106S9	V-S9-301018-501	280-116409-1	G	N	n-Butylbenzene	J	Surrogate Recovery
KAFB-106S9	V-S9-301018-501	280-116409-1	G	N	1,2-Dibromoethane	J	Surrogate Recovery
KAFB-106S9	V-S9-301018-501	280-116409-1	G	N	1,3,5-Trimethylbenzene	J	Surrogate Recovery
KAFB-106S9	V-S9-301018-501	280-116409-1	G	N	sec-Butylbenzene	J	Surrogate Recovery
KAFB-106S9	V-S9-301018-501	280-116409-1	G	N	Acetone	J	Surrogate Recovery
KAFB-106S9	V-S9-301018-501	280-116409-1	G	N	Naphthalene	J	Surrogate Recovery
KAFB-106S9	V-S9-301018-501	280-116409-1	G	N	Isopropylbenzene	J	Surrogate Recovery
KAFB-106S9	V-S9-301018-501	280-116409-1	G	N	4-Isopropyltoluene	J	Surrogate Recovery
KAFB-106S9	P-S9-091018-252	280-115432-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106S9	P-S9-191018-490	280-115896-1	G	N	TPH-GRO (C6-C10)	J	Field Duplicate
KAFB-106S9	P-S9-191058-490	280-115896-1	G	FD	TPH-GRO (C6-C10)	J	Field Duplicate
KAFB-106S9	V-S9-191018-490	280-115896-1	G	N	1,2-Dibromoethane	J	Field Duplicate
KAFB-106S9	V-S9-191018-490	280-115896-1	G	N	Ethylbenzene	J	Field Duplicate
KAFB-106S9	V-S9-191018-490	280-115896-1	G	N	n-Propylbenzene	J	Field Duplicate
KAFB-106S9	V-S9-191018-490	280-115896-1	G	N	n-Butylbenzene	J	Field Duplicate
KAFB-106S9	V-S9-191018-490	280-115896-1	G	N	1,3,5-Trimethylbenzene	J	Field Duplicate
KAFB-106S9	V-S9-191018-490	280-115896-1	G	N	Toluene	J	Field Duplicate
KAFB-106S9	V-S9-191018-490	280-115896-1	G	N	m- & p-Xylenes	J	Field Duplicate
KAFB-106S9	V-S9-191018-490	280-115896-1	G	N	Benzene	J	Field Duplicate
KAFB-106S9	V-S9-191018-490	280-115896-1	G	N	Naphthalene	J	Field Duplicate
KAFB-106S9	V-S9-191018-490	280-115896-1	G	N	o-Xylene	J	Field Duplicate
KAFB-106S9	V-S9-191018-490	280-115896-1	G	N	1,2,4-Trimethylbenzene	J	Field Duplicate
KAFB-106S9	V-S9-191018-490	280-115896-1	G	N	Isopropylbenzene	J	Field Duplicate
KAFB-106S9	V-S9-191018-490	280-115896-1	G	N	4-Isopropyltoluene	J	Field Duplicate

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
KAFB-106S9	V-S9-191058-490	280-115896-1	G	FD	1,2-Dibromoethane	J	Hold Time
KAFB-106S9	V-S9-191058-490	280-115896-1	G	FD	Ethylbenzene	J	Field Duplicate
KAFB-106S9	V-S9-191058-490	280-115896-1	G	FD	n-Propylbenzene	J	Field Duplicate
KAFB-106S9	V-S9-191058-490	280-115896-1	G	FD	n-Butylbenzene	J	Field Duplicate
KAFB-106S9	V-S9-191058-490	280-115896-1	G	FD	1,3,5-Trimethylbenzene	J	Field Duplicate
KAFB-106S9	V-S9-191058-490	280-115896-1	G	FD	Toluene	J	Field Duplicate
KAFB-106S9	V-S9-191058-490	280-115896-1	G	FD	m- & p-Xylenes	J	Field Duplicate
KAFB-106S9	V-S9-191058-490	280-115896-1	G	FD	Benzene	J	Field Duplicate
KAFB-106S9	V-S9-191058-490	280-115896-1	G	FD	Naphthalene	J	Field Duplicate
KAFB-106S9	V-S9-191058-490	280-115896-1	G	FD	o-Xylene	J	Field Duplicate
KAFB-106S9	V-S9-191058-490	280-115896-1	G	FD	1,2,4-Trimethylbenzene	J	Field Duplicate
KAFB-106S9	V-S9-191058-490	280-115896-1	G	FD	Isopropylbenzene	J	Field Duplicate
KAFB-106S9	V-S9-191058-490	280-115896-1	G	FD	4-Isopropyltoluene	J	Field Duplicate
KAFB-106V1	P-V1-161218-061	280-118336-1	G	N	TPH-GRO (C6-C10)	J	Hold Time
KAFB-106V1	P-V1-161218-115	280-118336-1	G	N	TPH-GRO (C6-C10)	J	Hold Time
KAFB-106V1	P-V1-171218-147	280-118336-1	G	N	TPH-GRO (C6-C10)	J	Hold Time
KAFB-106V1	P-V1-171218-161	280-118390-1	G	N	Motor Oil (C20-C38)	J	Hold Time
KAFB-106V1	P-V1-171218-161	280-118390-1	G	N	TPH-DRO (C10-C28)	J	Hold Time
KAFB-106V1	P-V1-191218-285	280-118473-1	G	N	TPH-GRO (C6-C10)	J	Field Duplicate
KAFB-106V1	P-V1-191258-285	280-118473-1	G	FD	TPH-GRO (C6-C10)	J	Field Duplicate
KAFB-106V1	V-V1-191258-285	280-118473-1	G	FD	1,2-Dibromoethane	UJ	MS/MSD - Percent recovery
KAFB-106V2	P-V2-121218-080	280-118266-1	G	N	TPH-GRO (C6-C10)	J	Hold Time
KAFB-106V2	P-V2-121218-103	280-118266-1	G	N	TPH-GRO (C6-C10)	J	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	Ethylbenzene	J	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	Styrene	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	cis-1,3-Dichloropropene	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	trans-1,3-Dichloropropene	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	n-Propylbenzene	J	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	n-Butylbenzene	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	4-Chlorotoluene	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	1,4-Dichlorobenzene	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	1,2-Dibromoethane	J	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	1,2-Dichloroethane	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	4-Methyl-2-pentanone	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	1,3,5-Trimethylbenzene	J	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	Bromobenzene	UJ	Hold Time

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	Toluene	J	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	Chlorobenzene	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	1,2,4-Trichlorobenzene	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	Dibromochloromethane	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	Tetrachloroethene	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	sec-Butylbenzene	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	1,3-Dichloropropane	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	cis-1,2-Dichloroethene	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	trans-1,2-Dichloroethene	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	Methyl tert-butyl ether	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	m- & p-Xylenes	J	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	1,2-Dichloroethene	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	1,3-Dichlorobenzene	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	Carbon tetrachloride	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	1,1-Dichloropropene	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	2-Hexanone	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	2,2-Dichloropropane	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	1,1,1,2-Tetrachloroethane	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	Acetone	J	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	Chloroform	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	Benzene	J	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	1,1,1-Trichloroethane	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	Bromomethane	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	Chloromethane	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	Dibromomethane	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	Bromochloromethane	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	Chloroethane	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	Vinyl Chloride	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	Methylene chloride	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	Carbon disulfide	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	Bromoform	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	Bromodichloromethane	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	1,1-Dichloroethane	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	1,1-Dichloroethene	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	Trichlorofluoromethane	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	Dichlorodifluoromethane	UJ	Hold Time

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	1,2-Dichloropropane	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	2-Butanone	J	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	1,1,2-Trichloroethane	J	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	Trichloroethene	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	1,1,2,2-Tetrachloroethane	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	1,2,3-Trichlorobenzene	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	Hexachloro-1,3-butadiene	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	Naphthalene	J	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	o-Xylene	J	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	2-Chlorotoluene	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	1,2-Dichlorobenzene	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	1,2,4-Trimethylbenzene	J	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	1,2-Dibromo-3-chloropropane	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	1,2,3-Trichloropropane	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	tert-Butylbenzene	UJ	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	Isopropylbenzene	J	Hold Time
KAFB-106V2	V-V2-121218-080	280-118266-1	G	N	4-Isopropyltoluene	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	Ethylbenzene	J	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	Styrene	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	cis-1,3-Dichloropropene	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	trans-1,3-Dichloropropene	J	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	n-Propylbenzene	J	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	n-Butylbenzene	J	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	4-Chlorotoluene	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	1,4-Dichlorobenzene	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	1,2-Dibromoethane	J	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	1,2-Dichloroethane	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	4-Methyl-2-pentanone	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	1,3,5-Trimethylbenzene	J	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	Bromobenzene	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	Toluene	J	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	Chlorobenzene	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	1,2,4-Trichlorobenzene	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	Dibromochloromethane	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	Tetrachloroethene	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	sec-Butylbenzene	J	Hold Time

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	1,3-Dichloropropane	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	cis-1,2-Dichloroethene	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	trans-1,2-Dichloroethene	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	Methyl tert-butyl ether	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	m- & p-Xylenes	J	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	1,2-Dichloroethene	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	1,3-Dichlorobenzene	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	Carbon tetrachloride	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	1,1-Dichloropropene	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	2-Hexanone	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	2,2-Dichloropropane	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	1,1,1,2-Tetrachloroethane	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	Acetone	J	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	Chloroform	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	Benzene	J	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	1,1,1-Trichloroethane	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	Bromomethane	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	Chloromethane	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	Dibromomethane	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	Bromochloromethane	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	Chloroethane	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	Vinyl Chloride	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	Methylene chloride	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	Carbon disulfide	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	Bromoform	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	Bromodichloromethane	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	1,1-Dichloroethane	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	1,1-Dichloroethene	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	Trichlorofluoromethane	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	Dichlorodifluoromethane	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	1,2-Dichloropropane	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	2-Butanone	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	1,1,2-Trichloroethane	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	Trichloroethene	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	1,1,2,2-Tetrachloroethane	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	1,2,3-Trichlorobenzene	UJ	Hold Time

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	Hexachloro-1,3-butadiene	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	Naphthalene	J	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	o-Xylene	J	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	2-Chlorotoluene	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	1,2-Dichlorobenzene	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	1,2,4-Trimethylbenzene	J	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	1,2-Dibromo-3-chloropropane	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	1,2,3-Trichloropropane	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	tert-Butylbenzene	UJ	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	Isopropylbenzene	J	Hold Time
KAFB-106V2	V-V2-121218-103	280-118266-1	G	N	4-Isopropyltoluene	J	Hold Time
KAFB-106V2	P-V2-141218-270	280-118291-1	G	N	Motor Oil (C20-C38)	UJ	Hold Time
KAFB-106V2	P-V2-141218-270	280-118291-1	G	N	TPH-DRO (C10-C28)	J	Hold Time
KAFB-106V2	P-V2-141258-270	280-118291-1	G	FD	Motor Oil (C20-C38)	UJ	Hold Time
KAFB-106V2	P-V2-141258-270	280-118291-1	G	FD	TPH-DRO (C10-C28)	J	Hold Time
KAFB-106V2	V-V2-141218-270	280-118291-1	G	N	Toluene	J	Field Duplicate
KAFB-106V2	V-V2-141258-270	280-118291-1	G	FD	Toluene	J	Field Duplicate
KAFB-106V2	P-V2-151218-287	280-118337-1	G	N	TPH-DRO (C10-C28)	U	Method Blank
KAFB-106V2	P-V2-151218-287	280-118337-1	G	N	TPH-GRO (C6-C10)	J	Hold Time
Rinsate Blank	RB-S3-271118-01	280-117400-1	G	RB	Acetone	U	Method Blank
Rinsate Blank	RB-S3-271118-01	280-117400-1	G	RB	Motor Oil (C20-C38)	J	Lab Control Spike
Rinsate Blank	RB-S3-271118-02	280-117400-1	G	RB	Acetone	U	Method Blank
Rinsate Blank	RB-S4-051118-01	280-116899-1	G	RB	Motor Oil (C20-C38)	J	Hold Time
Rinsate Blank	RB-S4-051118-01	280-116899-1	G	RB	TPH-DRO (C10-C28)	J	Hold Time
Rinsate Blank	RB-S4-051118-02	280-116899-1	G	RB	Motor Oil (C20-C38)	UJ	Hold Time
Rinsate Blank	RB-S4-051118-02	280-116899-1	G	RB	TPH-DRO (C10-C28)	J	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	1,1,1,2-Tetrachloroethane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	1,1,1-Trichloroethane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	1,1,2,2-Tetrachloroethane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	1,1,2-Trichloroethane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	1,1-Dichloroethane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	1,1-Dichloroethene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	1,1-Dichloropropene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	1,2,3-Trichlorobenzene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	1,2,3-Trichloropropane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	1,2,4-Trichlorobenzene	UJ	Hold Time

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	1,2,4-Trimethylbenzene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	1,2-Dibromo-3-Chloropropane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	1,2-Dibromoethane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	1,2-Dichlorobenzene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	1,2-Dichloroethane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	1,2-Dichloroethene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	1,2-Dichloropropane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	1,3,5-Trimethylbenzene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	1,3-Dichlorobenzene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	1,3-Dichloropropane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	1,4-Dichlorobenzene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	2,2-Dichloropropane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	2-Butanone	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	2-Chlorotoluene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	2-Hexanone	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	4-Chlorotoluene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	4-Isopropyltoluene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	4-Methyl-2-Pentanone	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	Acetone	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	Benzene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	Bromobenzene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	Bromochloromethane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	Bromodichloromethane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	Bromoform	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	Bromomethane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	Carbon Disulfide	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	Carbon Tetrachloride	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	Chlorobenzene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	Chloroethane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	Chloroform	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	Chloromethane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	Cis-1,2-Dichloroethene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	Cis-1,3-Dichloropropene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	Dibromochloromethane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	Dibromomethane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	Dichlorodifluoromethane	UJ	Hold Time

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	Ethylbenzene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	Hexachloro-1,3-Butadiene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	Isopropylbenzene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	m- & p-Xylenes	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	Methyl tert-Butyl Ether	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	Methylene Chloride	J	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	Naphthalene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	n-Butylbenzene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	n-Propylbenzene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	o-Xylene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	sec-Butylbenzene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	Styrene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	tert-Butylbenzene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	Tetrachloroethene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	Toluene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	Trans-1,2-Dichloroethene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	Trans-1,3-Dichloropropene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	Trichloroethene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	Trichlorofluoromethane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-01	280-116206-1	G	RB	Vinyl Chloride	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	1,1,1,2-Tetrachloroethane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	1,1,1-Trichloroethane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	1,1,2,2-Tetrachloroethane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	1,1,2-Trichloroethane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	1,1-Dichloroethane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	1,1-Dichloroethene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	1,1-Dichloropropene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	1,2,3-Trichlorobenzene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	1,2,3-Trichloropropane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	1,2,4-Trichlorobenzene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	1,2,4-Trimethylbenzene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	1,2-Dibromo-3-Chloropropane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	1,2-Dibromoethane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	1,2-Dichlorobenzene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	1,2-Dichloroethane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	1,2-Dichloroethene	UJ	Hold Time

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	1,2-Dichloropropane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	1,3,5-Trimethylbenzene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	1,3-Dichlorobenzene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	1,3-Dichloropropane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	1,4-Dichlorobenzene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	2,2-Dichloropropane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	2-Butanone	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	2-Chlorotoluene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	2-Hexanone	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	4-Chlorotoluene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	4-Isopropyltoluene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	4-Methyl-2-Pentanone	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	Acetone	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	Benzene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	Bromobenzene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	Bromochloromethane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	Bromodichloromethane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	Bromoform	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	Bromomethane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	Carbon Disulfide	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	Carbon Tetrachloride	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	Chlorobenzene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	Chloroethane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	Chloroform	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	Chloromethane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	Cis-1,2-Dichloroethene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	Cis-1,3-Dichloropropene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	Dibromochloromethane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	Dibromomethane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	Dichlorodifluoromethane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	Ethylbenzene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	Hexachloro-1,3-Butadiene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	Isopropylbenzene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	m- & p-Xylenes	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	Methyl tert-Butyl Ether	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	Methylene Chloride	UJ	Hold Time

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	Naphthalene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	n-Butylbenzene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	n-Propylbenzene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	o-Xylene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	sec-Butylbenzene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	Styrene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	tert-Butylbenzene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	Tetrachloroethene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	Toluene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	Trans-1,2-Dichloroethene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	Trans-1,3-Dichloropropene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	Trichloroethene	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	Trichlorofluoromethane	UJ	Hold Time
Rinsate Blank	RB-S5-251018-02	280-116206-1	G	RB	Vinyl Chloride	UJ	Hold Time
Rinsate Blank	RB-S9-011118-01	280-116567-1	G	RB	Naphthalene	U	Method Blank
Rinsate Blank	RB-S9-011118-02	280-116567-1	G	RB	Naphthalene	U	Method Blank
Rinsate Blank	RB-V1-151218-01	280-118337-1	G	RB	Methylene Chloride	J	Surrogate Recovery
Rinsate Blank	RB-V1-151218-01	280-118337-1	G	RB	TPH-GRO (C6-C10)	J	Hold Time
Rinsate Blank	RB-V2-111218-01	280-118241-1	G	RB	TPH-GRO (C6-C10)	J	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	Ethylbenzene	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	Styrene	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	cis-1,3-Dichloropropene	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	trans-1,3-Dichloropropene	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	n-Propylbenzene	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	n-Butylbenzene	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	4-Chlorotoluene	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	1,4-Dichlorobenzene	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	1,2-Dibromoethane	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	1,2-Dichloroethane	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	4-Methyl-2-pentanone	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	1,3,5-Trimethylbenzene	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	Bromobenzene	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	Toluene	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	Chlorobenzene	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	1,2,4-Trichlorobenzene	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	Dibromochloromethane	UJ	Hold Time

