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APPENDIX I

DESCRIPTION OF ORGANICS AND INORGANICS DATA REVIEW PACKAGES

The purpose of Appendix I is to familiarize the reader with a model for data review deliverables. This appendix consists of the following items:

- o A description of the data reporting format,
- o An example of a data review summary, and
- o Example data review forms.

Please note that the example forms are designed for the validation of Contract Laboratory Program (CLP) data packages. An example form is included for each analytical fraction (volatiles, semivolatiles, pesticide/Aroclors and metals) and for samples from soil/sediment and aqueous matrices. These forms nevertheless include the necessary information for the review of most types of data (analytical results, sample quantitation/detection limits, data qualifiers, etc.) not associated with the CLP.

1. DATA REPORTING FORMATS

Whenever an analytical laboratory is requested to analyze field samples for a specific site, the RPM (in consultation with the technical project team) must ensure that the laboratory will provide adequate documentation to support all current and future uses of the data. Potential uses of the data can include data validation, monitoring, modeling, risk assessment, site characterization, Record of Decision defense, enforcement, and litigation.

Data packages produced by analytical laboratories should contain all the documents that were produced or used by the laboratory for that particular analysis. The required documents should include a narrative (detailing the exact method performed, deviations from the method, problems encountered, and problem resolution), chain-of-custody records, laboratory logbook pages, and raw data and tabulated summary forms for all standards, quality control and field samples.

The documents should be organized in a logical manner and the entire data package should be paginated. Generally, the laboratory should be required to produce a data package with documents ordered in the following manner:

- 1) Narrative
- 2) Tabulated summary forms for laboratory standards and quality control samples (in chronological order by type of quality control sample/standard by date of analysis by instrument)
- 3) Tabulated summary forms for field sample results (in increasing RAS, SAS, or project sample number order)
- 4) Raw data for field samples (in increasing RAS, SAS, or project sample number order)
- 5) Raw data for laboratory standards and quality control samples (in chronological order by type of quality control sample/standard by date of analysis by instrument)
- 6) Laboratory logbook pages
- 7) Chain-of-custody records

APPENDIX I (continued)

It is often convenient to require that the laboratory data package resemble as closely as possible the data packages required by the current CLP RAS SOWs for organics and inorganics, that the tabulated summary forms provided in those SOWs be utilized and modified appropriately, and that the data qualifiers in those SOWs be applied to the data as appropriate. The following sections describe specific requirements for the content of each document contained in the laboratory data package.

NARRATIVE:

A narrative must be provided describing the analytical methods and exact procedures performed by the laboratory, as well as any deviations from the method. Problems encountered during analysis, problem resolution and any factors which may affect the validity of the data must be addressed. The narrative must include the laboratory name and RAS, SAS, or project sample numbers cross-referenced to the laboratory sample identification numbers, and must be signed and dated by the laboratory manager.

Any telephone communications between the laboratory and sampling personnel (or other parties outside of the laboratory) to resolve sampling discrepancies or analytical problems must be documented in detail on telephone communication logs. Those telephone logs must explicitly detail the problems requiring resolution, the agreed to resolution, and the names and affiliations of the communicating parties. All telephone logs must be appended to the narrative.

An example calculation of a positive hit and a detection/quantitation limit for each type of sample analysis must be provided. All equations, dilution factors and information required to reproduce the laboratory results must be provided.

TABULATED SUMMARY FORMS:

Laboratory Standards and Quality Control Samples

Tabulated summary forms must be provided for all laboratory standards, tunes, blanks, duplicates, spikes, and any other types of laboratory quality control samples/standards. The tabulated summary forms must contain information pertinent to the type of laboratory quality control sample/standard which was analyzed. Typical entries include: concentrations spiked, concentrations detected, spike compound names, results of statistical calculations (%R, %D, RPD, RSD, CV, RRF, SD, etc.), sample identification numbers, dates/times of analysis, instrument IDs, lab file IDs, and QC limits.

The exact format of each tabulated summary form will depend on the particular analysis method requested and the quality control procedures specified in that method. However, comprehensive tabulated summary forms must be prepared for all quality control samples/standards analyzed by the laboratory. For example, typical tabulated summary forms for volatile organics analyses include but are not limited to:

Surrogate results: Tabulate the sample identification numbers, surrogate compounds added, concentration added, percent recoveries, and QC limits for all standards, blanks, quality control samples and field samples. Flag outliers.

Matrix spike and matrix spike duplicate results: Tabulate the matrix spike compounds added, concentration added, percent recoveries and relative percent differences for the spiked compounds, and QC limits. Flag outliers. List the sample identification numbers. Results for

APPENDIX I (continued)

all non-spike compounds must be tabulated on the form used to summarize field sample results.

Method/laboratory blanks: Tabulate the sample identification numbers, lab file IDs, and time analyzed for field samples and matrix spike samples which pertain to each blank on a separate form. The form must also contain the GC column, instrument ID, laboratory sample identification number, lab file ID, and date/time of analysis for the blank itself. Results for each blank must also be tabulated on the form used to summarize field sample results.

Tuning results: Tabulate the m/e, ion abundance criteria, and percent relative abundances and list the tune compound name, instrument ID, lab file ID, and date/time of injection which pertain to each tune analysis on a separate form. The form must also contain tabulated sample identification numbers, lab file IDs, and date/time of analysis for all field samples, matrix spike samples, blanks, and standards which pertain to that tune. Flag outliers.

Initial calibration results: Tabulate the target compound names, relative response factors for each target and surrogate compound at each standard concentration, mean relative response factors and percent relative standard deviations for all target and surrogate compounds, and QC limits for each initial calibration on a separate form. The form must also contain the concentration of the calibration standards, instrument ID, lab file IDs, and dates/times of standard analyses for that initial calibration. Flag outliers.

