

Report of Analysis

Client:	Aramark Uniforms		Date Collected:	
Project:	Monthly 2025		Date Received:	
Client Sample ID:	PB166115BL		SDG No.:	Q1119
Lab Sample ID:	PB166115BL		Matrix:	TCLP
Analytical Method:	SW8270		% Solid:	0
Sample Wt/Vol:	1000	Units: mL	Final Vol:	1000 uL
Soil Aliquot Vol:		uL	Test:	TCLP BNA Group1
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
BF141213.D	1	01/17/25 12:00	01/20/25 11:03	PB166115

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
110-86-1	Pyridine	0.0016	U	0.0016	0.0050	mg/L
106-46-7	1,4-Dichlorobenzene	0.00084	U	0.00084	0.0050	mg/L
95-48-7	2-Methylphenol	0.0011	U	0.0011	0.0050	mg/L
65794-96-9	3+4-Methylphenols	0.0012	U	0.0012	0.010	mg/L
67-72-1	Hexachloroethane	0.0010	U	0.0010	0.0050	mg/L
98-95-3	Nitrobenzene	0.0013	U	0.0013	0.0050	mg/L
88-06-2	2,4,6-Trichlorophenol	0.00089	U	0.00089	0.0050	mg/L
95-95-4	2,4,5-Trichlorophenol	0.0010	U	0.0010	0.0050	mg/L
121-14-2	2,4-Dinitrotoluene	0.0015	U	0.0015	0.0050	mg/L
118-74-1	Hexachlorobenzene	0.0011	U	0.0011	0.0050	mg/L
87-86-5	Pentachlorophenol	0.0019	U	0.0019	0.010	mg/L
SURROGATES						
367-12-4	2-Fluorophenol	134		10 - 139	90%	SPK: 150
13127-88-3	Phenol-d6	132		10 - 134	88%	SPK: 150
4165-60-0	Nitrobenzene-d5	94.1		49 - 133	94%	SPK: 100
321-60-8	2-Fluorobiphenyl	94.3		52 - 132	94%	SPK: 100
118-79-6	2,4,6-Tribromophenol	153		44 - 137	102%	SPK: 150
1718-51-0	Terphenyl-d14	87.9		48 - 125	88%	SPK: 100
INTERNAL STANDARDS						
3855-82-1	1,4-Dichlorobenzene-d4	163000	6.81			
1146-65-2	Naphthalene-d8	650000	8.092			
15067-26-2	Acenaphthene-d10	351000	9.845			
1517-22-2	Phenanthrene-d10	609000	11.327			
1719-03-5	Chrysene-d12	443000	13.968			
1520-96-3	Perylene-d12	341000	15.433			

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CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
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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