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	Tetrachloroethene	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	sec-Butylbenzene	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	1,3-Dichloropropane	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	cis-1,2-Dichloroethene	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	trans-1,2-Dichloroethene	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	Methyl tert-butyl ether	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	m- & p-Xylenes	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	1,2-Dichloroethene	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	1,3-Dichlorobenzene	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	Carbon tetrachloride	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	1,1-Dichloropropene	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	2-Hexanone	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	2,2-Dichloropropane	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	1,1,1,2-Tetrachloroethane	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	Acetone	J	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	Chloroform	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	Benzene	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	1,1,1-Trichloroethane	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	Bromomethane	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	Chloromethane	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	Dibromomethane	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	Bromochloromethane	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	Chloroethane	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	Vinyl Chloride	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	Methylene chloride	J	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	Carbon disulfide	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	Bromoform	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	Bromodichloromethane	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	1,1-Dichloroethane	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	1,1-Dichloroethene	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	Trichlorofluoromethane	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	Dichlorodifluoromethane	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	1,2-Dichloropropane	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	2-Butanone	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	1,1,2-Trichloroethane	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	Trichloroethene	UJ	Hold Time

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	1,1,2,2-Tetrachloroethane	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	1,2,3-Trichlorobenzene	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	Hexachloro-1,3-butadiene	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	Naphthalene	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	o-Xylene	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	2-Chlorotoluene	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	1,2-Dichlorobenzene	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	1,2,4-Trimethylbenzene	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	1,2-Dibromo-3-chloropropane	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	1,2,3-Trichloropropane	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	tert-Butylbenzene	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	Isopropylbenzene	UJ	Hold Time
Trip Blank	TB-S9-161018-01	280-115708-1	G	TB	4-Isopropyltoluene	UJ	Hold Time
Trip Blank	TB-S4-051118-01	280-116615-2	G	TB	TPH-GRO (C6-C10)	U	Equipment Blank
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	TPH-GRO (C6-C10)	U	Equipment Blank
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	Ethylbenzene	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	Styrene	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	cis-1,3-Dichloropropene	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	trans-1,3-Dichloropropene	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	n-Propylbenzene	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	n-Butylbenzene	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	4-Chlorotoluene	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	1,4-Dichlorobenzene	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	1,2-Dibromoethane	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	1,2-Dichloroethane	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	4-Methyl-2-pentanone	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	1,3,5-Trimethylbenzene	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	Bromobenzene	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	Toluene	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	Chlorobenzene	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	1,2,4-Trichlorobenzene	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	Dibromochloromethane	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	Tetrachloroethene	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	sec-Butylbenzene	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	1,3-Dichloropropane	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	cis-1,2-Dichloroethene	UJ	Hold Time

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	trans-1,2-Dichloroethene	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	Methyl tert-butyl ether	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	m- & p-Xylenes	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	1,2-Dichloroethene	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	1,3-Dichlorobenzene	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	Carbon tetrachloride	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	1,1-Dichloropropene	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	2-Hexanone	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	2,2-Dichloropropane	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	1,1,1,2-Tetrachloroethane	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	Acetone	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	Chloroform	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	Benzene	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	1,1,1-Trichloroethane	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	Bromomethane	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	Chloromethane	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	Dibromomethane	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	Bromochloromethane	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	Chloroethane	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	Vinyl Chloride	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	Methylene chloride	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	Carbon disulfide	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	Bromoform	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	Bromodichloromethane	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	1,1-Dichloroethane	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	1,1-Dichloroethene	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	Trichlorofluoromethane	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	Dichlorodifluoromethane	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	1,2-Dichloropropane	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	2-Butanone	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	1,1,2-Trichloroethane	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	Trichloroethene	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	1,1,2,2-Tetrachloroethane	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	1,2,3-Trichlorobenzene	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	Hexachloro-1,3-butadiene	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	Naphthalene	UJ	Hold Time

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	o-Xylene	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	2-Chlorotoluene	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	1,2-Dichlorobenzene	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	1,2,4-Trimethylbenzene	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	1,2-Dibromo-3-chloropropane	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	1,2,3-Trichloropropane	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	tert-Butylbenzene	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	Isopropylbenzene	UJ	Hold Time
Trip Blank	TB-S5-221018-01	280-115956-1	G	TB	4-Isopropyltoluene	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	Ethylbenzene	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	Styrene	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	cis-1,3-Dichloropropene	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	trans-1,3-Dichloropropene	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	n-Propylbenzene	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	n-Butylbenzene	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	4-Chlorotoluene	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	1,4-Dichlorobenzene	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	1,2-Dibromoethane	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	1,2-Dichloroethane	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	4-Methyl-2-pentanone	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	1,3,5-Trimethylbenzene	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	Bromobenzene	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	Toluene	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	Chlorobenzene	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	1,2,4-Trichlorobenzene	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	Dibromochloromethane	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	Tetrachloroethene	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	sec-Butylbenzene	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	1,3-Dichloropropane	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	cis-1,2-Dichloroethene	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	trans-1,2-Dichloroethene	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	Methyl tert-butyl ether	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	m- & p-Xylenes	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	1,2-Dichloroethene	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	1,3-Dichlorobenzene	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	Carbon tetrachloride	UJ	Hold Time

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	1,1-Dichloropropene	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	2-Hexanone	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	2,2-Dichloropropane	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	1,1,1,2-Tetrachloroethane	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	Acetone	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	Chloroform	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	Benzene	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	1,1,1-Trichloroethane	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	Bromomethane	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	Chloromethane	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	Dibromomethane	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	Bromochloromethane	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	Chloroethane	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	Vinyl Chloride	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	Methylene chloride	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	Carbon disulfide	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	Bromoform	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	Bromodichloromethane	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	1,1-Dichloroethane	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	1,1-Dichloroethene	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	Trichlorofluoromethane	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	Dichlorodifluoromethane	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	1,2-Dichloropropane	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	2-Butanone	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	1,1,2-Trichloroethane	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	Trichloroethene	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	1,1,2,2-Tetrachloroethane	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	1,2,3-Trichlorobenzene	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	Hexachloro-1,3-butadiene	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	Naphthalene	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	o-Xylene	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	2-Chlorotoluene	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	1,2-Dichlorobenzene	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	1,2,4-Trimethylbenzene	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	1,2-Dibromo-3-chloropropane	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	1,2,3-Trichloropropane	UJ	Hold Time

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	tert-Butylbenzene	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	Isopropylbenzene	UJ	Hold Time
Trip Blank	TB-S9-221018-01	280-115951-1	G	TB	4-Isopropyltoluene	UJ	Hold Time
Trip Blank	TB-S5-231018-01	280-116044-1	G	TB	Methylene chloride	U	Method Blank
Trip Blank	TB-S3-211118-01	280-117359-1	G	TB	Acetone	U	Method Blank
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	Ethylbenzene	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	Styrene	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	cis-1,3-Dichloropropene	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	trans-1,3-Dichloropropene	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	n-Propylbenzene	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	n-Butylbenzene	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	4-Chlorotoluene	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	1,4-Dichlorobenzene	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	1,2-Dibromoethane	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	1,2-Dichloroethane	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	4-Methyl-2-pentanone	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	1,3,5-Trimethylbenzene	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	Bromobenzene	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	Toluene	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	Chlorobenzene	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	1,2,4-Trichlorobenzene	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	Dibromochloromethane	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	Tetrachloroethene	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	sec-Butylbenzene	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	1,3-Dichloropropane	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	cis-1,2-Dichloroethene	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	trans-1,2-Dichloroethene	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	Methyl tert-butyl ether	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	m- & p-Xylenes	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	1,2-Dichloroethene	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	1,3-Dichlorobenzene	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	Carbon tetrachloride	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	1,1-Dichloropropene	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	2-Hexanone	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	2,2-Dichloropropane	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	1,1,1,2-Tetrachloroethane	UJ	Hold Time

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	Acetone	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	Chloroform	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	Benzene	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	1,1,1-Trichloroethane	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	Bromomethane	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	Chloromethane	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	Dibromomethane	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	Bromochloromethane	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	Chloroethane	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	Vinyl Chloride	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	Methylene chloride	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	Carbon disulfide	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	Bromoform	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	Bromodichloromethane	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	1,1-Dichloroethane	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	1,1-Dichloroethene	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	Trichlorofluoromethane	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	Dichlorodifluoromethane	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	1,2-Dichloropropane	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	2-Butanone	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	1,1,2-Trichloroethane	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	Trichloroethene	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	1,1,2,2-Tetrachloroethane	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	1,2,3-Trichlorobenzene	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	Hexachloro-1,3-butadiene	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	Naphthalene	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	o-Xylene	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	2-Chlorotoluene	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	1,2-Dichlorobenzene	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	1,2,4-Trimethylbenzene	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	1,2-Dibromo-3-chloropropane	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	1,2,3-Trichloropropane	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	tert-Butylbenzene	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	Isopropylbenzene	UJ	Hold Time
Trip Blank	TB-S1-051218-02	280-117794-1	G	TB	4-Isopropyltoluene	UJ	Hold Time
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	Ethylbenzene	UJ	Surrogate Recovery

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	Styrene	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	cis-1,3-Dichloropropene	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	trans-1,3-Dichloropropene	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	n-Propylbenzene	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	n-Butylbenzene	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	4-Chlorotoluene	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	1,4-Dichlorobenzene	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	1,2-Dibromoethane	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	1,2-Dichloroethane	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	4-Methyl-2-pentanone	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	1,3,5-Trimethylbenzene	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	Bromobenzene	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	Toluene	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	Chlorobenzene	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	1,2,4-Trichlorobenzene	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	Dibromochloromethane	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	Tetrachloroethene	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	sec-Butylbenzene	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	1,3-Dichloropropane	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	cis-1,2-Dichloroethene	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	trans-1,2-Dichloroethene	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	Methyl tert-butyl ether	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	m- & p-Xylenes	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	1,2-Dichloroethene	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	1,3-Dichlorobenzene	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	Carbon tetrachloride	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	1,1-Dichloropropene	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	2-Hexanone	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	2,2-Dichloropropane	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	1,1,1,2-Tetrachloroethane	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	Acetone	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	Chloroform	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	Benzene	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	1,1,1-Trichloroethane	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	Bromomethane	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	Chloromethane	UJ	Surrogate Recovery

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	Dibromomethane	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	Bromochloromethane	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	Chloroethane	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	Vinyl Chloride	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	Methylene chloride	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	Carbon disulfide	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	Bromoform	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	Bromodichloromethane	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	1,1-Dichloroethane	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	1,1-Dichloroethene	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	Trichlorofluoromethane	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	Dichlorodifluoromethane	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	1,2-Dichloropropane	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	2-Butanone	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	1,1,2-Trichloroethane	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	Trichloroethene	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	1,1,2,2-Tetrachloroethane	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	1,2,3-Trichlorobenzene	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	Hexachloro-1,3-butadiene	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	Naphthalene	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	o-Xylene	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	2-Chlorotoluene	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	1,2-Dichlorobenzene	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	1,2,4-Trimethylbenzene	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	1,2-Dibromo-3-chloropropane	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	1,2,3-Trichloropropane	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	tert-Butylbenzene	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	Isopropylbenzene	UJ	Surrogate Recovery
Trip Blank	TB-V1-171218-01	280-118336-1	G	TB	4-Isopropyltoluene	UJ	Surrogate Recovery
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	Ethylbenzene	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	Styrene	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	cis-1,3-Dichloropropene	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	trans-1,3-Dichloropropene	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	n-Propylbenzene	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	n-Butylbenzene	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	4-Chlorotoluene	UJ	Hold Time

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	1,4-Dichlorobenzene	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	1,2-Dibromoethane	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	1,2-Dichloroethane	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	4-Methyl-2-pentanone	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	1,3,5-Trimethylbenzene	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	Bromobenzene	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	Toluene	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	Chlorobenzene	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	1,2,4-Trichlorobenzene	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	Dibromochloromethane	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	Tetrachloroethene	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	sec-Butylbenzene	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	1,3-Dichloropropane	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	cis-1,2-Dichloroethene	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	trans-1,2-Dichloroethene	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	Methyl tert-butyl ether	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	m- & p-Xylenes	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	1,2-Dichloroethene	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	1,3-Dichlorobenzene	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	Carbon tetrachloride	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	1,1-Dichloropropene	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	2-Hexanone	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	2,2-Dichloropropane	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	1,1,1,2-Tetrachloroethane	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	Acetone	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	Chloroform	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	Benzene	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	1,1,1-Trichloroethane	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	Bromomethane	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	Chloromethane	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	Dibromomethane	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	Bromochloromethane	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	Chloroethane	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	Vinyl Chloride	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	Methylene chloride	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	Carbon disulfide	UJ	Hold Time

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	Bromoform	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	Bromodichloromethane	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	1,1-Dichloroethane	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	1,1-Dichloroethene	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	Trichlorofluoromethane	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	Dichlorodifluoromethane	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	1,2-Dichloropropane	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	2-Butanone	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	1,1,2-Trichloroethane	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	Trichloroethene	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	1,1,2,2-Tetrachloroethane	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	1,2,3-Trichlorobenzene	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	Hexachloro-1,3-butadiene	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	Naphthalene	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	o-Xylene	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	2-Chlorotoluene	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	1,2-Dichlorobenzene	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	1,2,4-Trimethylbenzene	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	1,2-Dibromo-3-chloropropane	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	1,2,3-Trichloropropane	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	tert-Butylbenzene	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	Isopropylbenzene	UJ	Hold Time
Trip Blank	TB-V1-181218-01	280-118390-1	G	TB	4-Isopropyltoluene	UJ	Hold Time
Trip Blank	TB-V2-131218-01	280-118266-1	G	TB	TPH-GRO (C6-C10)	UJ	Hold Time
Trip Blank	TB-V2-141218-01	280-118290-1	G	TB	TPH-GRO (C6-C10)	UJ	Hold Time
Trip Blank	TB-V2-141218-02	280-118291-1	G	TB	TPH-GRO (C6-C10)	UJ	Hold Time
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	cis-1,3-Dichloropropene	UJ	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	trans-1,3-Dichloropropene	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	n-Propylbenzene	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	n-Butylbenzene	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	4-Chlorotoluene	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	1,4-Dichlorobenzene	UJ	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	1,2-Dichloroethane	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	4-Methyl-2-pentanone	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	1,3,5-Trimethylbenzene	UJ	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	Bromobenzene	UJ	Surrogate Recovery

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	Toluene	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	1,2,4-Trichlorobenzene	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	sec-Butylbenzene	UJ	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	cis-1,2-Dichloroethene	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	trans-1,2-Dichloroethene	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	Methyl tert-butyl ether	UJ	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	1,2-Dichloroethene	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	1,3-Dichlorobenzene	UJ	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	Carbon tetrachloride	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	1,1-Dichloropropene	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	2,2-Dichloropropane	UJ	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	Acetone	J	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	Chloroform	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	Benzene	UJ	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	1,1,1-Trichloroethane	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	Bromomethane	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	Chloromethane	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	Dibromomethane	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	Bromochloromethane	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	Chloroethane	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	Vinyl Chloride	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	Methylene chloride	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	Carbon disulfide	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	Bromodichloromethane	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	1,1-Dichloroethane	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	1,1-Dichloroethene	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	Trichlorofluoromethane	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	Dichlorodifluoromethane	UJ	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	1,2-Dichloropropane	UJ	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	2-Butanone	UJ	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	1,1,2-Trichloroethane	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	Trichloroethene	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	1,1,2,2-Tetrachloroethane	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	1,2,3-Trichlorobenzene	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	Hexachloro-1,3-butadiene	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	Naphthalene	U	Surrogate Recovery

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	2-Chlorotoluene	UJ	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	1,2-Dichlorobenzene	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	1,2,4-Trimethylbenzene	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	1,2-Dibromo-3-chloropropane	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	1,2,3-Trichloropropane	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	tert-Butylbenzene	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	Isopropylbenzene	U	Surrogate Recovery
Trip Blank	TB-V2-171218-01	280-118337-1	G	TB	4-Isopropyltoluene	U	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	Ethylbenzene	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	Styrene	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	cis-1,3-Dichloropropene	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	trans-1,3-Dichloropropene	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	n-Propylbenzene	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	n-Butylbenzene	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	4-Chlorotoluene	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	1,4-Dichlorobenzene	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	1,2-Dibromoethane	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	1,2-Dichloroethane	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	4-Methyl-2-pentanone	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	1,3,5-Trimethylbenzene	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	Bromobenzene	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	Toluene	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	Chlorobenzene	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	1,2,4-Trichlorobenzene	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	Dibromochloromethane	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	Tetrachloroethene	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	sec-Butylbenzene	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	1,3-Dichloropropane	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	cis-1,2-Dichloroethene	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	trans-1,2-Dichloroethene	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	Methyl tert-butyl ether	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	m- & p-Xylenes	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	1,2-Dichloroethene	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	1,3-Dichlorobenzene	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	Carbon tetrachloride	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	1,1-Dichloropropene	UJ	Surrogate Recovery

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	2-Hexanone	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	2,2-Dichloropropane	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	1,1,1,2-Tetrachloroethane	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	Acetone	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	Chloroform	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	Benzene	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	1,1,1-Trichloroethane	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	Bromomethane	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	Chloromethane	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	Dibromomethane	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	Bromochloromethane	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	Chloroethane	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	Vinyl Chloride	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	Methylene chloride	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	Carbon disulfide	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	Bromoform	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	Bromodichloromethane	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	1,1-Dichloroethane	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	1,1-Dichloroethene	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	Trichlorofluoromethane	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	Dichlorodifluoromethane	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	1,2-Dichloropropane	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	2-Butanone	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	1,1,2-Trichloroethane	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	Trichloroethene	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	1,1,2,2-Tetrachloroethane	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	1,2,3-Trichlorobenzene	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	Hexachloro-1,3-butadiene	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	Naphthalene	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	o-Xylene	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	2-Chlorotoluene	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	1,2-Dichlorobenzene	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	1,2,4-Trimethylbenzene	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	1,2-Dibromo-3-chloropropane	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	1,2,3-Trichloropropane	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	tert-Butylbenzene	UJ	Surrogate Recovery

Table H-3
Qualified Sample Results

Well Location ID	Sample Name	Sample Delivery Group	Collection Method	Sample Type	Analyte	Data Qualifier	Validation Reason Code
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	Isopropylbenzene	UJ	Surrogate Recovery
Trip Blank	TB-V1-191218-01	280-118473-1	G	TB	4-Isopropyltoluene	UJ	Surrogate Recovery
Trip Blank	TB-S8-180119-01	280-119347-1	G	TB	Acetone	J	Surrogate Recovery

CCV = continuing calibration verification

FD = field duplicate

G = grab sample

ID = identification

MS/MSD = matrix spike/matrix spike duplicate

N = normal field sample

RB = equipment rinse blank

TB = trip blank

Qualifiers:

J = Qualifier denotes the analyte was positively identified, but the associated numerical value is estimated.