Continuing calibration results: Tabulate the target compound names, mean relative response factors from initial calibration, relative response factors from continuing calibration, percent differences, and QC limits for all target and surrogate compounds for each continuing calibration on a separate form. The form must also contain the concentration of the continuing calibration standard, instrument ID, lab file ID, and dates/times of initial and continuing calibration standard analyses which pertain to that continuing calibration. Flag outliers.

Internal standard results: Tabulate the sample identification numbers, internal standard compound names, QC limits, retention times and area counts of the quantitation ion for each internal standard compound in the continuing calibration standard and all field samples, matrix spike samples, and blanks which pertain to that continuing calibration on a separate form. The form must also contain the instrument ID, lab file ID, and date/time of continuing calibration standard analysis. Flag outliers.

MDL study results: Tabulate the target compound names, concentrations spiked and detected for each MDL spike analysis, and the standard deviation and calculated MDL for each target compound. (Note: The narrative must explain the MDL procedure utilized to generate the values. The formula and associated constant values utilized in the calculation of the MDL for each analyte must be provided. The column, instrument ID, trap composition, and operating conditions must be clearly displayed on the raw data.)

Field Samples

The exact format of the tabulated summary form for each field sample will depend on the particular analysis method requested. However, comprehensive tabulated summary forms must be prepared for each field sample analyzed by the laboratory. At a minimum, the target compound names, concentration units, positive hits and numerical detection/quantitation limits and any laboratory qualifier flags for each target compound must be tabulated on a separate form. Definitions must be provided for all qualifier flags used by the laboratory. For each

APPENDIX I (continued)

sample, the tabulated form must also contain the RAS, SAS, or project sample identification number, laboratory name, laboratory sample ID, lab file ID, sample matrix type, and level of analysis (low, medium, high). The percent moisture/solids, weights and volumes of sample prepared/purged/extracted/digested/analyzed, initial and final extract/digest and extract clean-up volumes, injection volume, clean-ups performed, dilution factor, measured pH, and dates that sample was received/extracted/digested/analyzed should be included as appropriate to the analysis method.

RAW DATA:

Raw data must be provided by the laboratory for all laboratory quality control samples, blanks, spikes, duplicates, standards, and field samples. The exact format and content of the raw data will depend on the particular analysis method requested. However, any and all instrument printouts, strip chart recordings, chromatograms, quantitation reports, mass spectra and other types of raw data generated by the laboratory for a particular project must be provided in the data package. Typical raw data for organic GC/MS analyses includes but is not limited to:

- o Reconstructed total ion chromatograms,
- o Instrument quantitation reports containing the following information: laboratory sample identification number, RAS, SAS or project sample number, date and time of analysis, RT and/or scan number of quantitation ion with measured area, analyte concentration, copy of area table from data system, GC/MS instrument ID, lab file ID, column, trap composition, and operating conditions,
- o Raw and enhanced mass spectra for all positive field sample results and daily continuing calibration standard reference spectra for all positive field sample results,
- o Mass spectra and three library searched best-match mass spectra for all tentatively identified compounds reported, and
- o Instrument normalized mass listing and the mass spectrum for each tune.

Typical raw data for inorganic analyses includes but is not limited to:

- o Instrument printouts and strip chart recordings containing the following information: laboratory sample identification number, RAS, SAS or project sample number, date and time of analysis, absorbance/emissions values, analyte concentration, instrument ID, lab file ID, and operating conditions, and
- o Standard curve raw data, plotted standard curves, linear regression equations, and correlation coefficients.

LABORATORY LOGBOOK PAGES:

Copies of standards preparation logs, sample preparation/extraction/digestion logs, sample analysis run logs, personal logs, and any hand written project-specific notes must be included. The initial and final volumes of sample prepared/purged/extracted/digested, initial

APPENDIX I (continued)

and final extract/digest and extract clean-up volumes, injection volumes, and dilution factors must be clearly labelled.

CHAIN-OF-CUSTODY RECORDS:

All chain-of-custody records provided to the laboratory during sample shipment or generated by the laboratory during sample receipt, storage, preparation, and analysis must be included. Chain-of-custody records include but are not limited to: signed and dated field chain-of-custody forms, signed and dated shipping airbills, sample tags, SAS packing lists, RAS Traffic Reports, internal laboratory receiving records, and internal laboratory sample/extract/digest transfer records.

APPENDIX I (Continued)

2. DATA REVIEW SUMMARY

ORGANIC DATA SUMMARY FORMS UTILIZED BY REGION III IN THE CLP

DATE:

SUBJECT:

FROM:

TO:

THRU:

OVERVIEW

Case consisted of four (4) low level water and two (2) low level soil samples, submitted for full organic analyses. Included in this data set was one (1) equipment blank and one (1) trip blank. The trip blank was analyzed for volatiles only. The samples were analyzed as a Contract Laboratory Program (CLP) Routine Analytical Service (RAS).

SUMMARY

All samples were successfully analyzed for all target compounds with the exception of 2-Butanone and 2-Hexanone in the volatile fraction. All remaining instrument and method sensitivities were according to the Contract Laboratory Program (CLP) Routine Analytical Service (RAS) protocol.

MAJOR PROBLEM

The response factors (RF) for 2-Butanone and 2-Hexanone were less than 0.05 in one of the continuing volatile calibration. The quantitation limits for this compound in the affected samples were qualified unreliable, "R". (See Table I in Appendix F for the affected samples.)

MINOR PROBLEMS

Several compounds failed precision criteria for initial and/or continuing calibrations. Quantitation limits and the reported results for these compounds may be biased and, therefore, have been qualified estimated, "UJ" and "J", respectively. (See Table I in Appendix F for the affected samples).

