

Report of Analysis

Client:	PSEG	Date Collected:	05/30/24
Project:	Victory Laydown Yard	Date Received:	05/30/24
Client Sample ID:	WC-TP-1MS	SDG No.:	P2687
Lab Sample ID:	P2668-04MS	Matrix:	TCLP
Analytical Method:	SW8270	% 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
BP020517.D	1	06/03/24 10:50	06/05/24 14:24	PB161311

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
110-86-1	Pyridine	0.31		0.016	0.050	mg/L
106-46-7	1,4-Dichlorobenzene	0.40		0.0084	0.050	mg/L
95-48-7	2-Methylphenol	0.41		0.011	0.050	mg/L
65794-96-9	3+4-Methylphenols	0.41		0.012	0.10	mg/L
67-72-1	Hexachloroethane	0.39		0.010	0.050	mg/L
98-95-3	Nitrobenzene	0.41		0.013	0.050	mg/L
87-68-3	Hexachlorobutadiene	0.40		0.013	0.050	mg/L
88-06-2	2,4,6-Trichlorophenol	0.49		0.0089	0.050	mg/L
95-95-4	2,4,5-Trichlorophenol	0.46		0.010	0.050	mg/L
121-14-2	2,4-Dinitrotoluene	0.46		0.015	0.050	mg/L
118-74-1	Hexachlorobenzene	0.44		0.011	0.050	mg/L
87-86-5	Pentachlorophenol	0.77		0.019	0.10	mg/L
SURROGATES						
367-12-4	2-Fluorophenol	127		10 - 139	85%	SPK: 150
13127-88-3	Phenol-d6	112		10 - 134	75%	SPK: 150
4165-60-0	Nitrobenzene-d5	84.0		49 - 133	84%	SPK: 100
321-60-8	2-Fluorobiphenyl	79.4		52 - 132	79%	SPK: 100
118-79-6	2,4,6-Tribromophenol	147		32 - 145	98%	SPK: 150
1718-51-0	Terphenyl-d14	77.7		36 - 145	78%	SPK: 100
INTERNAL STANDARDS						
3855-82-1	1,4-Dichlorobenzene-d4	290000	7.763			
1146-65-2	Naphthalene-d8	1130000	10.534			
15067-26-2	Acenaphthene-d10	728000	14.392			
1517-22-2	Phenanthrene-d10	1490000	17.186			
1719-03-5	Chrysene-d12	1360000	21.651			
1520-96-3	Perylene-d12	1580000	25.004			

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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