U = Qualifier denotes the analyte was analyzed for but was not detected above the detection limit.

Table H-4
Field Quality Control Sample Results

Field Sample ID:			RB-S2-181118-01			RB-S2-181118-02			RB-S3-271118-01			RB-S3-271118-02			RB-S4-051118-01		
Sample Date:			11/18/2018			11/18/2018			11/27/2018			11/27/2018			11/5/2018		
Sample Type:			RB			RB			RB			RB			RB		
Parameter	Analytical Method	Analyte	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD
TPH	Method SW8015D (µg/L)	TPH-DRO (C10-C28)	0.096	J	0.11	0.075	J	0.13	0.12	J	0.12	0.058	J	0.12	0.11	J	0.12
		TPH-GRO (C6-C10)	0.071	J	0.11	ND	U	0.13	0.15	J	0.13	ND	U	0.12	0.068	J	0.13
		TPH-MRO (C20-C38)	37	--	25	35	--	25	23	J	25	36	--	25	14	J	25
EDB	Method SW8011 (µg/L)	1,2-Dibromoethane	0.052	--	0.014	0.061	--	0.014	0.09	--	0.014	0.094	--	0.014	ND	U	0.014
VOCS	Method SW8260B (µg/L)	1,1,1,2-Tetrachloroethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,1,1-Trichloroethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,1,2,2-Tetrachloroethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,1,2-Trichloroethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,1-Dichloroethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,1-Dichloroethene	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,1-Dichloropropene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,2,3-Trichlorobenzene	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,2,3-Trichloropropane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,2,4-Trichlorobenzene	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,2,4-Trimethylbenzene	0.2	J	0.4	0.22	J	0.4	0.35	J	0.4	0.62	J	0.4	ND	U	0.4
		1,2-Dibromo-3-Chloropropane	ND	U	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6
		1,2-Dibromoethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,2-Dichlorobenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,2-Dichloroethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	0.13	J	0.4	ND	U	0.4
		1,2-Dichloroethene	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2
		1,2-Dichloropropane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,3,5-Trimethylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	0.19	J	0.4	ND	U	0.4
		1,3-Dichlorobenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,3-Dichloropropane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,4-Dichlorobenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		2,2-Dichloropropane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		2-Butanone	ND	U	4	ND	U	4	ND	U	4	ND	U	4	ND	U	4
		2-Chlorotoluene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		2-Hexanone	ND	U	4	ND	U	4	ND	U	4	ND	U	4	ND	U	4
		4-Chlorotoluene	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		4-Isopropyltoluene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		4-Methyl-2-Pentanone	ND	U	3.2	ND	U	3.2	ND	U	3.2	ND	U	3.2	ND	U	3.2
		Acetone	ND	U	6.4	ND	U	6.4	ND	U	6.4	ND	U	6.4	ND	U	6.4
		Benzene	0.87	J	0.4	0.88	J	0.4	1.1	--	0.4	1.6	--	0.4	ND	U	0.4
		Bromobenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Bromochloromethane	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2
		Bromodichloromethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Bromoform	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Bromomethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Carbon Disulfide	ND	U	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6
		Carbon Tetrachloride	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Chlorobenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Chloroethane	ND	U	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6
		Chloroform	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4

Table H-4
Field Quality Control Sample Results

Field Sample ID:			RB-S2-181118-01			RB-S2-181118-02			RB-S3-271118-01			RB-S3-271118-02			RB-S4-051118-01		
Sample Date:			11/18/2018			11/18/2018			11/27/2018			11/27/2018			11/5/2018		
Sample Type:			RB			RB			RB			RB			RB		
Parameter	Analytical Method	Analyte	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD
		Chloromethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Cis-1,2-Dichloroethene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Cis-1,3-Dichloropropene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Dibromochloromethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Dibromomethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Dichlorodifluoromethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Ethylbenzene	0.96	J	0.4	0.98	J	0.4	1.2	--	0.4	1.6	--	0.4	ND	U	0.4
		Hexachloro-1,3-Butadiene	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Isopropylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		m- & p-Xylenes	2.7	--	0.8	2.9	--	0.8	2.7	--	0.8	4.1	--	0.8	ND	U	0.8
		Methyl tert-Butyl Ether	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Methylene Chloride	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Naphthalene	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		n-Butylbenzene	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		n-Propylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	0.23	J	0.4	ND	U	0.4
		o-Xylene	0.74	J	0.4	0.77	J	0.4	1	--	0.4	1.5	--	0.4	ND	U	0.4
		sec-Butylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Styrene	ND	U	0.4	0.18	J	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		tert-Butylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Tetrachloroethene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Toluene	6.9	--	0.4	7.1	--	0.4	7.3	--	0.4	10	--	0.4	0.81	J	0.4
		Trans-1,2-Dichloroethene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Trans-1,3-Dichloropropene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Trichloroethene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Trichlorofluoromethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Vinyl Chloride	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2

Table H-4
Field Quality Control Sample Results

Field Sample ID:			RB-S4-051118-02			RB-S5-251018-02			RB-S5-251018-01			RB-S7-200119-01			RB-S8-060119	
Sample Date:			11/5/2018			10/25/2018			10/25/2018			1/20/2019			1/6/2019	
Sample Type:			RB			RB			RB			RB			RB	
Parameter	Analytical Method	Analyte	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual
TPH	Method SW8015D (µg/L)	TPH-DRO (C10-C28)	0.079	J	0.12	0.1	J	0.11	0.086	J	0.12	0.048	J	0.12	0.1	J
		TPH-GRO (C6-C10)	ND	UJ	0.12	0.1	J	0.12	0.075	J	0.13	ND	U	0.13	ND	U
		TPH-MRO (C20-C38)	ND	U	25	ND	U	25	11	J	25	ND	U	25	ND	U
EDB	Method SW8011 (µg/L)	1,2-Dibromoethane	ND	U	0.014	ND	U	0.014	ND	U	0.014	ND	U	0.014	ND	U
VOCS	Method SW8260B (µg/L)	1,1,1,2-Tetrachloroethane	ND	U	0.8	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U
		1,1,1-Trichloroethane	ND	U	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U
		1,1,2,2-Tetrachloroethane	ND	U	0.8	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U
		1,1,2-Trichloroethane	ND	U	0.8	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U
		1,1-Dichloroethane	ND	U	0.8	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U
		1,1-Dichloroethene	ND	U	0.8	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U
		1,1-Dichloropropene	ND	U	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U
		1,2,3-Trichlorobenzene	ND	U	0.8	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U
		1,2,3-Trichloropropane	ND	U	0.8	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U
		1,2,4-Trichlorobenzene	ND	U	0.8	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U
		1,2,4-Trimethylbenzene	ND	U	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U
		1,2-Dibromo-3-Chloropropane	ND	U	1.6	ND	UJ	1.6	ND	UJ	1.6	ND	U	1.6	ND	U
		1,2-Dibromoethane	ND	U	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U
		1,2-Dichlorobenzene	ND	U	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U
		1,2-Dichloroethane	ND	U	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U
		1,2-Dichloroethene	ND	U	0.2	ND	UJ	0.2	ND	UJ	0.2	ND	U	0.2	ND	U
		1,2-Dichloropropane	ND	U	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U
		1,3,5-Trimethylbenzene	ND	U	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U
		1,3-Dichlorobenzene	ND	U	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U
		1,3-Dichloropropane	ND	U	0.8	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U
		1,4-Dichlorobenzene	ND	U	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U
		2,2-Dichloropropane	ND	U	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U
		2-Butanone	ND	U	4	ND	UJ	4	ND	UJ	4	ND	U	4	ND	U
		2-Chlorotoluene	ND	U	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U
		2-Hexanone	ND	U	4	ND	UJ	4	ND	UJ	4	ND	U	4	ND	U
		4-Chlorotoluene	ND	U	0.8	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U
		4-Isopropyltoluene	ND	U	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U
		4-Methyl-2-Pentanone	ND	U	3.2	ND	UJ	3.2	ND	UJ	3.2	ND	U	3.2	ND	U
		Acetone	ND	U	6.4	ND	UJ	6.4	ND	UJ	6.4	ND	U	6.4	ND	U
		Benzene	ND	U	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U
		Bromobenzene	ND	U	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U
		Bromochloromethane	ND	U	0.2	ND	UJ	0.2	ND	UJ	0.2	ND	U	0.2	ND	U
		Bromodichloromethane	ND	U	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U
		Bromoform	ND	U	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U
		Bromomethane	ND	U	0.8	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U
		Carbon Disulfide	ND	U	1.6	ND	UJ	1.6	ND	UJ	1.6	ND	U	1.6	ND	U
		Carbon Tetrachloride	ND	U	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U
		Chlorobenzene	ND	U	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U
		Chloroethane	ND	U	1.6	ND	UJ	1.6	ND	UJ	1.6	ND	U	1.6	ND	U
		Chloroform	ND	U	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U

Table H-4
Field Quality Control Sample Results

Field Sample ID:			RB-S4-051118-02			RB-S5-251018-02			RB-S5-251018-01			RB-S7-200119-01			RB-S8-060119-01	
Sample Date:			11/5/2018			10/25/2018			10/25/2018			1/20/2019			1/6/2019	
Sample Type:			RB			RB			RB			RB			RB	
Parameter	Analytical Method	Analyte	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual
		Chloromethane	ND	U	0.8	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U
		Cis-1,2-Dichloroethene	ND	U	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U
		Cis-1,3-Dichloropropene	ND	U	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U
		Dibromochloromethane	ND	U	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U
		Dibromomethane	ND	U	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U
		Dichlorodifluoromethane	ND	U	0.8	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U
		Ethylbenzene	ND	U	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U
		Hexachloro-1,3-Butadiene	ND	U	0.8	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U
		Isopropylbenzene	ND	U	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U
		m- & p-Xylenes	ND	U	0.8	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U
		Methyl tert-Butyl Ether	ND	U	0.8	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U
		Methylene Chloride	ND	U	0.8	ND	UJ	0.8	0.55	J	0.8	ND	U	0.8	ND	U
		Naphthalene	ND	U	0.8	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U
		n-Butylbenzene	ND	U	0.8	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U
		n-Propylbenzene	ND	U	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U
		o-Xylene	ND	U	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U
		sec-Butylbenzene	ND	U	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U
		Styrene	ND	U	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U
		tert-Butylbenzene	ND	U	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U
		Tetrachloroethene	ND	U	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U
		Toluene	0.65	J	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U
		Trans-1,2-Dichloroethene	ND	U	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U
		Trans-1,3-Dichloropropene	ND	U	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U
		Trichloroethene	ND	U	0.4	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U
		Trichlorofluoromethane	ND	U	0.8	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U
		Vinyl Chloride	ND	U	0.2	ND	UJ	0.2	ND	UJ	0.2	ND	U	0.2	ND	U

Table H-4
Field Quality Control Sample Results

Field Sample ID: J-01				RB-S9-201018-01			RB-S9-201018-02			RB-S9-011118-01			RB-S9-011118-02			RB-V1-151218-01		
Sample Date:				10/20/2018			10/20/2018			11/1/2018			11/1/2018			12/15/2018		
Sample Type:				RB			RB			RB			RB			RB		
Parameter	Analytical Method	Analyte	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD
TPH	Method SW8015D (µg/L)	TPH-DRO (C10-C28)	0.12	0.048	J	0.12	0.064	J	0.11	0.15	J	0.12	0.14	J	0.13	0.064	J	0.12
		TPH-GRO (C6-C10)	25	ND	U	0.12	ND	U	0.12	0.12	J	0.13	0.1	J	0.13	ND	U	0.12
		TPH-MRO (C20-C38)	0.12	25	--	25	13	J	25	ND	U	25	ND	U	25	21	J	25
EDB	Method SW8011 (µg/L)	1,2-Dibromoethane	0.014	ND	U	0.014	ND	U	0.014	ND	U	0.014	ND	U	0.014	ND	U	0.014
VOCS	Method SW8260B (µg/L)	1,1,1,2-Tetrachloroethane	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,1,1-Trichloroethane	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,1,2,2-Tetrachloroethane	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,1,2-Trichloroethane	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,1-Dichloroethane	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,1-Dichloroethene	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,1-Dichloropropene	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,2,3-Trichlorobenzene	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,2,3-Trichloropropane	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,2,4-Trichlorobenzene	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,2,4-Trimethylbenzene	0.4	ND	U	0.4	ND	U	0.4	2.8	--	0.4	0.55	J	0.4	ND	U	0.4
		1,2-Dibromo-3-Chloropropane	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6
		1,2-Dibromoethane	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,2-Dichlorobenzene	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,2-Dichloroethane	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,2-Dichloroethene	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2
		1,2-Dichloropropane	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,3,5-Trimethylbenzene	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,3-Dichlorobenzene	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,3-Dichloropropane	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,4-Dichlorobenzene	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		2,2-Dichloropropane	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		2-Butanone	4	ND	U	4	ND	U	4	ND	U	4	ND	U	4	ND	U	4
		2-Chlorotoluene	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		2-Hexanone	4	ND	U	4	ND	U	4	ND	U	4	ND	U	4	ND	U	4
		4-Chlorotoluene	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		4-Isopropyltoluene	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		4-Methyl-2-Pentanone	3.2	ND	U	3.2	ND	U	3.2	ND	U	3.2	ND	U	3.2	ND	U	3.2
		Acetone	6.4	2.8	J	6.4	2.8	J	6.4	ND	U	6.4	ND	U	6.4	ND	U	6.4
		Benzene	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Bromobenzene	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Bromochloromethane	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2
		Bromodichloromethane	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Bromoform	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Bromomethane	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Carbon Disulfide	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6
		Carbon Tetrachloride	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Chlorobenzene	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Chloroethane	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6
		Chloroform	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4

Table H-4
Field Quality Control Sample Results

Field Sample ID: J-01				RB-S9-201018-01			RB-S9-201018-02			RB-S9-011118-01			RB-S9-011118-02			RB-V1-151218-01		
Sample Date:				10/20/2018			10/20/2018			11/1/2018			11/1/2018			12/15/2018		
Sample Type:				RB			RB			RB			RB			RB		
Parameter	Analytical Method	Analyte	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD
		Chloromethane	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Cis-1,2-Dichloroethene	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Cis-1,3-Dichloropropene	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Dibromochloromethane	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Dibromomethane	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Dichlorodifluoromethane	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Ethylbenzene	0.4	ND	U	0.4	ND	U	0.4	1.6	--	0.4	0.33	J	0.4	ND	U	0.4
		Hexachloro-1,3-Butadiene	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Isopropylbenzene	0.4	ND	U	0.4	ND	U	0.4	0.22	J	0.4	ND	U	0.4	ND	U	0.4
		m- & p-Xylenes	0.8	ND	U	0.8	ND	U	0.8	2.5	--	0.8	0.52	J	0.8	ND	U	0.8
		Methyl tert-Butyl Ether	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Methylene Chloride	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	0.33	J	0.8
		Naphthalene	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		n-Butylbenzene	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		n-Propylbenzene	0.4	ND	U	0.4	ND	U	0.4	0.46	J	0.4	ND	U	0.4	ND	U	0.4
		o-Xylene	0.4	ND	U	0.4	ND	U	0.4	2	--	0.4	0.31	J	0.4	ND	U	0.4
		sec-Butylbenzene	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Styrene	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		tert-Butylbenzene	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Tetrachloroethene	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Toluene	0.4	ND	U	0.4	ND	U	0.4	0.71	J	0.4	0.19	J	0.4	ND	U	0.4
		Trans-1,2-Dichloroethene	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Trans-1,3-Dichloropropene	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Trichloroethene	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Trichlorofluoromethane	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Vinyl Chloride	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2