APPENDIX I (Continued)

2. DATA REVIEW SUMMARY

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NOTES

- o The soil semivolatile MS/MSD analyses were originally extracted within the technical and contractual holding times. Re-extractions were required because of surrogate recoveries, and these re-extractions were performed outside of holding times. Surrogate recoveries were again outside of the QC limits, therefore, original sample results are being reported.
- o The maximum concentration of compounds found in the trip blanks, field blanks, or method blanks are listed below. All samples with concentrations of common laboratory contaminants less than ten times (<10X) the blank concentration, and uncommon laboratory contaminants less than five times (<5X) the blank concentration have been qualified "B" in the data summary table. (See Appendix F).

<u>Compound</u>	<u>Concentration (ug/L)</u>
Methylene chloride *	7 J
Acetone *	9 J
Bis(2-ethylhexyl)phthalate *	10 J

* Common Laboratory Contaminant

- o The semivolatile MS/MSD analyses had compounds other than the spiking compounds present. The following is a table of results and precision estimates for the non-spiked compounds:

<u>MS/MSD Non-Spiked Compounds</u>	<u>Concentration (ug/L)</u>
------------------------------------	-----------------------------

<u>Compound</u>	<u>_____</u>	<u>_____</u>	<u>_____</u>	<u>%RSD</u>
Phenanthrene	150 J	190 J	140 J	16.5
Fluoranthene	340 J	470 J	440 J	16.3
Benzo(a)anthracene	290 J	310 J	320 J	5.0
Chrysene	290 J	330 J	300 J	6.8
Bis (2-ethylhexyl) phthalate	160 J	200 J	240 J	20.0
Benzo (b)pyrene	190 J	240 J	240 J	12.9
Benzo (k) pyrene	230 J	200 J	220 J	7.1
Benzo (a) pyrene	240 J	190 J	240 J	12.9

RSD= Relative Standard Deviation

APPENDIX I (Continued)

2. DATA REVIEW SUMMARY

Page 3 of 3

- o The pesticide/PCB analyses of all soil samples and associated QC samples had surrogate recoveries in excess of the QC limit. Since no positive results were reported for any pesticide or PCB compounds for any of the samples in this case no data was affected. (See Appendix F).
- o The reported Tentatively Identified Compounds (TIC's) in Appendix D have been reviewed and accepted or corrected.
- o All data for Case _____ were reviewed in accordance with the Functional Guidelines for Evaluating Organic Analyses with modifications for use within Region III. The text of this report addresses only those problems affecting usability.

ATTACHMENTS

APPENDIX A - Glossary of Data Qualifiers

APPENDIX B - Data Summary. These include:

- (a) All positive results for target compounds with qualifier codes where applicable.
- (b) All unusable detection limits (qualified "R").

APPENDIX C - Results as Reported by the Laboratory for All Target Compounds

APPENDIX D - Reviewed and Corrected Tentatively Identified Compounds

APPENDIX E - Organic Regional Data Assessment Summary

APPENDIX F - Support Documentation

APPENDIX I (Continued)

3. DATA REVIEW FORM

TABLE
CLP INORGANIC ANALYSIS
CERCLIS SITE NAME _____
CASE No. _____

TEMPA2-9

PAGE ____ of ____

AQUOUS SAMPLE DETECTION LIMITS ($\mu\text{g/l}$)

Sample Location	Inorganic Analytes	DOL ($\mu\text{g/l}$)	Identified in the quality control review (data validation).
Sample Number	Aluminum	P	
Traffic Report Number	Antimony	P	
Remarks	Arsenic	F	
Sampling date	Barium	P	
Analysis Date	Beryllium	P	
	Cadmium	P	
	Calcium	P	
	Chromium	P	
	Cobalt	P	
	Copper	P	
	Iron	P	
	Lead	F	
	Magnesium	P	
	Manganese	P	
	Mercury	CV	
	Nickel	P	
	Potassium	P	
	Selenium	F	
	Silver	P	
	Sodium	P	
	Thorium	F	
	Vanadium	P	
	Zinc	P	
	Cyanide	C	
			UJ The detection limit is approximated due to limitations
			control review (data validation).
			R Value is rejected.
			NA Not Analyzed.
	Analytical Method		
	F Furnace AA	NOTE:	
	P ICP/Flame AA		R
	CV Cold Vapor		NA
	C Colorimetric		
	Sample's wet weight (gms)		
	for ICP analysis		
	for ICP analysis		
	for furnace AA analysis		
	for Cyanide analysis		

TABLE I
CLY INORGANIC ANALYSIS
SOIL AND SEDIMENT SAMPLE DETECTION LIMITS (ppm)

APPENDIX I (Continued)

3. DATA REVIEW FORM

Sample Location	Sample Number	Traffic Report Number	Remarks	Sampling date	Analysis Date	Percent Solids	Inorganic Analytes	IDL (ug/L)	UJ The detection limit is approximated due to limitations identified in the quality control review (data validation).
							Aluminum	P	
							Antimony	P	
							Arsenic	F	
							Barium	P	
							Beryllium	P	
							Cadmium	P	
							Calcium	P	
							Chromium	P	
							Cobalt	P	
							Copper	P	
							Iron	P	
							Lead	F	
							Manganese	P	
							Mercury	CV	
							Nickel	P	
							Potassium	P	
							Selenium	F	
							Silver	P	
							Sodium	P	
							Thallium	F	
							Vanadium	P	
							Zinc	P	
							Cyanide	C	
							Analytical Method	UJ	The detection limit is approximated due to limitations identified in the quality control review (data validation).
							F Furnace AA	R	Value is rejected.
							P ICP/Flame AA	NA	Not Analyzed.
							CV Cold Vapor		
							C Colorimetric		
							Sample's wet weight (gms)		
							for Hg analysis		
							for ICP analysis		
							for furnace AA analysis		
							for cyanide analysis		

