

Report of Analysis

Client:	ENTACT		Date Collected:	05/19/25	
Project:	540 Degraw St, Brooklyn, NY - E9309		Date Received:	05/19/25	
Client Sample ID:	OR-640-COMP-66MS		SDG No.:	Q2078	
Lab Sample ID:	Q2084-05MS		Matrix:	TCLP	
Analytical Method:	8270E		% Solid:	0	
Sample Wt/Vol:	100	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 :	SW3510C				

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
BP024781.D	1	05/22/25 11:14	05/23/25 06:37	PB168131

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
110-86-1	Pyridine	0.31		0.013	0.050	mg/L
106-46-7	1,4-Dichlorobenzene	0.39		0.0053	0.050	mg/L
95-48-7	2-Methylphenol	0.40		0.011	0.050	mg/L
65794-96-9	3+4-Methylphenols	0.41		0.011	0.10	mg/L
67-72-1	Hexachloroethane	0.36		0.0065	0.050	mg/L
98-95-3	Nitrobenzene	0.39		0.0076	0.050	mg/L
87-68-3	Hexachlorobutadiene	0.45		0.0054	0.050	mg/L
88-06-2	2,4,6-Trichlorophenol	0.50		0.0051	0.050	mg/L
95-95-4	2,4,5-Trichlorophenol	0.50		0.0062	0.050	mg/L
121-14-2	2,4-Dinitrotoluene	0.50		0.012	0.050	mg/L
118-74-1	Hexachlorobenzene	0.51		0.0052	0.050	mg/L
87-86-5	Pentachlorophenol	1.10	E	0.016	0.10	mg/L
SURROGATES						
367-12-4	2-Fluorophenol	122		15 (10) - 110 (139)	82%	SPK: 150
13127-88-3	Phenol-d6	107		15 (10) - 110 (134)	71%	SPK: 150
4165-60-0	Nitrobenzene-d5	71.8		30 (49) - 130 (133)	72%	SPK: 100
321-60-8	2-Fluorobiphenyl	78.3		30 (52) - 130 (132)	78%	SPK: 100
118-79-6	2,4,6-Tribromophenol	163		15 (44) - 110 (137)	109%	SPK: 150
1718-51-0	Terphenyl-d14	82.9		30 (48) - 130 (125)	83%	SPK: 100
INTERNAL STANDARDS						
3855-82-1	1,4-Dichlorobenzene-d4	108000	7.651			
1146-65-2	Naphthalene-d8	434000	10.422			
15067-26-2	Acenaphthene-d10	286000	14.28			
1517-22-2	Phenanthrene-d10	578000	17.08			
1719-03-5	Chrysene-d12	676000	21.521			
1520-96-3	Perylene-d12	797000	24.798			

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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