Table H-4
Field Quality Control Sample Results

Field Sample ID:			RB-V2-111218-01			RB-247-290119-01			RB-247-310119-01			TB-S1-031218-01			TB-S1-041218-01		
Sample Date:			12/11/2018			1/29/2019			1/31/2019			12/3/2018			12/4/2018		
Sample Type:			RB			RB			RB			TB			TB		
Parameter	Analytical Method	Analyte	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD
TPH	Method SW8015D (µg/L)	TPH-DRO (C10-C28)	0.067	J	0.13	--	--	--	0.045	J	0.12	--	--	--	--	--	--
		TPH-GRO (C6-C10)	23	J	25	--	--	--	ND	U	0.13	20	J	25	25	--	25
		TPH-MRO (C20-C38)	ND	U	0.13	--	--	--	ND	U	25	--	--	--	--	--	--
EDB	Method SW8011 (µg/L)	1,2-Dibromoethane	0.15	--	0.014	--	--	--	ND	U	0.014	ND	U	0.014	ND	U	0.014
VOCS	Method SW8260B (µg/L)	1,1,1,2-Tetrachloroethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,1,1-Trichloroethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,1,2,2-Tetrachloroethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,1,2-Trichloroethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,1-Dichloroethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,1-Dichloroethene	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,1-Dichloropropene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,2,3-Trichlorobenzene	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,2,3-Trichloropropane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,2,4-Trichlorobenzene	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,2,4-Trimethylbenzene	0.31	J	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,2-Dibromo-3-Chloropropane	ND	U	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6
		1,2-Dibromoethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,2-Dichlorobenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,2-Dichloroethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,2-Dichloroethene	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2
		1,2-Dichloropropane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,3,5-Trimethylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,3-Dichlorobenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,3-Dichloropropane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,4-Dichlorobenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		2,2-Dichloropropane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		2-Butanone	ND	U	4	ND	U	4	ND	U	4	ND	U	4	ND	U	4
		2-Chlorotoluene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		2-Hexanone	ND	U	4	ND	U	4	ND	U	4	ND	U	4	ND	U	4
		4-Chlorotoluene	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		4-Isopropyltoluene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		4-Methyl-2-Pentanone	ND	U	3.2	ND	U	3.2	ND	U	3.2	ND	U	3.2	ND	U	3.2
		Acetone	3.8	J	6.4	ND	U	6.4	ND	U	6.4	ND	U	6.4	ND	U	6.4
		Benzene	1.2	--	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Bromobenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Bromochloromethane	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2
		Bromodichloromethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Bromoform	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Bromomethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Carbon Disulfide	ND	U	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6
		Carbon Tetrachloride	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Chlorobenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Chloroethane	ND	U	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6
		Chloroform	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4

Table H-4
Field Quality Control Sample Results

Field Sample ID:			RB-V2-111218-01			RB-247-290119-01			RB-247-310119-01			TB-S1-031218-01			TB-S1-041218-01		
Sample Date:			12/11/2018			1/29/2019			1/31/2019			12/3/2018			12/4/2018		
Sample Type:			RB			RB			RB			TB			TB		
Parameter	Analytical Method	Analyte	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD
		Chloromethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Cis-1,2-Dichloroethene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Cis-1,3-Dichloropropene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Dibromochloromethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Dibromomethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Dichlorodifluoromethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Ethylbenzene	0.86	J	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Hexachloro-1,3-Butadiene	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Isopropylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		m- & p-Xylenes	3.1	--	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Methyl tert-Butyl Ether	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Methylene Chloride	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Naphthalene	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		n-Butylbenzene	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		n-Propylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		o-Xylene	1.1	--	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		sec-Butylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Styrene	0.23	J	0.4	ND	U	0.4	0.22	J	0.4	ND	U	0.4	ND	U	0.4
		tert-Butylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Tetrachloroethene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Toluene	8.7	--	0.4	0.4	J	0.4	0.38	J	0.4	ND	U	0.4	ND	U	0.4
		Trans-1,2-Dichloroethene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Trans-1,3-Dichloropropene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Trichloroethene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Trichlorofluoromethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Vinyl Chloride	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2

Table H-4
Field Quality Control Sample Results

Field Sample ID:			TB-S1-051218-02			TB-S1-271118-01			TB-S1-291118-01			TB-S1-051218-01			TB-S2-061118-01		
Sample Date:			12/5/2018			11/27/2018			11/29/2018			12/5/2018			11/6/2018		
Sample Type:			TB			TB			TB			TB			TB		
Parameter	Analytical Method	Analyte	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD
TPH	Method SW8015D (µg/L)	TPH-DRO (C10-C28)	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--
		TPH-GRO (C6-C10)	47	--	25	ND	U	25	ND	U	25	17	J	25	ND	U	25
		TPH-MRO (C20-C38)	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--
EDB	Method SW8011 (µg/L)	1,2-Dibromoethane	ND	U	0.014	--	--	--	ND	U	0.014	ND	U	0.014	--	--	--
VOCS	Method SW8260B (µg/L)	1,1,1,2-Tetrachloroethane	ND	UJ	0.8	--	--	--	ND	U	0.8	ND	UJ	0.8	--	--	--
		1,1,1-Trichloroethane	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--
		1,1,2,2-Tetrachloroethane	ND	UJ	0.8	--	--	--	ND	U	0.8	ND	UJ	0.8	--	--	--
		1,1,2-Trichloroethane	ND	UJ	0.8	--	--	--	ND	U	0.8	ND	UJ	0.8	--	--	--
		1,1-Dichloroethane	ND	UJ	0.8	--	--	--	ND	U	0.8	ND	UJ	0.8	--	--	--
		1,1-Dichloroethene	ND	UJ	0.8	--	--	--	ND	U	0.8	ND	UJ	0.8	--	--	--
		1,1-Dichloropropene	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--
		1,2,3-Trichlorobenzene	ND	UJ	0.8	--	--	--	ND	U	0.8	ND	UJ	0.8	--	--	--
		1,2,3-Trichloropropane	ND	UJ	0.8	--	--	--	ND	U	0.8	ND	UJ	0.8	--	--	--
		1,2,4-Trichlorobenzene	ND	UJ	0.8	--	--	--	ND	U	0.8	ND	UJ	0.8	--	--	--
		1,2,4-Trimethylbenzene	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--
		1,2-Dibromo-3-Chloropropane	ND	UJ	1.6	--	--	--	ND	U	1.6	ND	UJ	1.6	--	--	--
		1,2-Dibromoethane	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--
		1,2-Dichlorobenzene	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--
		1,2-Dichloroethane	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--
		1,2-Dichloroethene	ND	UJ	0.2	--	--	--	ND	U	0.2	ND	UJ	0.2	--	--	--
		1,2-Dichloropropane	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--
		1,3,5-Trimethylbenzene	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--
		1,3-Dichlorobenzene	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--
		1,3-Dichloropropane	ND	UJ	0.8	--	--	--	ND	U	0.8	ND	UJ	0.8	--	--	--
		1,4-Dichlorobenzene	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--
		2,2-Dichloropropane	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--
		2-Butanone	ND	UJ	4	--	--	--	ND	U	4	ND	UJ	4	--	--	--
		2-Chlorotoluene	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--
		2-Hexanone	ND	UJ	4	--	--	--	ND	U	4	ND	UJ	4	--	--	--
		4-Chlorotoluene	ND	UJ	0.8	--	--	--	ND	U	0.8	ND	UJ	0.8	--	--	--
		4-Isopropyltoluene	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--
		4-Methyl-2-Pentanone	ND	UJ	3.2	--	--	--	ND	U	3.2	ND	UJ	3.2	--	--	--
		Acetone	ND	UJ	6.4	--	--	--	2.1	J	6.4	ND	UJ	6.4	--	--	--
		Benzene	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--
		Bromobenzene	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--
		Bromochloromethane	ND	UJ	0.2	--	--	--	ND	U	0.2	ND	UJ	0.2	--	--	--
		Bromodichloromethane	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--
		Bromoform	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--
		Bromomethane	ND	UJ	0.8	--	--	--	ND	U	0.8	ND	UJ	0.8	--	--	--
		Carbon Disulfide	ND	UJ	1.6	--	--	--	ND	U	1.6	ND	UJ	1.6	--	--	--
		Carbon Tetrachloride	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--
		Chlorobenzene	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--
		Chloroethane	ND	UJ	1.6	--	--	--	ND	U	1.6	ND	UJ	1.6	--	--	--
		Chloroform	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--

Table H-4
Field Quality Control Sample Results

Field Sample ID:			TB-S1-051218-02			TB-S1-271118-01			TB-S1-291118-01			TB-S1-051218-01			TB-S2-061118-01		
Sample Date:			12/5/2018			11/27/2018			11/29/2018			12/5/2018			11/6/2018		
Sample Type:			TB			TB			TB			TB			TB		
Parameter	Analytical Method	Analyte	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD
		Chloromethane	ND	UJ	0.8	--	--	--	ND	U	0.8	ND	UJ	0.8	--	--	--
		Cis-1,2-Dichloroethene	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--
		Cis-1,3-Dichloropropene	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--
		Dibromochloromethane	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--
		Dibromomethane	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--
		Dichlorodifluoromethane	ND	UJ	0.8	--	--	--	ND	U	0.8	ND	UJ	0.8	--	--	--
		Ethylbenzene	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--
		Hexachloro-1,3-Butadiene	ND	UJ	0.8	--	--	--	ND	U	0.8	ND	UJ	0.8	--	--	--
		Isopropylbenzene	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--
		m- & p-Xylenes	ND	UJ	0.8	--	--	--	ND	U	0.8	ND	UJ	0.8	--	--	--
		Methyl tert-Butyl Ether	ND	UJ	0.8	--	--	--	ND	U	0.8	ND	UJ	0.8	--	--	--
		Methylene Chloride	ND	UJ	0.8	--	--	--	0.48	J	0.8	ND	UJ	0.8	--	--	--
		Naphthalene	ND	UJ	0.8	--	--	--	ND	U	0.8	ND	UJ	0.8	--	--	--
		n-Butylbenzene	ND	UJ	0.8	--	--	--	ND	U	0.8	ND	UJ	0.8	--	--	--
		n-Propylbenzene	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--
		o-Xylene	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--
		sec-Butylbenzene	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--
		Styrene	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--
		tert-Butylbenzene	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--
		Tetrachloroethene	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--
		Toluene	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--
		Trans-1,2-Dichloroethene	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--
		Trans-1,3-Dichloropropene	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--
		Trichloroethene	ND	UJ	0.4	--	--	--	ND	U	0.4	ND	UJ	0.4	--	--	--
		Trichlorofluoromethane	ND	UJ	0.8	--	--	--	ND	U	0.8	ND	UJ	0.8	--	--	--
		Vinyl Chloride	ND	UJ	0.2	--	--	--	ND	U	0.2	ND	UJ	0.2	--	--	--

Table H-4
Field Quality Control Sample Results

Field Sample ID:			TB-S2-151118-01			TB-S2-161118-01			TB-S2-161118-02			TB-S3-051118-01			TB-S3-191018-01		
Sample Date:			11/15/2018			11/16/2018			11/16/2018			11/5/2018			10/19/2018		
Sample Type:			TB			TB			TB			TB			TB		
Parameter	Analytical Method	Analyte	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD
TPH	Method SW8015D (µg/L)	TPH-DRO (C10-C28)	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--
		TPH-GRO (C6-C10)	ND	U	25	47	--	25	12	J	25	ND	U	25	19	J	25
		TPH-MRO (C20-C38)	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--
EDB	Method SW8011 (µg/L)	1,2-Dibromoethane	ND	U	0.014	ND	U	0.014	ND	U	0.014	--	--	--	--	--	--
VOCS	Method SW8260B (µg/L)	1,1,1,2-Tetrachloroethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--
		1,1,1-Trichloroethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--
		1,1,2,2-Tetrachloroethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--
		1,1,2-Trichloroethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--
		1,1-Dichloroethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--
		1,1-Dichloroethene	ND	U	0.8	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--
		1,1-Dichloropropene	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--
		1,2,3-Trichlorobenzene	ND	U	0.8	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--
		1,2,3-Trichloropropane	ND	U	0.8	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--
		1,2,4-Trichlorobenzene	ND	U	0.8	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--
		1,2,4-Trimethylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--
		1,2-Dibromo-3-Chloropropane	ND	U	1.6	ND	U	1.6	ND	U	1.6	--	--	--	--	--	--
		1,2-Dibromoethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--
		1,2-Dichlorobenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--
		1,2-Dichloroethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--
		1,2-Dichloroethene	ND	U	0.2	ND	U	0.2	ND	U	0.2	--	--	--	--	--	--
		1,2-Dichloropropane	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--
		1,3,5-Trimethylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--
		1,3-Dichlorobenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--
		1,3-Dichloropropane	ND	U	0.8	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--
		1,4-Dichlorobenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--
		2,2-Dichloropropane	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--
		2-Butanone	ND	U	4	ND	U	4	ND	U	4	--	--	--	--	--	--
		2-Chlorotoluene	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--
		2-Hexanone	ND	U	4	ND	U	4	ND	U	4	--	--	--	--	--	--
		4-Chlorotoluene	ND	U	0.8	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--
		4-Isopropyltoluene	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--
		4-Methyl-2-Pentanone	ND	U	3.2	ND	U	3.2	ND	U	3.2	--	--	--	--	--	--
		Acetone	ND	U	6.4	ND	U	6.4	ND	U	6.4	--	--	--	--	--	--
		Benzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--
		Bromobenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--
		Bromochloromethane	ND	U	0.2	ND	U	0.2	ND	U	0.2	--	--	--	--	--	--
		Bromodichloromethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--
		Bromoform	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--
		Bromomethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--
		Carbon Disulfide	ND	U	1.6	ND	U	1.6	ND	U	1.6	--	--	--	--	--	--
		Carbon Tetrachloride	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--
		Chlorobenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--
		Chloroethane	ND	U	1.6	ND	U	1.6	ND	U	1.6	--	--	--	--	--	--
		Chloroform	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--

Table H-4
Field Quality Control Sample Results

Field Sample ID:			TB-S2-151118-01			TB-S2-161118-01			TB-S2-161118-02			TB-S3-051118-01			TB-S3-191018-01		
Sample Date:			11/15/2018			11/16/2018			11/16/2018			11/5/2018			10/19/2018		
Sample Type:			TB			TB			TB			TB			TB		
Parameter	Analytical Method	Analyte	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD
		Chloromethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--
		Cis-1,2-Dichloroethene	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--
		Cis-1,3-Dichloropropene	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--
		Dibromochloromethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--
		Dibromomethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--
		Dichlorodifluoromethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--
		Ethylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--
		Hexachloro-1,3-Butadiene	ND	U	0.8	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--
		Isopropylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--
		m- & p-Xylenes	ND	U	0.8	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--
		Methyl tert-Butyl Ether	ND	U	0.8	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--
		Methylene Chloride	ND	U	0.8	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--
		Naphthalene	ND	U	0.8	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--
		n-Butylbenzene	ND	U	0.8	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--
		n-Propylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--
		o-Xylene	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--
		sec-Butylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--
		Styrene	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--
		tert-Butylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--
		Tetrachloroethene	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--
		Toluene	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--
		Trans-1,2-Dichloroethene	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--
		Trans-1,3-Dichloropropene	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--
		Trichloroethene	ND	U	0.4	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--
		Trichlorofluoromethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--
		Vinyl Chloride	ND	U	0.2	ND	U	0.2	ND	U	0.2	--	--	--	--	--	--

Table H-4
Field Quality Control Sample Results

Field Sample ID:			TB-S3-191118-01			TB-S3-201118-02			TB-S3-2018-01			TB-S3-211118-01			TB-S4-021118-01		
Sample Date:			11/19/2018			11/20/2018			11/20/2018			11/21/2018			11/2/2018		
Sample Type:			TB			TB			TB			TB			TB		
Parameter	Analytical Method	Analyte	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD
TPH	Method SW8015D (µg/L)	TPH-DRO (C10-C28)	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--
		TPH-GRO (C6-C10)	ND	U	25	ND	U	25	ND	U	25	ND	U	25	ND	U	25
		TPH-MRO (C20-C38)	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--
EDB	Method SW8011 (µg/L)	1,2-Dibromoethane	ND	U	0.014	ND	U	0.014	ND	U	0.014	ND	U	0.015	ND	U	0.014
VOCS	Method SW8260B (µg/L)	1,1,1,2-Tetrachloroethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,1,1-Trichloroethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,1,2,2-Tetrachloroethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,1,2-Trichloroethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,1-Dichloroethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,1-Dichloroethene	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,1-Dichloropropene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,2,3-Trichlorobenzene	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,2,3-Trichloropropane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,2,4-Trichlorobenzene	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,2,4-Trimethylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,2-Dibromo-3-Chloropropane	ND	U	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6
		1,2-Dibromoethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,2-Dichlorobenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,2-Dichloroethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,2-Dichloroethene	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2
		1,2-Dichloropropane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,3,5-Trimethylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,3-Dichlorobenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,3-Dichloropropane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,4-Dichlorobenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		2,2-Dichloropropane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		2-Butanone	ND	U	4	ND	U	4	ND	U	4	ND	U	4	ND	U	4
		2-Chlorotoluene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		2-Hexanone	ND	U	4	ND	U	4	ND	U	4	ND	U	4	ND	U	4
		4-Chlorotoluene	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		4-Isopropyltoluene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		4-Methyl-2-Pentanone	ND	U	3.2	ND	U	3.2	ND	U	3.2	ND	U	3.2	ND	U	3.2
		Acetone	ND	U	6.4	ND	U	6.4	ND	U	6.4	ND	U	6.4	ND	U	6.4
		Benzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Bromobenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Bromochloromethane	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2
		Bromodichloromethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Bromoform	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Bromomethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Carbon Disulfide	ND	U	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6
		Carbon Tetrachloride	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Chlorobenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Chloroethane	ND	U	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6
		Chloroform	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4