APPENDIX I (Continued)

3. DATA REVIEW FORM

TEMP 2+1

TABLE CLP INORGANIC ANALYSIS

CURLIS SITE NAME: _____
CASE NO.: _____SOIL ANALYTICAL RESULTS (w/w)
SDO NO. _____

Sample Location	Sample Number	Traffic Report Number	Remarks	Sampling date	CRDL	Inorganic analytes
					P	Aluminum
					P	Antimony
					F	Arsenic
					P	Barium
					P	Beryllium
					P	Cadmium
					P	Calcium
					P	Chromium
					P	Cobalt
					P	Copper
					P	Iron
					P	Lead
					P	Magnesium
					P	Manganese
					CV	Mercury
					P	Nickel
					P	Potassium
					F	Selenium
					P	Silver
					P	Sodium
					F	Thallium
					P	Vanadium
					P	Zinc
					C	Cyanide

Analytical Method
 F Furnace
 P ICP Flame AA
 CV Cold Vapor
 C Colorimetric

J Quantitation is approximated due to limitations identified during the quality control review.
 R Value is rejected.
 U Revised Sample Quantitation Limit.
 UJ Quantitation limit is approximate due to limitations identified in the quality control review.
 NA Not Analyzed.

Sample results are reported on a dry weight basis.

APPENDIX I (Continued)

3. DATA REVIEW FORM

TABLE I
AQUEOUS ANALYTICAL RESULTS (CONT.)

TEMP12-7

Sample Location	Sample Number	Traffic Report Number	Remarks	Sampling date	CRDL	Inorganic analytes	Quantitation	Method
					P	Aluminum	200	Furnace
					P	Antimony	60	P
					F	Arsenic	10	ICP/Flame AA
					P	Barium	200	CV
					P	Beryllium	5	Colorimetric
					P	Cadmium	5	
					P	Calcium	5000	
					P	Chromium	10	
					P	Cobalt	50	
					P	Copper	25	
					P	Iron	100	
					P	Lead	3	
					P	Magnesium	5000	
					P	Manganese	15	
					CV	Mercury	0.2	
					P	Nickel	40	
					P	Potassium	5000	
					F	Selenium	5	
					P	Silver	10	
					P	Sodium	5000	
					F	Thallium	10	
					P	Vanadium	50	
					P	Zinc	20	
					C	Cyanide	10	
								J Quantitation is approximated due to limitations identified during the quality control review.
								R Value is rejected.
								U Revised Sample Quantitation Limit.
								Q Quantitation limit is approximate due to limitations identified in the quality control review.
								NA Not Analyzed.
								Sample results are reported on a dry weight basis.

TABLE
CLP VOLATILE ORGANIC ANALYSIS AERONICS ANALYTICAL RESULTS (ppm)

CERCUS SITE NAME: _____ SDO NO. _____
CASE NO. _____

Sample Location	Sample Number	Traffic Report Number	Remarks	Sampling Date	Analysis Date	Volatile Organic Compound (CRA)	CRQL
						Chloromethane	10
						Bromomethane	10
						Vinyl Chloride	10
						Chloroethane	10
						Methylene Chloride	5
						Acetone	10
						Carbon Disulfide	5
						1,1-Dichloroethene	5
						1,1-Dichloroethane	5
						1,2-Dichloroethane (Total)	5
						Chloroform	5
						1,2-Dichloroethane	5
						2-Butanone	10
						1,1,1-Trichloroethane	5
						Carbon Tetrachloride	5
						Vinyl Acetate	10
						Bromodichloromethane	5
						1,2-Dichloropropane	5
						cis-1,3-Dichloropropene	5
						Trichloroethene	5
						Dibromochloromethane	5
						1,1,2-Trichloroethane	5
						Benzene	5
						trans-1,3-Dichloropropene	5
						Bromoform	10
						4-Methyl-2-pentanone	10
						2-Hexanone	10
						Tetrachloroethene	5
						1,1,2,2-Tetrachloroethane	5
						Toluene	5
						Chlorobenzene	5
						Ethylbenzene	5
						Styrene	5
						Xylene (Total)	5

APPENDIX I (Continued)

3. DATA REVIEW FORM

CRQL Contract Required Quantitation Limit.
 J Quantitation is approximate due to limitations identified during the quality control review.
 U Quantitation limit is approximated due to limitations identified in the quality control review.
 R Value is rejected.

APPENDIX I (Continued)

3. DATA REVIEW FORM

PAGE _____ of _____

TABLE I CLP VOLATILE ORGANIC ANALYSIS SOIL ANALYTICAL RESULTS (ug/g)

Sample Location	Sample Number	Traffic Report Number	Remarks	Sampling Date	Analysis Date	Volatile Organic Compound	CRQL	Contract Required Quantitation Limit
						Chloromethane	10	J
						Bromoethane	10	J
						Vinyl Chloride	10	J
						Chloroethane	10	J
						Methylene Chloride	5	J
						Acetone	10	J
						Carbon Disulfide	5	J
						1,1-Dichloroethene	5	J
						1,1-Dichloroethane	5	J
						1,2-Dichloroethene (Total)	5	J
						Chloroform	5	J
						1,2-Dichloroethane	5	J
						2-Butalone	10	J
						1,1,1-Trichloroethane	5	J
						Carbon Tetrachloride	5	J
						Vinyl Acetate	10	J
						Bromodichloromethane	5	J
						1,2-Dichloropropane	5	J
						cis-1,3-Dichloropropene	5	J
						Trichloroethene	5	J
						Dibromochloromethane	5	J
						1,1,2-Trichloroethane	5	J
						Benzene	5	J
						trans-1,3-Dichloropropene	5	J
						Bromoform	5	J
						4-Methyl-2-pentanone	10	J
						2-Hexanone	10	J
						Tetrachloroethene	5	J
						1,1,2,2-Tetrachloroethane	5	J
						Toluene	5	J
						Chlorobenzene	5	J
						Ethylbenzene	5	J
						Styrene	5	J
						Xylene (Total)	5	J

CRQL Contract Required Quantitation Limit
 J Quantitation limit is approximate due to limitations identified during the quality control review.
 U Quantitation limit is approximate due to limitations identified in the quality control review.
 R Value is rejected.

