

Report of Analysis

Client:	Tully Construction Co., Inc.		Date Collected:	
Project:	DOT Harper Street Yard - North Side		Date Received:	
Client Sample ID:	PB167310BS		SDG No.:	Q1627
Lab Sample ID:	PB167310BS		Matrix:	TCLP
Analytical Method:	SW8270		% Solid:	0
Sample Wt/Vol:	1000	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
BF142096.D	1	03/25/25 12:25	03/26/25 14:36	PB167310

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
110-86-1	Pyridine	46.9		1.30	5.00	ug/L
106-46-7	1,4-Dichlorobenzene	50.6		0.53	5.00	ug/L
95-48-7	2-Methylphenol	49.5		1.10	5.00	ug/L
65794-96-9	3+4-Methylphenols	48.0		1.10	10.0	ug/L
67-72-1	Hexachloroethane	50.1		0.65	5.00	ug/L
98-95-3	Nitrobenzene	50.0		0.76	5.00	ug/L
87-68-3	Hexachlorobutadiene	53.3		0.54	5.00	ug/L
88-06-2	2,4,6-Trichlorophenol	50.3		0.51	5.00	ug/L
95-95-4	2,4,5-Trichlorophenol	53.2		0.62	5.00	ug/L
121-14-2	2,4-Dinitrotoluene	57.8		1.20	5.00	ug/L
118-74-1	Hexachlorobenzene	53.3		0.52	5.00	ug/L
87-86-5	Pentachlorophenol	100	E	1.60	10.0	ug/L
SURROGATES						
367-12-4	2-Fluorophenol	146		10 - 139	97%	SPK: 150
13127-88-3	Phenol-d6	140		10 - 134	93%	SPK: 150
4165-60-0	Nitrobenzene-d5	104		49 - 133	104%	SPK: 100
321-60-8	2-Fluorobiphenyl	103		52 - 132	103%	SPK: 100
118-79-6	2,4,6-Tribromophenol	179		44 - 137	120%	SPK: 150
1718-51-0	Terphenyl-d14	107		48 - 125	107%	SPK: 100
INTERNAL STANDARDS						
3855-82-1	1,4-Dichlorobenzene-d4	126000		6.869		
1146-65-2	Naphthalene-d8	477000		8.151		
15067-26-2	Acenaphthene-d10	273000		9.91		
1517-22-2	Phenanthrene-d10	497000		11.398		
1719-03-5	Chrysene-d12	338000		14.039		
1520-96-3	Perylene-d12	324000		15.51		

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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