Table H-4
Field Quality Control Sample Results

Field Sample ID:			TB-S3-191118-01			TB-S3-201118-02			TB-S3-2018-01			TB-S3-211118-01			TB-S4-021118-01		
Sample Date:			11/19/2018			11/20/2018			11/20/2018			11/21/2018			11/2/2018		
Sample Type:			TB			TB			TB			TB			TB		
Parameter	Analytical Method	Analyte	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD
		Chloromethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Cis-1,2-Dichloroethene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Cis-1,3-Dichloropropene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Dibromochloromethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Dibromomethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Dichlorodifluoromethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Ethylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Hexachloro-1,3-Butadiene	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Isopropylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		m- & p-Xylenes	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Methyl tert-Butyl Ether	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Methylene Chloride	0.58	J	0.8	0.61	J	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Naphthalene	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		n-Butylbenzene	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		n-Propylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		o-Xylene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		sec-Butylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Styrene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		tert-Butylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Tetrachloroethene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Toluene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Trans-1,2-Dichloroethene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Trans-1,3-Dichloropropene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Trichloroethene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Trichlorofluoromethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Vinyl Chloride	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2

Table H-4
Field Quality Control Sample Results

Field Sample ID:			TB-S4-041118-01			TB-S4-051118-01			TB-S4-171018-01			TB-S5-091018-01			TB-S5-221018-01		
Sample Date:			11/4/2018			11/5/2018			10/17/2018			10/9/2018			10/22/2018		
Sample Type:			TB			TB			TB			TB			TB		
Parameter	Analytical Method	Analyte	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD
TPH	Method SW8015D (µg/L)	TPH-DRO (C10-C28)	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--
		TPH-GRO (C6-C10)	15	J	25	ND	U	25	14	J	25	ND	U	25	ND	U	25
		TPH-MRO (C20-C38)	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--
EDB	Method SW8011 (µg/L)	1,2-Dibromoethane	ND	U	0.014	ND	U	0.014	--	--	--	--	--	--	--	--	--
VOCS	Method SW8260B (µg/L)	1,1,1,2-Tetrachloroethane	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--	ND	UJ	0.8
		1,1,1-Trichloroethane	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4
		1,1,2,2-Tetrachloroethane	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--	ND	UJ	0.8
		1,1,2-Trichloroethane	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--	ND	UJ	0.8
		1,1-Dichloroethane	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--	ND	UJ	0.8
		1,1-Dichloroethene	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--	ND	UJ	0.8
		1,1-Dichloropropene	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4
		1,2,3-Trichlorobenzene	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--	ND	UJ	0.8
		1,2,3-Trichloropropane	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--	ND	UJ	0.8
		1,2,4-Trichlorobenzene	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--	ND	UJ	0.8
		1,2,4-Trimethylbenzene	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4
		1,2-Dibromo-3-Chloropropane	ND	U	1.6	ND	U	1.6	--	--	--	--	--	--	ND	UJ	1.6
		1,2-Dibromoethane	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4
		1,2-Dichlorobenzene	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4
		1,2-Dichloroethane	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4
		1,2-Dichloroethene	ND	U	0.2	ND	U	0.2	--	--	--	--	--	--	ND	UJ	0.2
		1,2-Dichloropropane	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4
		1,3,5-Trimethylbenzene	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4
		1,3-Dichlorobenzene	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4
		1,3-Dichloropropane	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--	ND	UJ	0.8
		1,4-Dichlorobenzene	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4
		2,2-Dichloropropane	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4
		2-Butanone	ND	U	4	ND	U	4	--	--	--	--	--	--	ND	UJ	4
		2-Chlorotoluene	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4
		2-Hexanone	ND	U	4	ND	U	4	--	--	--	--	--	--	ND	UJ	4
		4-Chlorotoluene	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--	ND	UJ	0.8
		4-Isopropyltoluene	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4
		4-Methyl-2-Pentanone	ND	U	3.2	ND	U	3.2	--	--	--	--	--	--	ND	UJ	3.2
		Acetone	ND	U	6.4	ND	U	6.4	--	--	--	--	--	--	ND	UJ	6.4
		Benzene	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4
		Bromobenzene	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4
		Bromochloromethane	ND	U	0.2	ND	U	0.2	--	--	--	--	--	--	ND	UJ	0.2
		Bromodichloromethane	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4
		Bromoform	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4
		Bromomethane	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--	ND	UJ	0.8
		Carbon Disulfide	ND	U	1.6	ND	U	1.6	--	--	--	--	--	--	ND	UJ	1.6
		Carbon Tetrachloride	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4
		Chlorobenzene	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4
		Chloroethane	ND	U	1.6	ND	U	1.6	--	--	--	--	--	--	ND	UJ	1.6
		Chloroform	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4

Table H-4
Field Quality Control Sample Results

Field Sample ID:			TB-S4-041118-01			TB-S4-051118-01			TB-S4-171018-01			TB-S5-091018-01			TB-S5-221018-01		
Sample Date:			11/4/2018			11/5/2018			10/17/2018			10/9/2018			10/22/2018		
Sample Type:			TB			TB			TB			TB			TB		
Parameter	Analytical Method	Analyte	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD
		Chloromethane	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--	ND	UJ	0.8
		Cis-1,2-Dichloroethene	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4
		Cis-1,3-Dichloropropene	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4
		Dibromochloromethane	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4
		Dibromomethane	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4
		Dichlorodifluoromethane	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--	ND	UJ	0.8
		Ethylbenzene	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4
		Hexachloro-1,3-Butadiene	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--	ND	UJ	0.8
		Isopropylbenzene	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4
		m- & p-Xylenes	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--	ND	UJ	0.8
		Methyl tert-Butyl Ether	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--	ND	UJ	0.8
		Methylene Chloride	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--	ND	UJ	0.8
		Naphthalene	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--	ND	UJ	0.8
		n-Butylbenzene	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--	ND	UJ	0.8
		n-Propylbenzene	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4
		o-Xylene	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4
		sec-Butylbenzene	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4
		Styrene	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4
		tert-Butylbenzene	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4
		Tetrachloroethene	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4
		Toluene	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4
		Trans-1,2-Dichloroethene	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4
		Trans-1,3-Dichloropropene	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4
		Trichloroethene	ND	U	0.4	ND	U	0.4	--	--	--	--	--	--	ND	UJ	0.4
		Trichlorofluoromethane	ND	U	0.8	ND	U	0.8	--	--	--	--	--	--	ND	UJ	0.8
		Vinyl Chloride	ND	U	0.2	ND	U	0.2	--	--	--	--	--	--	ND	UJ	0.2

Table H-4
Field Quality Control Sample Results

Field Sample ID:			TB-S5-231018-01			TB-S5-231018-02			TB-S7-210119-01			TB-S7-220119-01			TB-S8-160119-01		
Sample Date:			10/23/2018			10/23/2018			1/21/2019			1/22/2019			1/16/2019		
Sample Type:			TB			TB			TB			TB			TB		
Parameter	Analytical Method	Analyte	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD
TPH	Method SW8015D (µg/L)	TPH-DRO (C10-C28)	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--
		TPH-GRO (C6-C10)	17	J	25	24	J	25	ND	U	25	ND	U	25	ND	U	25
		TPH-MRO (C20-C38)	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--
EDB	Method SW8011 (µg/L)	1,2-Dibromoethane	--	--	--	--	--	--	ND	U	0.014	ND	U	0.014	ND	U	0.014
VOCS	Method SW8260B (µg/L)	1,1,1,2-Tetrachloroethane	ND	U	0.8	--	--	--	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,1,1-Trichloroethane	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,1,2,2-Tetrachloroethane	ND	U	0.8	--	--	--	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,1,2-Trichloroethane	ND	U	0.8	--	--	--	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,1-Dichloroethane	ND	U	0.8	--	--	--	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,1-Dichloroethene	ND	U	0.8	--	--	--	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,1-Dichloropropene	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,2,3-Trichlorobenzene	ND	U	0.8	--	--	--	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,2,3-Trichloropropane	ND	U	0.8	--	--	--	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,2,4-Trichlorobenzene	ND	U	0.8	--	--	--	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,2,4-Trimethylbenzene	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,2-Dibromo-3-Chloropropane	ND	U	1.6	--	--	--	ND	U	1.6	ND	U	1.6	ND	U	1.6
		1,2-Dibromoethane	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,2-Dichlorobenzene	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,2-Dichloroethane	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,2-Dichloroethene	ND	U	0.2	--	--	--	ND	U	0.2	ND	U	0.2	ND	U	0.2
		1,2-Dichloropropane	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,3,5-Trimethylbenzene	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,3-Dichlorobenzene	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,3-Dichloropropane	ND	U	0.8	--	--	--	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,4-Dichlorobenzene	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4
		2,2-Dichloropropane	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4
		2-Butanone	ND	U	4	--	--	--	ND	U	4	ND	U	4	ND	U	4
		2-Chlorotoluene	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4
		2-Hexanone	ND	U	4	--	--	--	ND	U	4	ND	U	4	ND	U	4
		4-Chlorotoluene	ND	U	0.8	--	--	--	ND	U	0.8	ND	U	0.8	ND	U	0.8
		4-Isopropyltoluene	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4
		4-Methyl-2-Pentanone	ND	U	3.2	--	--	--	ND	U	3.2	ND	U	3.2	ND	U	3.2
		Acetone	ND	U	6.4	--	--	--	ND	U	6.4	ND	U	6.4	ND	U	6.4
		Benzene	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Bromobenzene	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Bromochloromethane	ND	U	0.2	--	--	--	ND	U	0.2	ND	U	0.2	ND	U	0.2
		Bromodichloromethane	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Bromoform	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Bromomethane	ND	U	0.8	--	--	--	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Carbon Disulfide	ND	U	1.6	--	--	--	ND	U	1.6	ND	U	1.6	ND	U	1.6
		Carbon Tetrachloride	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Chlorobenzene	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Chloroethane	ND	U	1.6	--	--	--	ND	U	1.6	ND	U	1.6	ND	U	1.6
		Chloroform	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4

Table H-4
Field Quality Control Sample Results

Field Sample ID:			TB-S5-231018-01			TB-S5-231018-02			TB-S7-210119-01			TB-S7-220119-01			TB-S8-160119-01		
Sample Date:			10/23/2018			10/23/2018			1/21/2019			1/22/2019			1/16/2019		
Sample Type:			TB			TB			TB			TB			TB		
Parameter	Analytical Method	Analyte	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD
		Chloromethane	ND	U	0.8	--	--	--	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Cis-1,2-Dichloroethene	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Cis-1,3-Dichloropropene	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Dibromochloromethane	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Dibromomethane	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Dichlorodifluoromethane	ND	U	0.8	--	--	--	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Ethylbenzene	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Hexachloro-1,3-Butadiene	ND	U	0.8	--	--	--	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Isopropylbenzene	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4
		m- & p-Xylenes	ND	U	0.8	--	--	--	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Methyl tert-Butyl Ether	ND	U	0.8	--	--	--	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Methylene Chloride	ND	U	0.8	--	--	--	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Naphthalene	ND	U	0.8	--	--	--	ND	U	0.8	ND	U	0.8	ND	U	0.8
		n-Butylbenzene	ND	U	0.8	--	--	--	ND	U	0.8	ND	U	0.8	ND	U	0.8
		n-Propylbenzene	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4
		o-Xylene	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4
		sec-Butylbenzene	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Styrene	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4
		tert-Butylbenzene	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Tetrachloroethene	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Toluene	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Trans-1,2-Dichloroethene	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Trans-1,3-Dichloropropene	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Trichloroethene	ND	U	0.4	--	--	--	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Trichlorofluoromethane	ND	U	0.8	--	--	--	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Vinyl Chloride	ND	U	0.2	--	--	--	ND	U	0.2	ND	U	0.2	ND	U	0.2

Table H-4
Field Quality Control Sample Results

Field Sample ID:			TB-S8-180119-01			TB-S9-091018-01			TB-S9-101018-01			TB-S9-111018-01			TB-S9-161018-01		
Sample Date:			1/18/2019			10/9/2018			10/10/2018			10/11/2018			10/16/2018		
Sample Type:			TB			TB			TB			TB			TB		
Parameter	Analytical Method	Analyte	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD
TPH	Method SW8015D (µg/L)	TPH-DRO (C10-C28)	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--
		TPH-GRO (C6-C10)	ND	U	25	16	J	25	ND	U	25	13	J	25	ND	U	25
		TPH-MRO (C20-C38)	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--
EDB	Method SW8011 (µg/L)	1,2-Dibromoethane	ND	U	0.014	--	--	--	--	--	--	--	--	--	--	--	--
VOCS	Method SW8260B (µg/L)	1,1,1,2-Tetrachloroethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	UJ	0.8
		1,1,1-Trichloroethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4
		1,1,2,2-Tetrachloroethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	UJ	0.8
		1,1,2-Trichloroethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	UJ	0.8
		1,1-Dichloroethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	UJ	0.8
		1,1-Dichloroethene	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	UJ	0.8
		1,1-Dichloropropene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4
		1,2,3-Trichlorobenzene	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	UJ	0.8
		1,2,3-Trichloropropane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	UJ	0.8
		1,2,4-Trichlorobenzene	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	UJ	0.8
		1,2,4-Trimethylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4
		1,2-Dibromo-3-Chloropropane	ND	U	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6	ND	UJ	1.6
		1,2-Dibromoethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4
		1,2-Dichlorobenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4
		1,2-Dichloroethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4
		1,2-Dichloroethene	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	UJ	0.2
		1,2-Dichloropropane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4
		1,3,5-Trimethylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4
		1,3-Dichlorobenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4
		1,3-Dichloropropane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	UJ	0.8
		1,4-Dichlorobenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4
		2,2-Dichloropropane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4
		2-Butanone	ND	U	4	ND	U	4	ND	U	4	ND	U	4	ND	UJ	4
		2-Chlorotoluene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4
		2-Hexanone	ND	U	4	ND	U	4	ND	U	4	ND	U	4	ND	UJ	4
		4-Chlorotoluene	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	UJ	0.8
		4-Isopropyltoluene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4
		4-Methyl-2-Pentanone	ND	U	3.2	ND	U	3.2	ND	U	3.2	ND	U	3.2	ND	UJ	3.2
		Acetone	4.5	J	6.4	3	J	6.4	ND	U	6.4	3.2	J	6.4	2.5	J	6.4
		Benzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4
		Bromobenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4
		Bromochloromethane	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	UJ	0.2
		Bromodichloromethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4
		Bromoform	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4
		Bromomethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	UJ	0.8
		Carbon Disulfide	ND	U	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6	ND	UJ	1.6
		Carbon Tetrachloride	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4
		Chlorobenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4
		Chloroethane	ND	U	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6	ND	UJ	1.6
		Chloroform	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4

Table H-4
Field Quality Control Sample Results

Field Sample ID:			TB-S8-180119-01			TB-S9-091018-01			TB-S9-101018-01			TB-S9-111018-01			TB-S9-161018-01		
Sample Date:			1/18/2019			10/9/2018			10/10/2018			10/11/2018			10/16/2018		
Sample Type:			TB			TB			TB			TB			TB		
Parameter	Analytical Method	Analyte	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD
		Chloromethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	UJ	0.8
		Cis-1,2-Dichloroethene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4
		Cis-1,3-Dichloropropene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4
		Dibromochloromethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4
		Dibromomethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4
		Dichlorodifluoromethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	UJ	0.8
		Ethylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4
		Hexachloro-1,3-Butadiene	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	UJ	0.8
		Isopropylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4
		m- & p-Xylenes	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	UJ	0.8
		Methyl tert-Butyl Ether	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	UJ	0.8
		Methylene Chloride	ND	U	0.8	ND	U	0.8	0.62	J	0.8	0.51	J	0.8	0.53	J	0.8
		Naphthalene	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	UJ	0.8
		n-Butylbenzene	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	UJ	0.8
		n-Propylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4
		o-Xylene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4
		sec-Butylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4
		Styrene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4
		tert-Butylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4
		Tetrachloroethene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4
		Toluene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4
		Trans-1,2-Dichloroethene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4
		Trans-1,3-Dichloropropene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4
		Trichloroethene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	UJ	0.4
		Trichlorofluoromethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	UJ	0.8
		Vinyl Chloride	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	UJ	0.2