APPENDIX I (Continued)

3. DATA REVIEW FORM

TABLE CLP VOLATILE ORGANIC ANALYSIS
SOIL SAMPLE QUANTITATION LIMITS (ppm)

CERCLIS SITE NAME: _____
CASE NO.: _____ SDG NO. _____

Sample Location	Sample Number	Traffic Report Number	Remarks	Sampling Date	Dilution Factor	Percent Solids	Volatile Organic Compound
							Chloromethane
							Bromomethane
							Vinyl Chloride
							Chloroethane
							Methylene Chloride
							Acetone
							Carbon Disulfide
							1,1-Dichloroethene
							1,1-Dichloroethane
							1,2-Dichloroethene (Total)
							Chloroform
							1,2-Dichloroethane
							2-Butene
							1,1,1-Trichloroethane
							Carbon Tetrachloride
							Vinyl Acetate
							Bromodichloromethane
							1,2-Dichloropropane
							cis-1,3-Dichloropropene
							Trichloroethene
							Dibromoacetonemethane
							1,1,2-Trichloroethane
							Benzene
							trans-1,3-Dichloropropene
							Bromoform
							6-Methyl-2-pentanone
							2-Hexanone
							Tetrachloroethene
							1,1,2,2-Tetrachloroethene
							Toluene
							Chlorobenzene
							Ethylbenzene
							Styrene
							Xylene (Total)

Sample Quantitation limits are reported on a dry weight basis.
U Quantitation limit is approximated due to limitations during the quality control review.
R Value is rejected.

CLP VOLATILE ORGANIC ANALYSIS

AQUEOUS SAMPLE QUANTITATION LIMITS (ug/l)

CERC/LIS SITE NAME _____

Sample Location	CASE No.	SDG No.
Sample Number		
Traffic Report Number		
Remarks		
Sampling Date		
Dilution Factor		
Volatile Organic Compound		
Chloromethane		
Bromomethane		
Vinyl Chloride		
Chloroethane		
Methylene Chloride		
Acetone		
Carbon Disulfide		
1,1-Dichloroethene		
1,1-Dichloroethane		
1,2-Dichloroethene (Total)		
Chloroform		
1,2-Dichloroethane		
2-Butanone		
1,1,1-Trichloroethane		
Carbon Tetrachloride		
Vinyl Acetate		
Bromodichloromethane		
1,2-Dichloropropene		
cis-1,3-Dichloropropene		
Trichloroethene		
Dibromochloromethane		
1,1,2-Trichloroethane		
Benzene		
trans-1,3-Dichloropropene		
Bromoform		
4-Methyl-1-Pentanone		
2-Hexanone		
Tetrachloroethene		
1,1,2,2-Tetrachloroethane		
Toluene		
Chlorobenzene		
Ethylbenzene		
Styrene		
Xylene (Total)		

Sample Quantitation limits are reported on a dry weight basis.
 UJ Quantitation limit is approximated due to limitations during the quality control review.
 R Value is rejected.

APPENDIX I (Continued)

3. DATA REVIEW FORM

PAGE ____ OF ____

TABLE _____
CLP EXTRACTABLE ORGANIC ANALYSIS
CERCLIS SITE NAME: _____
CASE NO.: _____ SDG NO. _____
AQUEOUS ANALYTICAL RESULTS (ug/l)

Sample Location	Sample Number	Traffic Report Number	Remarks	Sampling Date	Extraction Date	Analysis Date	Semi-Volatile Compound	CRQL
							Phenol	10
							bis (2-Chloroethyl) ether	10
							2-Chlorophenoil	10
							1,3-Dichlorobenzene	10
							1,4-Dichlorobenzene	10
							Benzyl Alcohol	10
							1,2-Dichlorobenzene	10
							2-Methylphenol	10
							bis (2-Chloro)isopropyl)ether	10
							4-Methylphenol	10
							N-Nitroso-di-n-propyl amine	10
							Hexachlorobutane	10
							Nitrobenzene	10
							Isophorone	10
							2-Nitrophenol	10
							2,4-Dimethylphenol	10
							Benzoic acid	50
							bis (2-Chloroethoxy) methane	10
							2,4-Dichlorophenol	10
							1,2,4-Trichlorobenzene	10
							Phthalene	10
							4-Chloroaniline	10
							Hexachlorobutadiene	10
							4-Chloro-3-methylphenol	10
							2-Methylnaphthalene	10
							Hexachlorocyclopentadiene	10
							2,4,6-Trichlorophenol	10
							2,4,5-Trichlorophenol	50
							1,2-Chloronaphthalene	10
							2-Nitroaniline	50
							Dimethylphthalate	10
							Acenaphthylene	10
							2,6-Dinitrotoluene	10

CRL Contract Required Quantitation Limit.
 J Quantitation is approximate due to limitations identified during the quality control review.
 UJ Quantitation limit is approximated due to limitations identified in the quality control review.
 R Value is rejected.

