

Report of Analysis

Client:	Environmental Restoration, LLC		Date Collected:	06/27/25	
Project:	CC2-16 Analytical		Date Received:	06/27/25	
Client Sample ID:	CC0627-OXL		SDG No.:	Q2481	
Lab Sample ID:	Q2481-18		Matrix:	TCLP	
Analytical Method:	8270E		% Solid:	0	
Sample Wt/Vol:	10	Units: mL	Final Vol:	2000	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
BP025246.D	10	07/23/25 11:15	07/24/25 17:31	PB168983

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
110-86-1	Pyridine	2600	U	2600	10000	ug/L
106-46-7	1,4-Dichlorobenzene	1100	U	1100	10000	ug/L
95-48-7	2-Methylphenol	2200	U	2200	10000	ug/L
65794-96-9	3+4-Methylphenols	2200	U	2200	20000	ug/L
67-72-1	Hexachloroethane	1300	U	1300	10000	ug/L
98-95-3	Nitrobenzene	1500	U	1500	10000	ug/L
87-68-3	Hexachlorobutadiene	1100	U	1100	10000	ug/L
88-06-2	2,4,6-Trichlorophenol	1000	U	1000	10000	ug/L
95-95-4	2,4,5-Trichlorophenol	1200	U	1200	10000	ug/L
121-14-2	2,4-Dinitrotoluene	2400	U	2400	10000	ug/L
118-74-1	Hexachlorobenzene	1000	U	1000	10000	ug/L
87-86-5	Pentachlorophenol	3200	U	3200	20000	ug/L
SURROGATES						
367-12-4	2-Fluorophenol	0.14	*	23 - 138	0%	SPK: 150
13127-88-3	Phenol-d6	0	*	10 - 134	0%	SPK: 150
4165-60-0	Nitrobenzene-d5	489	*	67 - 132	48%	SPK: 100
321-60-8	2-Fluorobiphenyl	100000	*	52 - 132	10034%	SPK: 100
118-79-6	2,4,6-Tribromophenol	122000	*	44 - 137	8108%	SPK: 150
1718-51-0	Terphenyl-d14	415		42 - 152	41%	SPK: 100
INTERNAL STANDARDS						
3855-82-1	1,4-Dichlorobenzene-d4	491000		7.437		
1146-65-2	Naphthalene-d8	179000		10.178		
15067-26-2	Acenaphthene-d10	511		14.084		
1517-22-2	Phenanthrene-d10	64500		16.884		
1719-03-5	Chrysene-d12	56200		21.336		
1520-96-3	Perylene-d12	1450		24.436		

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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