Table H-4
Field Quality Control Sample Results

Field Sample ID:			TB-S9-181018-01			TB-S9-191018-01			TB-S9-221018-01			TB-S9-311018-01			TB-V1-171218-01		
Sample Date:			10/18/2018			10/19/2018			10/22/2018			10/31/2018			12/17/2018		
Sample Type:			TB			TB			TB			TB			TB		
Parameter	Analytical Method	Analyte	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD
TPH	Method SW8015D (µg/L)	TPH-DRO (C10-C28)	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--
		TPH-GRO (C6-C10)	65	--	25	16	J	25	17	J	25	ND	U	25	ND	U	25
		TPH-MRO (C20-C38)	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--
EDB	Method SW8011 (µg/L)	1,2-Dibromoethane	--	--	--	--	--	--	ND	U	0.014	ND	U	0.014	ND	U	0.014
VOCS	Method SW8260B (µg/L)	1,1,1,2-Tetrachloroethane	ND	U	0.8	ND	U	0.8	ND	UJ	0.8	ND	U	0.8	ND	UJ	0.8
		1,1,1-Trichloroethane	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4
		1,1,2,2-Tetrachloroethane	ND	U	0.8	ND	U	0.8	ND	UJ	0.8	ND	U	0.8	ND	UJ	0.8
		1,1,2-Trichloroethane	ND	U	0.8	ND	U	0.8	ND	UJ	0.8	ND	U	0.8	ND	UJ	0.8
		1,1-Dichloroethane	ND	U	0.8	ND	U	0.8	ND	UJ	0.8	ND	U	0.8	ND	UJ	0.8
		1,1-Dichloroethene	ND	U	0.8	ND	U	0.8	ND	UJ	0.8	ND	U	0.8	ND	UJ	0.8
		1,1-Dichloropropene	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4
		1,2,3-Trichlorobenzene	ND	U	0.8	ND	U	0.8	ND	UJ	0.8	ND	U	0.8	ND	UJ	0.8
		1,2,3-Trichloropropane	ND	U	0.8	ND	U	0.8	ND	UJ	0.8	ND	U	0.8	ND	UJ	0.8
		1,2,4-Trichlorobenzene	ND	U	0.8	ND	U	0.8	ND	UJ	0.8	ND	U	0.8	ND	UJ	0.8
		1,2,4-Trimethylbenzene	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4
		1,2-Dibromo-3-Chloropropane	ND	U	1.6	ND	U	1.6	ND	UJ	1.6	ND	U	1.6	ND	UJ	1.6
		1,2-Dibromoethane	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4
		1,2-Dichlorobenzene	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4
		1,2-Dichloroethane	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4
		1,2-Dichloroethene	ND	U	0.2	ND	U	0.2	ND	UJ	0.2	ND	U	0.2	ND	UJ	0.2
		1,2-Dichloropropane	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4
		1,3,5-Trimethylbenzene	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4
		1,3-Dichlorobenzene	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4
		1,3-Dichloropropane	ND	U	0.8	ND	U	0.8	ND	UJ	0.8	ND	U	0.8	ND	UJ	0.8
		1,4-Dichlorobenzene	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4
		2,2-Dichloropropane	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4
		2-Butanone	ND	U	4	ND	U	4	ND	UJ	4	ND	U	4	ND	UJ	4
		2-Chlorotoluene	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4
		2-Hexanone	ND	U	4	ND	U	4	ND	UJ	4	ND	U	4	ND	UJ	4
		4-Chlorotoluene	ND	U	0.8	ND	U	0.8	ND	UJ	0.8	ND	U	0.8	ND	UJ	0.8
		4-Isopropyltoluene	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4
		4-Methyl-2-Pentanone	ND	U	3.2	ND	U	3.2	ND	UJ	3.2	ND	U	3.2	ND	UJ	3.2
		Acetone	2.7	J	6.4	2.8	J	6.4	ND	UJ	6.4	ND	U	6.4	ND	UJ	6.4
		Benzene	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4
		Bromobenzene	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4
		Bromochloromethane	ND	U	0.2	ND	U	0.2	ND	UJ	0.2	ND	U	0.2	ND	UJ	0.2
		Bromodichloromethane	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4
		Bromoform	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4
		Bromomethane	ND	U	0.8	ND	U	0.8	ND	UJ	0.8	ND	U	0.8	ND	UJ	0.8
		Carbon Disulfide	ND	U	1.6	ND	U	1.6	ND	UJ	1.6	ND	U	1.6	ND	UJ	1.6
		Carbon Tetrachloride	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4
		Chlorobenzene	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4
		Chloroethane	ND	U	1.6	ND	U	1.6	ND	UJ	1.6	ND	U	1.6	ND	UJ	1.6
		Chloroform	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4

Table H-4
Field Quality Control Sample Results

Field Sample ID:			TB-S9-181018-01			TB-S9-191018-01			TB-S9-221018-01			TB-S9-311018-01			TB-V1-171218-01		
Sample Date:			10/18/2018			10/19/2018			10/22/2018			10/31/2018			12/17/2018		
Sample Type:			TB			TB			TB			TB			TB		
Parameter	Analytical Method	Analyte	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD
		Chloromethane	ND	U	0.8	ND	U	0.8	ND	UJ	0.8	ND	U	0.8	ND	UJ	0.8
		Cis-1,2-Dichloroethene	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4
		Cis-1,3-Dichloropropene	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4
		Dibromochloromethane	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4
		Dibromomethane	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4
		Dichlorodifluoromethane	ND	U	0.8	ND	U	0.8	ND	UJ	0.8	ND	U	0.8	ND	UJ	0.8
		Ethylbenzene	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4
		Hexachloro-1,3-Butadiene	ND	U	0.8	ND	U	0.8	ND	UJ	0.8	ND	U	0.8	ND	UJ	0.8
		Isopropylbenzene	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4
		m- & p-Xylenes	ND	U	0.8	ND	U	0.8	ND	UJ	0.8	ND	U	0.8	ND	UJ	0.8
		Methyl tert-Butyl Ether	ND	U	0.8	ND	U	0.8	ND	UJ	0.8	ND	U	0.8	ND	UJ	0.8
		Methylene Chloride	0.59	J	0.8	0.77	J	0.8	ND	UJ	0.8	ND	U	0.8	ND	UJ	0.8
		Naphthalene	ND	U	0.8	ND	U	0.8	ND	UJ	0.8	ND	U	0.8	ND	UJ	0.8
		n-Butylbenzene	ND	U	0.8	ND	U	0.8	ND	UJ	0.8	ND	U	0.8	ND	UJ	0.8
		n-Propylbenzene	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4
		o-Xylene	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4
		sec-Butylbenzene	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4
		Styrene	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4
		tert-Butylbenzene	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4
		Tetrachloroethene	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4
		Toluene	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4
		Trans-1,2-Dichloroethene	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4
		Trans-1,3-Dichloropropene	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4
		Trichloroethene	ND	U	0.4	ND	U	0.4	ND	UJ	0.4	ND	U	0.4	ND	UJ	0.4
		Trichlorofluoromethane	ND	U	0.8	ND	U	0.8	ND	UJ	0.8	ND	U	0.8	ND	UJ	0.8
		Vinyl Chloride	ND	U	0.2	ND	U	0.2	ND	UJ	0.2	ND	U	0.2	ND	UJ	0.2

Table H-4
Field Quality Control Sample Results

Field Sample ID:			TB-V1-181218-01			TB-V1-191218-01			TB-V2-131218-01			TB-V2-141218-01			TB-V2-141218-02		
Sample Date:			12/18/2018			12/19/2018			12/13/2018			12/14/2018			12/14/2018		
Sample Type:			TB			TB			TB			TB			TB		
Parameter	Analytical Method	Analyte	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD
TPH	Method SW8015D (µg/L)	TPH-DRO (C10-C28)	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--
		TPH-GRO (C6-C10)	ND	U	25	ND	U	25	ND	UJ	25	ND	UJ	25	ND	UJ	25
		TPH-MRO (C20-C38)	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--
EDB	Method SW8011 (µg/L)	1,2-Dibromoethane	ND	U	0.014	ND	U	0.014	ND	U	0.014	ND	U	0.014	--	--	--
VOCS	Method SW8260B (µg/L)	1,1,1,2-Tetrachloroethane	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,1,1-Trichloroethane	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,1,2,2-Tetrachloroethane	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,1,2-Trichloroethane	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,1-Dichloroethane	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,1-Dichloroethene	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,1-Dichloropropene	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,2,3-Trichlorobenzene	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,2,3-Trichloropropane	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,2,4-Trichlorobenzene	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,2,4-Trimethylbenzene	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,2-Dibromo-3-Chloropropane	ND	UJ	1.6	ND	UJ	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6
		1,2-Dibromoethane	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,2-Dichlorobenzene	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,2-Dichloroethane	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,2-Dichloroethene	ND	UJ	0.2	ND	UJ	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2
		1,2-Dichloropropane	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,3,5-Trimethylbenzene	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,3-Dichlorobenzene	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,3-Dichloropropane	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,4-Dichlorobenzene	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		2,2-Dichloropropane	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		2-Butanone	ND	UJ	4	ND	UJ	4	ND	U	4	ND	U	4	ND	U	4
		2-Chlorotoluene	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		2-Hexanone	ND	UJ	4	ND	UJ	4	ND	U	4	ND	U	4	ND	U	4
		4-Chlorotoluene	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		4-Isopropyltoluene	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		4-Methyl-2-Pentanone	ND	UJ	3.2	ND	UJ	3.2	ND	U	3.2	ND	U	3.2	ND	U	3.2
		Acetone	ND	UJ	6.4	ND	UJ	6.4	ND	U	6.4	ND	U	6.4	ND	U	6.4
		Benzene	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Bromobenzene	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Bromochloromethane	ND	UJ	0.2	ND	UJ	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2
		Bromodichloromethane	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Bromoform	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Bromomethane	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Carbon Disulfide	ND	UJ	1.6	ND	UJ	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6
		Carbon Tetrachloride	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Chlorobenzene	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Chloroethane	ND	UJ	1.6	ND	UJ	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6
		Chloroform	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4

Table H-4
Field Quality Control Sample Results

Field Sample ID:			TB-V1-181218-01			TB-V1-191218-01			TB-V2-131218-01			TB-V2-141218-01			TB-V2-141218-02		
Sample Date:			12/18/2018			12/19/2018			12/13/2018			12/14/2018			12/14/2018		
Sample Type:			TB			TB			TB			TB			TB		
Parameter	Analytical Method	Analyte	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD
		Chloromethane	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Cis-1,2-Dichloroethene	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Cis-1,3-Dichloropropene	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Dibromochloromethane	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Dibromomethane	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Dichlorodifluoromethane	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Ethylbenzene	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Hexachloro-1,3-Butadiene	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Isopropylbenzene	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		m- & p-Xylenes	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Methyl tert-Butyl Ether	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Methylene Chloride	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Naphthalene	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		n-Butylbenzene	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		n-Propylbenzene	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		o-Xylene	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		sec-Butylbenzene	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Styrene	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		tert-Butylbenzene	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Tetrachloroethene	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Toluene	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Trans-1,2-Dichloroethene	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Trans-1,3-Dichloropropene	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Trichloroethene	ND	UJ	0.4	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Trichlorofluoromethane	ND	UJ	0.8	ND	UJ	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Vinyl Chloride	ND	UJ	0.2	ND	UJ	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2

Table H-4
Field Quality Control Sample Results

Field Sample ID:			TB-V2-171218-01			TB-247-050219-01			TB-247-300119-01			TB-247-310119-01			SWB-S9-201018-01		
Sample Date:			12/17/2018			2/5/2019			1/30/2019			1/31/2019			10/20/2018		
Sample Type:			TB			TB			TB			TB			SB		
Parameter	Analytical Method	Analyte	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD
TPH	Method SW8015D (µg/L)	TPH-DRO (C10-C28)	--	--	--	--	--	--	--	--	--	--	--	--	0.049	J	0.12
		TPH-GRO (C6-C10)	ND	U	25	ND	U	25	ND	U	25	ND	U	25	ND	U	0.12
		TPH-MRO (C20-C38)	--	--	--	--	--	--	--	--	--	--	--	--	23	J	25
EDB	Method SW8011 (µg/L)	1,2-Dibromoethane	ND	U	0.014	ND	U	0.014	ND	U	0.014	ND	U	0.014	ND	U	0.014
VOCS	Method SW8260B (µg/L)	1,1,1,2-Tetrachloroethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,1,1-Trichloroethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,1,2,2-Tetrachloroethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,1,2-Trichloroethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,1-Dichloroethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,1-Dichloroethene	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,1-Dichloropropene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,2,3-Trichlorobenzene	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,2,3-Trichloropropane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,2,4-Trichlorobenzene	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,2,4-Trimethylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,2-Dibromo-3-Chloropropane	ND	U	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6
		1,2-Dibromoethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,2-Dichlorobenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,2-Dichloroethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,2-Dichloroethene	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2
		1,2-Dichloropropane	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,3,5-Trimethylbenzene	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,3-Dichlorobenzene	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		1,3-Dichloropropane	ND	U	0.8	ND	U	0.2	ND	U	0.8	ND	U	0.8	ND	U	0.8
		1,4-Dichlorobenzene	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		2,2-Dichloropropane	ND	UJ	0.4	ND	U	0.8	ND	U	0.4	ND	U	0.4	ND	U	0.4
		2-Butanone	ND	UJ	4	ND	U	4	ND	U	4	ND	U	4	ND	U	4
		2-Chlorotoluene	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		2-Hexanone	ND	U	4	ND	U	4	ND	U	4	ND	U	4	ND	U	4
		4-Chlorotoluene	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		4-Isopropyltoluene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		4-Methyl-2-Pentanone	ND	U	3.2	ND	U	3.2	ND	U	3.2	ND	U	3.2	ND	U	3.2
		Acetone	2.5	J	6.4	ND	U	6.4	ND	U	6.4	ND	U	6.4	2.6	J	6.4
		Benzene	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Bromobenzene	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Bromochloromethane	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2
		Bromodichloromethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	0.89	J	0.4
		Bromoform	ND	U	0.4	ND	U	1	ND	U	0.4	ND	U	0.4	4.6	--	0.4
		Bromomethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Carbon Disulfide	ND	U	1.6	ND	U	0.8	ND	U	1.6	ND	U	1.6	ND	U	1.6
		Carbon Tetrachloride	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Chlorobenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Chloroethane	ND	U	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6	ND	U	1.6
		Chloroform	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	0.39	J	0.4

Table H-4
Field Quality Control Sample Results

Field Sample ID:			TB-V2-171218-01			TB-247-050219-01			TB-247-300119-01			TB-247-310119-01			SWB-S9-201018-01		
Sample Date:			12/17/2018			2/5/2019			1/30/2019			1/31/2019			10/20/2018		
Sample Type:			TB			TB			TB			TB			SB		
Parameter	Analytical Method	Analyte	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD
		Chloromethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Cis-1,2-Dichloroethene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Cis-1,3-Dichloropropene	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Dibromochloromethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	2.3	--	0.4
		Dibromomethane	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Dichlorodifluoromethane	ND	UJ	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Ethylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Hexachloro-1,3-Butadiene	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Isopropylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		m- & p-Xylenes	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Methyl tert-Butyl Ether	ND	UJ	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Methylene Chloride	ND	U	0.8	ND	U	2	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Naphthalene	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		n-Butylbenzene	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		n-Propylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		o-Xylene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		sec-Butylbenzene	ND	UJ	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Styrene	ND	U	0.4	ND	U	0.8	ND	U	0.4	ND	U	0.4	ND	U	0.4
		tert-Butylbenzene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Tetrachloroethene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Toluene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Trans-1,2-Dichloroethene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Trans-1,3-Dichloropropene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Trichloroethene	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4	ND	U	0.4
		Trichlorofluoromethane	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8	ND	U	0.8
		Vinyl Chloride	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2	ND	U	0.2

Table H-4
Field Quality Control Sample Results

µg/L = microgram per liter.
DRO = diesel range organics
EDB = ethylene dibromide (1,2-dibromoethane)
GRO = gasoline range organics
ID = identification
LOD = limit of detection
MRO = motor oil organics
ND = not detected above the detection limit
RB = equipment rinse blank
SW = source water blank
TB = trip blank
TPH = total petroleum hydrocarbon
Val Qual = validation qualifier
VOC = volatile organic compound
Shading = detected concentrations above the detection limit

Qualifiers:
Val Qual based on independent data validation.
J = Qualifier denotes the analyte was positively identified, but the associated numerical value is estimated.
U = Qualifier denotes the analyte was analyzed but not detected above the detection limit. The value associated with the U-qualifier is the LOD.