TABLE CLP EXTRACTABLE ORGANIC ANALYSIS AQUEOUS ANALYTICAL RESULTS (ug)

Sample Location	Sample Number	Traffic Report Number	Remarks	Sampling Date	Extraction Date	Analysis Date	Semi-Volatile Compound	CRL	Required Quantitation Limit
							3-Nitroaniline	50	J
							Acenaphthene	10	J
							2,4-Dinitrophenol	50	J
							4-Nitrophorhol	50	J
							Dibenzofuran	10	J
							2,4-Dinitrotoluene	10	J
							Diethylphthalate	10	J
							4-Chlorophenyl-phenylether	10	J
							Fluorene	10	J
							4-Nitroaniline	50	J
							4,6-Dinitro-2-methylphenol	50	J
							N-Nitrosodiphenylamine	10	J
							4-Bromophenyl phenylether	10	J
							Hexachlorobenzene	10	J
							Pentachlorophenol	50	J
							Phenanthrene	10	J
							Anthracene	10	J
							Di-n-butyl phthalate	10	J
							Fluoranthene	10	J
							Pyrene	10	J
							Butylbenzylphthalate	10	J
							3,3'-Dichlorobenzidine	20	J
							Benz(a)anthracene	10	J
							Chrysene	10	J
							Bis(2-Ethylhexyl)phthalate	10	J
							Di-n-octyl phthalate	10	J
							Benz(b)fluoranthene	10	J
							Benz(k)fluoranthene	10	J
							Benzo(a)pyrene	10	J
							Indeno(1,2,3-cd)pyrene	10	J
							Dibenz(a,h)anthracene	10	J
							Benzo(g,h,i)perylene	10	J

J Quantitation is approximate due to limitations identified during the quality control review.
 UJ Quantitation limit is approximated due to limitations identified in the quality control review.
 R Value is rejected.

APPENDIX I (Continued)

3. DATA REVIEW FORM

TABLE
CLP EXTRACTABLE ORGANIC ANALYSIS
AQUEOUS ANALYTICAL RESULTS (ug/l)

Sample Location	Sample Number	Traffic Report Number	Remarks	Sampling Date	Extraction Date	Analysis Date	Pesticide/PCB Compound	CRAT
							alpha-BHC	0.05
							beta-BHC	0.05
							delta-BHC	0.05
							gamma-BHC (l-indane)	0.05
							Heptachlor	0.05
							Aldrin	0.05
							Heptachlor epoxide	0.05
							Endosulfan I	0.05
							Dieldrin	0.10
							4,4'-DDE	0.10
							Endrin	0.10
							Endosulfan II	0.10
							4,4'-DDD	0.10
							Endosulfan sulfate	0.10
							4,4'-DDT	0.10
							Methoxychlor	0.5
							Endrin ketone	0.10
							alpha-Chlordane	0.5
							gamma-Chlordane	0.5
							Toxaphene	1.0
							Aroclor-1016	0.5
							Aroclor-1221	0.5
							Aroclor-1232	0.5
							Aroclor-1242	0.5
							Aroclor-1248	0.5
							Aroclor-1254	1.0
							Aroclor-1260	1.0

CRAT: Contract Required Quantitation Limit.
 J: Quantitation is approximate due to limitations identified during the quality control review.
 U: Quantitation limit is approximated due to limitations identified in the quality control review.
 R: Value is rejected.

TABLE I
CIP EXTRACTABLE ORGANIC ANALYSIS
AQUEOUS SAMPLE QUANTITATION LIMITS (ug/L)

APPENDIX I (Continued)

3. DATA REVIEW FORM

Sample Location	Sample Number	Traffic Report Number	Sampling Date	Dilution Factor	Percent Solids	Semi-Volatile Compound
						3-Nitroaniline
						Aceanaphthene
						2,4,6-trinitrophenol
						4-Nitrostrophenol
						Dibenzofuran
						2,4-Dinitrotoluene
						Diethylphthalate
						4-Chlorophenyl-phenylether
						Fluorene
						4-Nitroaniline
						4,6-Dinitro-2-methylphenol
						N-Nitrosodiphenylamine
						4-Bromophenyl-phenylether
						Hexachlorobenzene
						Pentachlorophenol
						Phenanthrene
						Anthracene
						Di-n-butylphthalate
						Fluoranthene
						Pyrene
						Burylbenzylphthalate
						3,3'-Dichlorobenzidine
						Benzocycloanthracene
						Chrysene
						bis(2-Ethylhexyl)phthalate
						Di-n-octyl phthalate
						Benzoc(b)fluoranthene
						Benzoc(b)pyrene
						Indeno(1,2,3-cd)pyrene
						Dibenz(a,h)anthracene
						Benzoc(9,h,i)perylene

Sample Quantitation limits are reported on dry weight basis. Quantitation limits are approximate due to limitations identified during the quality control review.

APPENDIX I (Continued)

3. DATA REVIEW FORM

PAGE _____ of _____

TABLE CDP EXTRACTABLE ORGANIC ANALYSIS AQUEOUS SAMPLE QUANTITATION LIMITS (ppb)

Sample Location	Sample Number	Traffic Report Number	Remarks	Sampling Date	Dilution Factor	Percent Solids	Semi-Volatile Compound
							Phenol
							bis (2-Chloroethyl) ether
							2-Chlorophenol
							1,3-Dichlorobenzene
							1,4-Dichlorobenzene
							Benzyl Alcohol
							1,2-Dichlorobenzene
							2-Ethylphenol
							bis (2-Chloroisopropyl)ether
							4-Ethylphenol
							N-Nitroso-di-n-propylamine
							Hexachloroethane
							Nitrobenzene
							Isophorone
							2-Nitrophenol
							2,4-Dimethylphenol
							Benzoic acid
							bis (2-Chloroethoxy) methane
							2,4-Dichlorophenol
							1,2,4-Trichlorobenzene
							Naphthalene
							6-Chloroniline
							Hexachlorobutadiene
							4-Chloro-3-methylphenol
							2-Methylnaphthalene
							Hexachlorocyclopentadiene
							2,4,6-Trichlorophenol
							2,4,5-Trichlorophenol
							2-Chloronaphthalene
							2-Nitroaniline
							Dimethylphthalate
							Acenaphthylene
							2,6-Dinitrotoluene

Quantitation limits are reported on a dry weight basis.
 U Quantitation limit is approximated due to limitations identified in the quality control review.
 R Value is rejected.

