

Report of Analysis

Client:	Environmental Restoration, LLC		Date Collected:	06/27/25	
Project:	CC2-16 Analytical		Date Received:	06/27/25	
Client Sample ID:	CC0625-OXBL		SDG No.:	Q2481	
Lab Sample ID:	Q2481-14		Matrix:	TCLP	
Analytical Method:	8270E		% Solid:	0	
Sample Wt/Vol:	10	Units: mL	Final Vol:	1000	uL
Soil Aliquot Vol:		uL	Test:	TCLP BNA	
Extraction Type :		Decanted : N	Level :	LOW	
Injection Volume :		GPC Factor : 1.0	GPC Cleanup :	N	PH :
Prep Method :	SW3541				

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
BP025241.D	10	07/23/25 11:15	07/24/25 14:05	PB168983

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
110-86-1	Pyridine	1300	U	1300	5000	ug/L
106-46-7	1,4-Dichlorobenzene	530	U	530	5000	ug/L
95-48-7	2-Methylphenol	1100	U	1100	5000	ug/L
65794-96-9	3+4-Methylphenols	1100	U	1100	10000	ug/L
67-72-1	Hexachloroethane	650	U	650	5000	ug/L
98-95-3	Nitrobenzene	760	U	760	5000	ug/L
87-68-3	Hexachlorobutadiene	540	U	540	5000	ug/L
88-06-2	2,4,6-Trichlorophenol	510	U	510	5000	ug/L
95-95-4	2,4,5-Trichlorophenol	620	U	620	5000	ug/L
121-14-2	2,4-Dinitrotoluene	1200	U	1200	5000	ug/L
118-74-1	Hexachlorobenzene	520	U	520	5000	ug/L
87-86-5	Pentachlorophenol	1600	U	1600	10000	ug/L
SURROGATES						
367-12-4	2-Fluorophenol	117	*	23 - 138	7%	SPK: 150
13127-88-3	Phenol-d6	68.0	*	10 - 134	4%	SPK: 150
4165-60-0	Nitrobenzene-d5	86.3	*	67 - 132	8%	SPK: 100
321-60-8	2-Fluorobiphenyl	602		52 - 132	60%	SPK: 100
118-79-6	2,4,6-Tribromophenol	701		44 - 137	46%	SPK: 150
1718-51-0	Terphenyl-d14	112	*	42 - 152	11%	SPK: 100
INTERNAL STANDARDS						
3855-82-1	1,4-Dichlorobenzene-d4	488000		7.431		
1146-65-2	Naphthalene-d8	2050000		10.178		
15067-26-2	Acenaphthene-d10	203000		14.078		
1517-22-2	Phenanthrene-d10	2660000		16.907		
1719-03-5	Chrysene-d12	2520000		21.324		
1520-96-3	Perylene-d12	760000		24.477		

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U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products