Table H-5
Field Duplicate Sample Results

		Boring ID:	KAFB-106S1			KAFB-106S1			KAFB-106S2			KAFB-106S2			KAFB-106S3			KAFB-106S3			KAFB-106S4		
		Sample Date:	12/5/2018			12/5/2018			11/16/2018			11/16/2018			11/20/2018			11/20/2018			11/5/2018		
		Sample Depth (ft bgs):	510			510			510			510			459			459			504		
		Sample Type:	REG			Field Duplicate			REG			Field Duplicate			REG			Field Duplicate			REG		
Parameter	Analytical Method	Chemical Name	CAS	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD		
TPH	Method SW8015D (mg/kg)	TPH-MRO (C20-C38)	MOIL	ND	U	24	ND	U	24	ND	U	25	ND	U	23	ND	U	22	ND	U	26		
		TPH-DRO (C10-C28)	TPH-DRO	ND	U	4.1	ND	U	4.2	ND	U	4.2	ND	U	3.9	ND	U	3.8	ND	U	4.4		
		TPH-GRO (C6-C10)	TPH-GRO	ND	U	1.2	ND	U	1.6	2.9	J	1.2	7.6	J	1.2	ND	U	1.4	ND	U	1.2		
Moisture	ASTM D2216 (%) TPH	NA	NA	10.9	--	--	11.2	--	--	10.9	--	--	9.1	--	--	3.2	--	--	2.9	--	--		
EDB	Method SW8011 (mg/kg) ^a	1,2-Dibromoethane	106-93-4	ND	U	0.00043	ND	U	0.00042	ND	U	0.00042	ND	U	0.00043	ND	UJ	0.00039	ND	U	0.00047		
Moisture	ASTM D2216 (%) EDB	NA	NA	11.2	--	--	9.3	--	--	10.1	--	--	11.3	--	--	3.6	--	--	3.6	--	--		
VOCs	Method SW8260B (mg/kg) ^a	1,1,1,2-Tetrachloroethane	630-20-6	ND	U	0.0017	ND	U	0.0015	ND	U	0.0023	ND	U	0.0017	ND	U	0.0018	ND	U	0.0018		
		1,1,1-Trichloroethane	71-55-6	ND	U	0.0017	ND	U	0.0015	ND	U	0.0023	ND	U	0.0017	ND	U	0.0018	ND	U	0.0018		
		1,1,2,2-Tetrachloroethane	79-34-5	ND	U	0.0017	ND	U	0.0015	ND	U	0.0023	ND	U	0.0017	ND	U	0.0018	ND	U	0.0018		
		1,1,2-Trichloroethane	79-00-5	ND	U	0.0035	ND	U	0.003	ND	U	0.0046	ND	U	0.0034	ND	U	0.0036	ND	U	0.0036		
		1,1-Dichloroethane	75-34-3	ND	U	0.00087	ND	U	0.00075	ND	U	0.0012	ND	U	0.00084	ND	U	0.00091	ND	U	0.00089		
		1,1-Dichloroethene	75-35-4	ND	U	0.0017	ND	U	0.0015	ND	U	0.0023	ND	U	0.0017	ND	U	0.0018	ND	U	0.0018		
		1,1-Dichloropropene	563-58-6	ND	U	0.0017	ND	U	0.0015	ND	U	0.0023	ND	U	0.0017	ND	U	0.0018	ND	U	0.0018		
		1,2,3-Trichlorobenzene	87-61-6	ND	U	0.0017	ND	U	0.0015	ND	U	0.0023	ND	U	0.0017	ND	U	0.0018	ND	U	0.0018		
		1,2,3-Trichloropropane	96-18-4	ND	U	0.0035	ND	U	0.003	ND	U	0.0046	ND	U	0.0034	ND	U	0.0036	ND	U	0.0036		
		1,2,4-Trichlorobenzene	120-82-1	ND	U	0.0017	ND	U	0.0015	ND	U	0.0023	ND	U	0.0017	ND	U	0.0018	ND	U	0.0018		
		1,2,4-Trimethylbenzene	95-63-6	ND	U	0.0035	ND	U	0.003	0.051	J	0.0046	0.0034	J	0.0034	ND	U	0.0036	ND	U	0.0036		
		1,2-Dibromo-3-Chloropropane	96-12-8	ND	U	0.0017	ND	U	0.0015	ND	U	0.0023	ND	U	0.0017	ND	U	0.0018	ND	U	0.0018		
		1,2-Dibromoethane	106-93-4	ND	U	0.0017	ND	U	0.0015	ND	U	0.0023	ND	U	0.0017	ND	U	0.0018	ND	U	0.0018		
		1,2-Dichlorobenzene	95-50-1	ND	U	0.0017	ND	U	0.0015	ND	U	0.0023	ND	U	0.0017	ND	U	0.0018	ND	U	0.0018		
		1,2-Dichloroethane	107-06-2	ND	U	0.0017	ND	U	0.0015	ND	U	0.0023	ND	U	0.0017	ND	U	0.0018	ND	U	0.0018		
		1,2-Dichloroethene	540-59-0	ND	U	0.0017	ND	U	0.0015	ND	U	0.0023	ND	U	0.0017	ND	U	0.0018	ND	U	0.0018		
		1,2-Dichloropropane	78-87-5	ND	U	0.0017	ND	U	0.0015	ND	U	0.0023	ND	U	0.0017	ND	U	0.0018	ND	U	0.0018		
		1,3,5-Trimethylbenzene	108-67-8	ND	U	0.0017	ND	U	0.0015	0.024	J	0.0023	0.0014	J	0.0017	ND	U	0.0018	ND	U	0.0018		
		1,3-Dichlorobenzene	541-73-1	ND	U	0.0017	ND	U	0.0015	ND	U	0.0023	ND	U	0.0017	ND	U	0.0018	ND	U	0.0018		
		1,3-Dichloropropane	142-28-9	ND	U	0.0017	ND	U	0.0015	ND	U	0.0023	ND	U	0.0017	ND	U	0.0018	ND	U	0.0018		
		1,4-Dichlorobenzene	106-46-7	ND	U	0.0017	ND	U	0.0015	ND	U	0.0023	ND	U	0.0017	ND	U	0.0018	ND	U	0.0018		
		2,2-Dichloropropane	594-20-7	ND	U	0.0017	ND	U	0.0015	ND	U	0.0023	ND	U	0.0017	ND	U	0.0018	ND	U	0.0018		
		2-Butanone	78-93-3	ND	U	0.0069	ND	U	0.006	ND	U	0.0093	0.017	J	0.0067	0.035	J	0.0073	0.02	J	0.0071		
		2-Chlorotoluene	95-49-8	ND	U	0.0017	ND	U	0.0015	ND	U	0.0023	ND	U	0.0017	ND	U	0.0018	ND	U	0.0018		
		2-Hexanone	591-78-6	ND	U	0.014	ND	U	0.012	ND	U	0.019	ND	U	0.013	ND	U	0.015	ND	U	0.014		
		4-Chlorotoluene	106-43-4	ND	U	0.0017	ND	U	0.0015	ND	U	0.0023	ND	U	0.0017	ND	U	0.0018	ND	U	0.0018		
		4-Isopropyltoluene	99-87-6	ND	U	0.0017	ND	U	0.0015	0.0059	J	0.0023	0.00077	J	0.0017	ND	U	0.0018	ND	U	0.0018		
		4-Methyl-2-Pentanone	108-10-1	ND	U	0.014	ND	U	0.012	ND	U	0.019	ND	U	0.013	ND	U	0.015	ND	U	0.014		
		Acetone	67-64-1	ND	U	0.014	ND	U	0.012	0.091	--	0.019	0.077	--	0.013	0.021	J	0.015	0.017	J	0.014		
		Benzene	71-43-2	0.0012	J	0.0017	0.00063	J	0.0015	0.05	J	0.0023	0.076	J	0.0017	ND	U	0.0018	ND	U	0.0018		
		Bromobenzene	108-86-1	ND	U	0.0017	ND	U	0.0015	ND	U	0.0023	ND	U	0.0017	ND	U	0.0018	ND	U	0.0018		
		Bromochloromethane	74-97-5	ND	U	0.00087	ND	U	0.00075	ND	U	0.0012	ND	U	0.00084	ND	U	0.00091	ND	U	0.00089		
		Bromodichloromethane	75-27-4	ND	U	0.00087	ND	U	0.00075	ND	U	0.0012	ND	U	0.00084	ND	U	0.00091	ND	U	0.00089		
		Bromoform	75-25-2	ND	U	0.00087	ND	U	0.00075	ND	U	0.0012	ND	U	0.00084	ND	U	0.00091	ND	U	0.00089		
		Bromomethane	74-83-9	ND	U	0.0017	ND	U	0.0015	ND	U	0.0023	ND	U	0.0017	ND	U	0.0018	ND	U	0.0018		
		Carbon Disulfide	75-15-0	ND	U	0.0017	ND	U	0.0015	ND	U	0.0023	ND	U	0.0017	ND	U	0.0018	ND	U	0.0018		
		Carbon Tetrachloride	56-23-5	ND	U	0.0017	ND	U	0.0015	ND	U	0.0023	ND	U	0.0017	ND	U	0.0018	ND	U	0.0018		
		Chlorobenzene	108-90-7	ND	U	0.0017	ND	U	0.0015	ND	U	0.0023	ND	U	0.0017	ND	U	0.0018	ND	U	0.0018		
		Chloroethane	75-00-3	ND	U	0.0035	ND	U	0.003	ND	U	0.0046	ND	U	0.0034	ND	U	0.0036	ND	U	0.0036		
		Chloroform	67-66-3	ND	U	0.00087	ND	U	0.00075	ND	U	0.0012	ND	U	0.00084	ND	U	0.00091	ND	U	0.00089		
		Chloromethane	74-87-3	ND	U	0.0017	ND	U	0.0015	ND	U	0.0023	ND	U	0.0017	ND	U	0.0018	ND	U	0.0018		
		Cis-1,2-Dichloroethene	156-59-2	ND	U	0.0017	ND	U	0.0015	ND	U	0.0023	ND	U	0.0017	ND	U	0.0018	ND	U	0.0018		
		Cis-1,3-Dichloropropene	10061-01-5	ND	U	0.0035	ND	U	0.003	ND	U	0.0046	ND	U	0.0034	ND	U	0.0036	ND	U	0.0036		
		Dibromochloromethane	124-48-1	ND	U	0.0017	ND	U	0.0015	ND	U	0.0023	ND	U	0.0017	ND	U	0.0018	ND	U	0.0018		
		Dibromomethane	74-95-3	ND	U	0.0035	ND	U	0.003	ND	U	0.0046	ND	U	0.0034	ND	U	0.0036	ND	U	0.0036		
		Dichlorodifluoromethane	75-71-8	ND	U	0.0017	ND	U	0.0015	ND	U	0.0023	ND	U	0.0017	ND	U	0.0018	ND	U	0.0018		
		Ethylbenzene	100-41-4	ND	U	0.0017	ND	U	0.0015	0.032	J	0.0023	0.005	J	0.0017	ND	U	0.0018	ND	U	0.0018		
		Hexachloro-1,3-Butadiene	87-68-3	ND	U	0.0017	ND	U	0.0015	ND	U	0.0023	ND	U	0.0017	ND	U	0.0018	ND	U	0.0018		
		Isopropylbenzene	98-82-8	ND	U	0.0017	ND	U	0.0015	0.011	--	0.0023	ND	U	0.0017	ND	U	0.0018	ND	U	0.0018		
		m- & p-Xylenes	179601-23-1	0.0019	J	0.0035	0.0011	J	0.003	0.083	J	0.0046	0.0056	J	0.0034	ND	U	0.0036	ND	U	0.0036		
		Methyl tert-Butyl Ether	1634-04-4																				

Table H-5
Field Duplicate Sample Results

				Boring ID:			KAFB-106S1			KAFB-106S1			KAFB-106S2			KAFB-106S2			KAFB-106S3			KAFB-106S3			KAFB-106S4		
				Sample Date:			12/5/2018			12/5/2018			11/16/2018			11/16/2018			11/20/2018			11/20/2018			11/5/2018		
				Sample Depth (ft bgs):			510			510			510			510			459			459			504		
				Sample Type:			REG			Field Duplicate			REG			Field Duplicate			REG			Field Duplicate			REG		
Parameter	Analytical Method	Chemical Name	CAS	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD			
Moisture	ASTM D2216 (%) VOCs	Vinyl Chloride	75-01-4	ND	U	0.0035	ND	U	0.003	ND	U	0.0046	ND	U	0.0034	ND	U	0.0036	ND	U	0.0036	ND	UJ	0.0031			
		NA	NA	11.2	--	--	9.3	--	--	10.1	--	--	11.3	--	--	3.3	--	--	3.6	--	--	19.7	--	--			

Table H-5
Field Duplicate Sample Results

		Boring ID:		KAFB-106S4			KAFB-106S5			KAFB-106S5			KAFB-106S7			KAFB-106S7			KAFB-106S8			KAFB-106S8			KAFB-106S9			KAFB-106S9					
		Sample Date:		11/5/2018			10/23/2018			10/23/2018			1/22/2019			1/22/2019			1/18/2019			1/18/2019			10/19/2018			10/19/2018					
		Sample Depth (ft bgs):		504			467			467			506			506			514			514			490			490					
		Sample Type:		Field Duplicate			REG			Field Duplicate			REG			Field Duplicate			REG			Field Duplicate			REG			Field Duplicate					
Parameter	Analytical Method	Chemical Name	CAS	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD						
TPH	Method SW8015D (mg/kg)	TPH-MRO (C20-C38)	MOIL	ND	U	25	ND	U	22	ND	U	23	ND	U	23	ND	U	24	ND	U	26	ND	U	27	12	J	24	12	J	25			
		TPH-DRO (C10-C28)	TPH-DRO	ND	U	4.3	5.6	J	3.8	4.7	J	4	ND	U	4	ND	U	4.1	ND	U	4.4	ND	U	4.5	1,900	J	4.1	1,900	--	4.3			
		TPH-GRO (C6-C10)	TPH-GRO	ND	U	1.1	ND	U	1.6	ND	UJ	1.3	ND	U	1.3	ND	U	1.1	ND	U	1.4	ND	U	1.4	2,300	J	210	1,400	J	120			
Moisture	ASTM D2216 (%) TPH	NA	NA	11.3	--	--	3.8	--	--	3.7	--	--	9.4	--	--	8.8	--	--	11.6	--	--	12.5	--	--	9.9	--	--	9.6	--	--			
EDB	Method SW8011 (mg/kg) ^a	1,2-Dibromoethane	106-93-4	ND	U	0.000043	ND	U	0.000039	ND	U	0.00004	ND	U	0.000042	ND	U	0.000042	ND	U	0.000045	ND	U	0.000044	0.29	J	0.0084	0.13	J	0.0085			
Moisture	ASTM D2216 (%) EDB	NA	NA	11.3	--	--	4.1	--	--	4.2	--	--	10.3	--	--	10.2	--	--	16.5	--	--	14.6	--	--	9.9	--	--	10.4	--	--			
VOCs	Method SW8260B (mg/kg) ^a	1,1,1,2-Tetrachloroethane	630-20-6	ND	U	0.0013	ND	U	0.0017	ND	U	0.0017	ND	U	0.0015	ND	U	0.0017	ND	U	0.0016	ND	U	0.0014	ND	U	4.3	ND	U	1.3			
		1,1,1-Trichloroethane	71-55-6	ND	U	0.0013	ND	U	0.0017	ND	U	0.0017	ND	U	0.0015	ND	U	0.0017	ND	U	0.0016	ND	U	0.0014	ND	U	2.2	ND	U	0.67			
		1,1,2,2-Tetrachloroethane	79-34-5	ND	U	0.0013	ND	U	0.0017	ND	U	0.0017	ND	U	0.0015	ND	U	0.0017	ND	U	0.0016	ND	U	0.0014	ND	U	4.3	ND	U	1.3			
		1,1,2-Trichloroethane	79-00-5	ND	U	0.0026	ND	U	0.0034	ND	U	0.0034	ND	U	0.0031	ND	U	0.0034	ND	U	0.0032	ND	U	0.0029	ND	U	4.3	ND	U	1.3			
		1,1-Dichloroethane	75-34-3	ND	U	0.00066	ND	U	0.00085	ND	U	0.00085	ND	U	0.00077	ND	U	0.00085	ND	U	0.00081	ND	U	0.00072	ND	U	8.6	ND	U	2.7			
		1,1-Dichloroethene	75-35-4	ND	U	0.0013	ND	U	0.0017	ND	U	0.0017	ND	U	0.0015	ND	U	0.0015	ND	U	0.0016	ND	U	0.0014	ND	U	8.6	ND	U	2.7			
		1,1-Dichloropropene	563-58-6	ND	U	0.0013	ND	U	0.0017	ND	U	0.0017	ND	U	0.0015	ND	U	0.0017	ND	U	0.0016	ND	U	0.0014	ND	U	8.6	ND	U	2.7			
		1,2,3-Trichlorobenzene	87-61-6	ND	U	0.0013	ND	U	0.0017	ND	U	0.0017	ND	U	0.0015	ND	U	0.0017	ND	U	0.0016	ND	U	0.0014	ND	U	4.3	ND	U	1.3			
		1,2,3-Trichloropropane	96-18-4	ND	U	0.0026	ND	U	0.0034	ND	U	0.0034	ND	U	0.0031	ND	U	0.0034	0.00099	J	0.0032	ND	U	0.0029	ND	U	8.6	ND	U	2.7			
		1,2,4-Trichlorobenzene	120-82-1	ND	U	0.0013	ND	U	0.0017	ND	U	0.0017	ND	U	0.0015	ND	U	0.0017	ND	U	0.0016	ND	U	0.0014	ND	U	8.6	ND	U	2.7			
		1,2,4-Trimethylbenzene	95-63-6	0.0029	J	0.0026	ND	U	0.0034	ND	U	0.0034	0.009	--	0.0031	0.003	J	0.0034	ND	U	0.0032	ND	U	0.0029	68	J	4.3	42	J	1.3			
		1,2-Dibromo-3-Chloropropane	96-12-8	ND	U	0.0013	ND	U	0.0017	ND	U	0.0017	ND	U	0.0015	ND	U	0.0017	ND	U	0.0016	ND	U	0.0014	ND	U	17	ND	U	5.4			
		1,2-Dibromoethane	106-93-4	ND	U	0.0013	ND	U	0.0017	ND	U	0.0017	ND	U	0.0015	ND	U	0.0017	ND	U	0.0016	ND	U	0.0014	ND	U	4.3	ND	U	1.3			
		1,2-Dichlorobenzene	95-50-1	ND	U	0.0013	ND	U	0.0017	ND	U	0.0017	ND	U	0.0015	ND	U	0.0017	ND	U	0.0016	ND	U	0.0014	ND	U	17	ND	U	5.4			
		1,2-Dichloroethane	107-06-2	ND	U	0.0013	ND	U	0.0017	ND	U	0.0017	ND	U	0.0015	ND	U	0.0017	ND	U	0.0016	ND	U	0.0014	ND	U	4.3	ND	U	1.3			
		1,2-Dichloroethene	540-59-0	ND	U	0.0013	ND	U	0.0017	ND	U	0.0017	ND	U	0.0015	ND	U	0.0017	ND	U	0.0016	ND	U	0.0014	ND	U	4.3	ND	U	1.3			
		1,2-Dichloropropane	78-87-5	ND	U	0.0013	ND	U	0.0017	ND	U	0.0017	ND	U	0.0015	ND	U	0.0017	ND	U	0.0016	ND	U	0.0014	ND	U	8.6	ND	U	2.7			
		1,3,5-Trimethylbenzene	108-67-8	0.0013	J	0.0013	ND	U	0.0017	ND	U	0.0017	ND	U	0.0017	0.0033	J	0.0015	ND	U	0.0017	ND	U	0.0016	ND	U	0.0014	31	J	4.3	17	J	1.3
		1,3-Dichlorobenzene	541-73-1	ND	U	0.0013	ND	U	0.0017	ND	U	0.0017	ND	U	0.0015	ND	U	0.0017	ND	U	0.0016	ND	U	0.0014	ND	U	8.6	ND	U	2.7			
		1,3-Dichloropropane	142-28-9	ND	U	0.0013	ND	U	0.0017	ND	U	0.0017	ND	U	0.0015	ND	U	0.0017	ND	U	0.0016	ND	U	0.0014	ND	U	4.3	ND	U	1.3			
		1,4-Dichlorobenzene	106-46-7	ND	U	0.0013	ND	U	0.0017	ND	U	0.0017	ND	U	0.0015	ND	U	0.0017	ND	U	0.0016	ND	U	0.0014	ND	U	4.3	ND	U	1.3			
		2,2-Dichloropropane	594-20-7	ND	U	0.0013	ND	U	0.0017	ND	U	0.0017	ND	U	0.0015	ND	U	0.0017	ND	U	0.0016	ND	U	0.0014	ND	U	4.3	ND	U	1.3			
		2-Butanone	78-93-3	ND	U	0.0052	ND	U	0.0068	0.0042	J	0.0068	ND	U	0.0062	ND	U	0.0068	ND	U	0.0065	ND	U	0.0058	ND	U	34	ND	U	11			
		2-Chlorotoluene	95-49-8	ND	U	0.0013	ND	U	0.0017	ND	U	0.0017	ND	U	0.0015	ND	U	0.0017	ND	U	0.0016	ND	U	0.0014	ND	U	4.3	ND	U	1.3			
		2-Hexanone	591-78-6	ND	U	0.01	ND	U	0.014	ND	U	0.014	ND	U	0.012	ND	U	0.014	0.015	J	0.013	ND	U	0.012	ND	U	34	ND	U	11			
		4-Chlorotoluene	106-43-4	ND	U	0.0013	ND	U	0.0017	ND	U	0.0017	ND	U	0.0015	ND	U	0.0017	ND	U	0.0016	ND	U	0.0014	ND	U	4.3	ND	U	1.3			
		4-Isopropyltoluene	99-87-6	0.00043	J	0.0013	ND	U	0.0017	ND	U	0.0017	ND	U	0.0017	0.0011	J	0.0015	0.00091	J	0.0017	ND	U	0.0016	ND	U	0.0014	5.4	J	4.3	3.4	J	1.3
		4-Methyl-2-Pentanone	108-10-1	ND	U	0.01	ND	U	0.014	ND	U	0.014	ND	U	0.012	ND	U	0.012	ND	U	0.014	ND	U	0.013	ND	U	0.012	ND	U	34	ND	U	11
		Acetone	67-64-1	0.089	--	0.01	ND	U	0.014	ND	U	0.014	ND	U	0.012	ND	U	0.012	0.036	--	0.014	ND	U	0.013	ND	U	0.012	ND	U	69	50	--	22
		Benzene	71-43-2	0.0033	J	0.0013	ND	U	0.0017	0.0006	J	0.0017	0.0011	J	0.0015	0.00088	J	0.0017	0.0014	J	0.0016	0.0014	J	0.0014	J	0.0014	57	J	8.6	27	J	2.7	
		Bromobenzene	108-86-1	ND	U	0.0013	ND	U	0.0017	ND	U	0.0017	ND	U	0.0015	ND	U	0.0017	ND	U	0.0016	ND	U	0.0014	ND	U	2.2	ND	U	0.67			
		Bromochloromethane	74-97-5	ND	U	0.00066	ND	U	0.00085	ND	U	0.00085																					