TABLE I
CLP EXTRACTABLE ORGANIC ANALYSIS AQUEOUS SAMPLE QUANTITATION LIMITS (ug/l)

CERCLIS SITE NAME: _____
CASE NO.: _____ SDG NO.: _____

Sample Location	Sample Number	Traffic Report Number	Remarks	Sampling Date	Dilution Factor	Percent Solids	Pesticide/PCB Compound
							alpha-BHC
							Delta-BHC
							delta-BHC
							gamma-BHC (Lindane)
							Heptachlor
							Aldrin
							Heptachlor epoxide
							Endosulfan I
							Endosulfan II
							Endosulfan sulfate
							4,4'-DDT
							Methoxychlor
							Endrin ketone
							alpha-Chlordane
							gamma-Chlordane
							Toxaphene
							Aroclor-1016
							Aroclor-1221
							Aroclor-1232
							Aroclor-1242
							Aroclor-1248
							Aroclor-1254
							Aroclor-1260

Sample Quantitation Limits are reported on dry weight basis.
 UW Quantitation Limits are approximate due to limitations identified during the quality control review.
 R Value is rejected.

APPENDIX I (Continued)

3. DATA REVIEW FORM

APPENDIX I (Continued)

3. DATA REVIEW FORM

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TABLE CIP EXTRACTABLE ORGANIC ANALYSIS

TMRP2-5-1

CIRCLE SITE NAME: _____
CASE NO.: _____

SOIL ANALYTICAL RESULTS (ug/g)

Sample Location	Semi-Volatile Compound	CQL	Contract Required Quantitation Limit
Sample Number	Phenol	330	J
Traffic Report Number	bis (2-Chloroethyl) ether	330	Quantitation is approximate due to limitations identified during the quality control review.
Remarks	2-Chlorophenol	330	
Sampling Date	1,3-Dichlorobenzene	330	
Extraction Date	1,4-Dichlorobenzene	330	
Analysis Date	Benzyl Alcohol	330	
	1,2-Dichlorobenzene	330	
	2-Methylphenol	330	
	bis (2-Chloroisopropyl) ether	330	
	4-Methylphenol	330	
	N-Nitroso-di-n-propylamine	330	
	Hexachloroethane	330	
	Nitrobenzene	330	
	Isophorone	330	
	2-Nitrophenol	330	
	2,4-Dimethylphenol	330	
	Benzoic Acid	1600	J
	bis (2-Chloroethoxy) methane	330	Quantitation is approximate due to limitations identified during the quality control review.
	2,4-Dichlorophenol	330	
	1,2,4-Trichlorobenzene	330	
	Naphthalene	330	
	4-Chloroniline	330	
	Hexachlorobutadiene	330	
	4-Chloro-3-methylphenol	330	
	2-Methylnaphthalene	330	
	Heptachlorocyclopentadiene	330	
	2,4,6-Trichlorophenol	330	
	2,4,5-Trichlorophenol	1600	R
	2-Chloronaphthalene	330	Value is rejected.
	2-Nitroaniline	1600	
	Dinethyl phthalate	330	
	Acenaphthylene	330	
	2,6-Dinitrotoluene	330	

TABLE CLP EXTRACTABLE ORGANIC ANALYSIS
CIRCLE/S SITE NAME: _____
CASE NO.: _____ SDNO.: _____

Sample Location	Traffic Report Number	Sampling Date	Extraction Date	Semi-volatile compound	CRDL	SOIL ANALYTICAL RESULTS (ug/g)
				3-Nitroaniline	1600	
				Acenaphthene	330	
				2,4-Dinitrophenol	1600	
				4-Nitrophenol	1600	
				Oibenzenoturan	330	
				2,4-Dinitrotoluene	330	
				Diethyl phthalate	330	
				4-Chlorophenyl-phenyl ether	330	
				Fluorene	330	
				4-Nitroaniline	1600	
				4,6-Dinitro-2-methylphenol	1600	
				N-Nitrosodiphenylamine	330	
				4-Aromophenyl-phenyl ether	330	
				Pentachlorobenzene	330	
				Pentachlorophenol	1600	
				Phenanthrene	330	
				Anthracene	330	
				Di-n-butylphthalate	330	
				Fluoranthene	330	
				Pyrene	330	
				Benzylbenzylphthalate	330	
				3,3'-Dichlorobenzidine	660	
				Benz(a)anthracene	330	
				Chrysene	330	
				bis(2-Ethyhexyl)phthalate	330	
				Di-n-octyl phthalate	330	
				Benz(b)fluoranthene	330	
				Benz(a)fluoranthene	330	
				Benz(a)pyrene	330	
				Indeno (1,2,3-cd)pyrene	330	
				Dibenz(a,h)anthracene	330	
				Benzot(g,h,i)perylene	330	

CRDL Contract Required Detection Limit.
 J Quantitation is approximate due to limitations identified during the quality control review.
 UJ Quantitation limit is approximate due to limitations identified in the quality control review.
 R Value is rejected.

APPENDIX I (Continued)

3. DATA REVIEW FORM

TABLE
CLP EXTRACTABLE ORGANIC ANALYSIS
CERCLIS SITE NAME: _____
CASE NO.: _____

TMR2-S-3

SOIL ANALYTICAL RESULTS (ug/g)

PAGE ____ of ____

Sample Location	Sample Number	Traffic Report Number	Remarks	Sampling Date	Extraction Date	Pesticide/PCB Compound	CRL
						alpha-BHC	8.0
						beta-BHC	8.0
						delta-BHC	8.0
						gamma-BHC (lindane)	8.0
						Heptachlor	8.0
						Aldrin	8.0
						Heptachlor epoxide	8.0
						Endosulfan I	8.0
						Dieldrin	16.0
						6,4'-DD	16.0
						Endrin	16.0
						Endosulfan II	16.0
						4,4'-DDD	16.0
						Endosulfan sulfate	16.0
						6,4'-DDT	16.0
						Methoxychlor	80.0
						Endrin ketone	16.0
						alpha-Chlordane	80.0
						gamma-Chlordane	80.0
						Toxaphene	160.0
						Aroclor-1016	80.0
						Aroclor-1221	80.0
						Aroclor-1232	80.0
						Aroclor-1242	80.0
						Aroclor-1248	160.0
						Aroclor-1254	160.0
						Aroclor-1260	160.0

CRL = Contract Required Quantitation Limit.
 J = Quantitation is approximate due to limitations identified during the quality control review.
 U = Quantitation is approximate due to limitations identified in the quality control review.
 R = Value is rejected.

3. DATA REVIEW FORM

Sample Location	Sample Number	Traffic Report Number	Remarks	Sampling Date	Dilution factor	Percent Solids	Semi-Volatile Compound
							Phenol
							bis (2-Chloroethyl) ether
							2-Chlorophenol
							1,3-Dichlorobenzene
							1,4-Dichlorobenzene
							Benzyl Alcohol
							1,2-Dichlorobenzene
							2-Methylphenol
							bis (2-Chloroisopropyl)ether
							4-Nitrophenol
							N,N-Nitroso-di-n-propylamine
							Hexachloroethane
							Nitrobenzene
							Isophorone
							2-Nitrophenol
							2,4-Dimethylphenol
							Benzoic acid
							bis (2-Chloroethoxy) methane
							1,2,4-Trichlorobenzene
							Naphthalene
							4-Chloroaniline
							Hexachlorobutadiene
							4-Chloro-3-methylphenol
							2-Methylnaphthalene
							Hexachlorocyclohexadiene
							2,4,6-Trichlorophenol
							2,4,5-Trichlorophenol
							2-Chlorophthalalene
							2-Nitroaniline
							Dimethylphthalate
							Acenaphthylene
							2,6-Dinitrotoluene

Sample Quantitation Limits are reported on a dry weight basis.
 UJ Quantitation limit is approximated due to limitations identified in the quality control review.
 R Value is rejected.

Sample Location	Sample Number	Traffic Report Number	Sampling Date	Dilution Factor	Percent Solids	Semi-Volatile Compound	SD No.	CASE NO.	CHLUS SITE NAME	TABLE
						3-Nitroaniline				
						Acenaphthene				
						2,4-Dinitrophenol				
						4-Nitrophenol				
						Dibenzofuran				
						2,4-Dinitrotoluene				
						Diethylphthalate				
						4-Chlorophenyl-phenylether				
						Fluorene				
						4-Nitroaniline				
						4,6-Dinitro-2-methyl phenol				
						N-Nitrosodiphenylamine				
						4-Bromophenyl-phenylether				
						Hexachlorobenzene				
						Pentachlorophenol				
						Phenanthrene				
						Anthracene				
						Di-n-butylphthalate				
						Fluoranthene				
						Pyrene				
						Butylbenzylphthalate				
						3,5'-Dichlorobenzidine				
						Benz(a)anthracene				
						Chrysene				
						bis(2-Ethylhexyl)phthalate				
						Di-n-octyl phthalate				
						Benz(b)fluoranthene				
						Benz(k)fluoranthene				
						Benz(a)pyrene				
						Indeno (1,2,3-cd)pyrene				
						Dibenz(a,h)anthracene				
						Benz(g,h,i)perylene				

Sample Quantitation Limits are reported on dry weight basis.
 UJ Quantitation Limits are approximate due to limitations identified during the quality control review.
 R Value is rejected.

APPENDIX I (Continued)

3. DATA REVIEW FORM

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TABLE _____ CCLP EXTRACTABLE ORGANIC ANALYSIS
CIRCLE IS STR. NAME(S) SOIL SAMPLE QUANTITATION LIMITS (ug/g)
CASE NO. _____ SDNO. _____

Sample Location	Sample Number	Traffic Report Number	Remarks	Sampling Date	Dilution Factor	Percent Solids	Pesticide/PCB Compound
							alpha-BHC
							beta-BHC
							delta-BHC
							gamma-BHC (Lindane)
							Heptachlor
							Aldrin
							Heptachlor epoxide
							Endosulfan I
							Heptdrin
							4,4'-DDT
							Endrin
							Endosulfan II
							4,4'-DDD
							Endosulfan sulfate
							4,4'-DDT
							Methoxychlor
							Endrin ketone
							alpha-Chlordane
							gamma-Chlordane
							Toxaphene
							Aroclor-1016
							Aroclor-1221
							Aroclor-1232
							Aroclor-1242
							Aroclor-1248
							Aroclor-1254
							Aroclor-1260

Sample Quantitation Limits are reported on dry weight basis.
 UJ Quantitation limits are approximate due to limitations identified during the quality control review.