Table H-5
Field Duplicate Sample Results

		Boring ID:			KAFB-106S4			KAFB-106S5			KAFB-106S5			KAFB-106S7			KAFB-106S7			KAFB-106S8			KAFB-106S8			KAFB-106S9			KAFB-106S9		
		Sample Date:			11/5/2018			10/23/2018			10/23/2018			1/22/2019			1/22/2019			1/18/2019			1/18/2019			10/19/2018			10/19/2018		
		Sample Depth (ft bgs):			504			467			467			506			506			514			514			490			490		
		Sample Type:			Field Duplicate			REG			Field Duplicate			REG			Field Duplicate			REG			Field Duplicate			REG			Field Duplicate		
Parameter	Analytical Method	Chemical Name			CAS	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD		
		Vinyl Chloride			75-01-4	ND	U	0.0026	ND	U	0.0034	ND	U	0.0034	ND	U	0.0031	ND	U	0.0034	ND	U	0.0032	ND	U	0.0029	ND	U	4.3		
Moisture	ASTM D2216 (%) VOCs	NA			NA	11.3	--	--	4.1	--	--	4.2	--	--	10.3	--	--	10.2	--	--	16.5	--	--	14.6	--	--	9.9	--	--	10.4	

Table H-5
Field Duplicate Sample Results

Parameter	Analytical Method	Chemical Name	CAS	Boring ID:			KAFB-106V1			KAFB-106V1			KAFB-106V2			KAFB-106V2		
				Sample Date:			12/19/2018			12/19/2018			12/14/2018			12/14/2018		
				Sample Depth (ft bgs):			285			285			270			270		
				Sample Type:			REG			Field Duplicate			REG			Field Duplicate		
				Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD
TPH	Method SW8015D (mg/kg)	TPH-MRO (C20-C38)	MOIL	ND	U	24	ND	U	22	ND	UJ	29	ND	UJ	29	ND	UJ	29
		TPH-DRO (C10-C28)	TPH-DRO	270	--	4.1	290	--	3.8	8.8	J	4.9	7.5	J	5			
		TPH-GRO (C6-C10)	TPH-GRO	380	J	140	75	J	2.6	8.2	--	1.7	6.5	--	1.8			
Moisture	ASTM D2216 (%) TPH	NA	NA	3.3	--	--	3.8	--	--	19.9	--	--	20.3	--	--			
EDB	Method SW8011 (mg/kg) ^a	1,2-Dibromoethane	106-93-4	ND	U	0.00039	ND	UJ	0.00079	0.0006	J	0.00048	0.00046	--	0.000047			
Moisture	ASTM D2216 (%) EDB	NA	NA	3.3	--	--	3.6	--	--	20.8	--	--	19.6	--	--			
VOCs	Method SW8260B (mg/kg) ^a	1,1,1,2-Tetrachloroethane	630-20-6	ND	U	1.4	ND	U	0.95	ND	U	1.1	ND	U	0.11			
		1,1,1-Trichloroethane	71-55-6	ND	U	0.7	ND	U	0.48	ND	U	0.53	ND	U	0.055			
		1,1,2,2-Tetrachloroethane	79-34-5	ND	U	1.4	ND	U	0.95	ND	U	1.1	ND	U	0.11			
		1,1,2-Trichloroethane	79-00-5	ND	U	1.4	ND	U	0.95	ND	U	1.1	0.043	J	0.11			
		1,1-Dichloroethane	75-34-3	ND	U	2.8	ND	U	1.9	ND	U	2.1	ND	U	0.22			
		1,1-Dichloroethene	75-35-4	ND	U	2.8	ND	U	1.9	ND	U	2.1	ND	U	0.22			
		1,1-Dichloropropene	563-58-6	ND	U	2.8	ND	U	1.9	ND	U	2.1	ND	U	0.22			
		1,2,3-Trichlorobenzene	87-61-6	ND	U	1.4	ND	U	0.95	ND	U	1.1	0.055	J	0.11			
		1,2,3-Trichloropropane	96-18-4	ND	U	2.8	ND	U	1.9	ND	U	2.1	ND	U	0.22			
		1,2,4-Trichlorobenzene	120-82-1	ND	U	2.8	ND	U	1.9	ND	U	2.1	ND	U	0.22			
		1,2,4-Trimethylbenzene	95-63-6	1.7	J	1.4	0.87	J	0.95	ND	U	1.1	0.048	J	0.11			
		1,2-Dibromo-3-Chloropropane	96-12-8	ND	U	5.6	ND	U	3.8	ND	U	4.2	ND	U	0.44			
		1,2-Dibromoethane	106-93-4	ND	U	1.4	ND	U	0.95	ND	U	1.1	0.046	J	0.11			
		1,2-Dichlorobenzene	95-50-1	ND	U	5.6	ND	U	3.8	ND	U	4.2	ND	U	0.44			
		1,2-Dichloroethane	107-06-2	ND	U	1.4	ND	U	0.95	ND	U	1.1	ND	U	0.11			
		1,2-Dichloroethene	540-59-0	ND	U	1.4	ND	U	0.95	ND	U	1.1	0.03	J	0.11			
		1,2-Dichloropropane	78-87-5	ND	U	2.8	ND	U	1.9	ND	U	2.1	ND	U	0.22			
		1,3,5-Trimethylbenzene	108-67-8	0.94	J	1.4	0.53	J	0.95	ND	U	1.1	ND	U	0.11			
		1,3-Dichlorobenzene	541-73-1	ND	U	2.8	ND	U	1.9	ND	U	2.1	ND	U	0.22			
		1,3-Dichloropropane	142-28-9	ND	U	1.4	ND	U	0.95	ND	U	1.1	ND	U	0.11			
		1,4-Dichlorobenzene	106-46-7	ND	U	1.4	ND	U	0.95	ND	U	1.1	0.034	J	0.11			
		2,2-Dichloropropane	594-20-7	ND	U	1.4	ND	U	0.95	ND	U	1.1	ND	U	0.11			
		2-Butanone	78-93-3	ND	U	11	ND	U	7.6	ND	U	8.4	7.3	--	0.87			
		2-Chlorotoluene	95-49-8	ND	U	1.4	ND	U	0.95	ND	U	1.1	ND	U	0.11			
		2-Hexanone	591-78-6	ND	U	11	ND	U	7.6	ND	U	8.4	ND	U	0.87			
		4-Chlorotoluene	106-43-4	ND	U	1.4	ND	U	0.95	ND	U	1.1	0.031	J	0.11			
		4-Isopropyltoluene	99-87-6	ND	U	1.4	ND	U	0.95	ND	U	1.1	ND	U	0.11			
		4-Methyl-2-Pentanone	108-10-1	ND	U	11	ND	U	7.6	ND	U	8.4	0.34	J	0.87			
		Acetone	67-64-1	ND	U	22	ND	U	15	76	--	17	78	--	17			
		Benzene	71-43-2	ND	U	2.8	ND	U	1.9	ND	U	2.1	0.26	J	0.22			
		Bromobenzene	108-86-1	ND	U	0.7	ND	U	0.48	ND	U	0.53	0.039	J	0.055			
		Bromochloromethane	74-97-5	ND	U	2.8	ND	U	1.9	ND	U	2.1	ND	U	0.22			
		Bromodichloromethane	75-27-4	ND	U	1.4	ND	U	0.95	ND	U	1.1	ND	U	0.11			
		Bromoform	75-25-2	ND	U	2.8	ND	U	1.9	ND	U	2.1	ND	U	0.22			
		Bromomethane	74-83-9	ND	U	2.8	ND	U	1.9	ND	U	2.1	ND	U	0.22			
		Carbon Disulfide	75-15-0	ND	U	2.8	ND	U	1.9	ND	U	2.1	ND	U	0.22			
		Carbon Tetrachloride	56-23-5	ND	U	0.7	ND	U	0.48	ND	U	0.53	0.026	J	0.055			
		Chlorobenzene	108-90-7	ND	U	1.4	ND	U	0.95	ND	U	1.1	ND	U	0.11			
		Chloroethane	75-00-3	ND	U	2.8	ND	U	1.9	ND	U	2.1	ND	U	0.22			
		Chloroform	67-66-3	ND	U	2.8	ND	U	1.9	ND	U	2.1	ND	U	0.22			
		Chloromethane	74-87-3	ND	U	2.8	ND	U	1.9	ND	U	2.1	ND	U	0.22			
		Cis-1,2-Dichloroethene	156-59-2	ND	U	1.4	ND	U	0.95	ND	U	1.1	0.03	J	0.11			
		Cis-1,3-Dichloropropene	10061-01-5	ND	U	1.4	ND	U	0.95	ND	U	1.1	ND	U	0.11			
		Dibromochloromethane	124-48-1	ND	U	1.4	ND	U	0.95	ND	U	1.1	ND	U	0.11			
		Dibromomethane	74-95-3	ND	U	1.4	ND	U	0.95	ND	U	1.1	ND	U	0.11			
		Dichlorodifluoromethane	75-71-8	ND	U	1.4	ND	U	0.95	ND	U	1.1	ND	U	0.11			
		Ethylbenzene	100-41-4	ND	U	1.4	ND	U	0.95	ND	U	1.1	0.06	J	0.11			
		Hexachloro-1,3-Butadiene	87-68-3	ND	U	2.8	ND	U	1.9	ND	U	2.1	ND	U	0.22			
		Isopropylbenzene	98-82-8	ND	U	1.4	ND	U	0.95	ND	U	1.1	ND	U	0.11			
		m- & p-Xylenes	179601-23-1	2.5	J	2.8	1.3	J	1.9	ND	U	2.1	ND	U	0.22			
		Methyl tert-Butyl Ether	1634-04-4	ND	U	2.8	ND	U	1.9	ND	U	2.1	ND	U	0.22			
		Methylene Chloride	75-09-2	ND	U	2.8	ND	U	1.9	ND	U	2.1	ND	U	0.22			
		Naphthalene	91-20-3	1.9	J	1.4	1.4	J	0.95	0.69	J	1.1	ND	U	0.11			
		n-Butylbenzene	104-51-8	ND	U	1.4	ND	U	0.95	ND	U	1.1	ND	U	0.11			
		n-Propylbenzene	103-65-1	ND	U	2.8	ND	U	1.9	ND	U	2.1	ND	U	0.22			
		o-Xylene	95-47-6	1.2	J	1.4	0.69	J	0.95	ND	U	1.1	0.057	J	0.11			
		sec-Butylbenzene	135-98-8	ND	U	1.4	0.5	J	0.95	ND	U	1.1	ND	U	0.11			
		Styrene	100-42-5	ND	U	1.4	ND	U	0.95	ND	U	1.1	ND	U	0.11			
		tert-Butylbenzene	98-06-6	ND	U	1.4	ND	U	0.95	ND	U	1.1	0.039	J	0.11			
		Tetrachloroethene	127-18-4	ND	U	1.4	ND	U	0.95	ND	U	1.1	ND	U	0.11			
		Toluene	108-88-3	0.99	J	1.4	0.51	J	0.95	1.7	J	1.1	0.69	J	0.11			
		Trans-1,2-Dichloroethene	156-60-5	ND	U	2.8	ND	U	1.9	ND	U	2.1	ND	U	0.22			
		Trans-1,3-Dichloropropene	10061-02-6	ND	U	2.8	ND	U	1.9	ND	U	2.1	ND	U	0.22			
		Trichloroethene	79-01-6	ND	U	1.4	ND	U	0.95	ND	U	1.1	ND	U	0.11			
		Trichlorofluoromethane	75-69-4	ND	U	2.8	ND	U	1.9	ND	U	2.1	ND	U	0.22			

Table H-5
Field Duplicate Sample Results

		Boring ID:			KAFB-106V1			KAFB-106V1			KAFB-106V2			KAFB-106V2		
		Sample Date:			12/19/2018			12/19/2018			12/14/2018			12/14/2018		
		Sample Depth (ft bgs):			285			285			270			270		
		Sample Type:			REG			Field Duplicate			REG			Field Duplicate		
		Parameter	Analytical Method	Chemical Name	CAS	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual	LOD	Result	Val Qual
		Vinyl Chloride	75-01-4	ND	U	1.4	ND	U	0.95	ND	U	1.1	ND	U	0.11	
Moisture	ASTM D2216 (%) VOCs	NA	NA	3.3	--	--	3.6	--	--	20.8	--	--	19.6	--	--	

Table H-5
Field Duplicate Sample Results

a Results reported by the laboratory in microgram per kilogram but converted to mg/kg for presentation.
% = percent
AFB = Air Force Base
BFF = Bulk Fuels Facility
bgs = below ground surface
CAS = chemical abstract service
DRO = diesel range organics
EDB = ethylene dibromide (1,2-dibromoethane)
ft = foot/feet
GRO = gasoline range organics
ID = identification
KAFB = Kirtland Air Force Base
LOD = limit of detection
mg/kg = milligram per kilogram
MRO = motor oil organics
NA = not applicable
ND = not detected
NMED= New Mexico Environment Department
REG = normal field sample
RSL = regional screening level
SWMU = Solid Waste Management Unit
TPH = total petroleum hydrocarbons
Val Qual = validation qualifier
VOC = volatile organic compound
Shading = detected concentrations above the detection limit
Bold/Shading = reported concentrations exceed either the EPA RSL or NMED soil screening level.
Val Quals based on independent data validation
J = Qualifier denotes the analyte was positively identified, but the associated numerical value is estimated.
U = Qualifier denotes the analyte was analyzed but not detected above the detection limit. The value associated with the U-qualifier is the LOD.
-- = Validation qualifier not assigned.
— = Compound not analyzed for.

Table H-6
Technical Data Completeness - Source Zone Characterization

Analytical Parameter	Field/Field Duplicate Sample Analytes	Quality Control Sample Analytes (TB, RB)	Qualified Analytes	Percent Technical Completeness^a
VOCs (SW8260B)	5,478	3,762	632	100
Ethylene dibromide (SW8011)	83	47	7	100
Total Petroleum Hydrocarbons (SW8015D) ^b	291	97	96	100

a. Percent technical completeness including analytes qualified as estimated data. No data were rejected.

b. Total petroleum hydrocarbons (SW8015D) = gasoline range organics, diesel range organics, and motor oil. Trip blanks analyzed for gasoline range organics only.

RB = equipment rinse blank

TB = trip blank

VOC = volatile organic compound