Alliance TECHNICAL GROUP
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Report of Analysis

Client:	Aramark Uniforms			Date Collected:	
Project:	Monthly 2025			Date Received:	
Client Sample ID:	PB170216BL			SDG No.:	Q3368
Lab Sample ID:	PB170216BL			Matrix:	TCLP
Analytical Method:	8270E	Level :	LOW	% Solid:	0
Sample Wt/Vol:	1000 mL	Final Vol:	1000 uL	Test:	TCLP BNA Group1
Prep Method :		Prep Date:	10/22/25		

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	Prep BatchID
TARGETS									
110-86-1	Pyridine	1.30	U	1	1.30	5.00	ug/L	10/22/25 17:14	PB170216
106-46-7	1,4-Dichlorobenzene	0.53	U	1	0.53	5.00	ug/L	10/22/25 17:14	PB170216
95-48-7	2-Methylphenol	1.10	U	1	1.10	5.00	ug/L	10/22/25 17:14	PB170216
65794-96-9	3+4-Methylphenols	1.10	U	1	1.10	10.0	ug/L	10/22/25 17:14	PB170216
67-72-1	Hexachloroethane	0.65	U	1	0.65	5.00	ug/L	10/22/25 17:14	PB170216
98-95-3	Nitrobenzene	0.76	U	1	0.76	5.00	ug/L	10/22/25 17:14	PB170216
88-06-2	2,4,6-Trichlorophenol	0.51	U	1	0.51	5.00	ug/L	10/22/25 17:14	PB170216
95-95-4	2,4,5-Trichlorophenol	0.62	U	1	0.62	5.00	ug/L	10/22/25 17:14	PB170216
121-14-2	2,4-Dinitrotoluene	1.20	U	1	1.20	5.00	ug/L	10/22/25 17:14	PB170216
118-74-1	Hexachlorobenzene	0.52	U	1	0.52	5.00	ug/L	10/22/25 17:14	PB170216
87-86-5	Pentachlorophenol	1.60	U	1	1.60	10.0	ug/L	10/22/25 17:14	PB170216
SURROGATES									
367-12-4	2-Fluorophenol	151			23 - 138	101%	SPK: 150		
13127-88-3	Phenol-d6	142			10 - 134	95%	SPK: 150		
4165-60-0	Nitrobenzene-d5	78.7			67 - 132	79%	SPK: 100		
321-60-8	2-Fluorobiphenyl	76.6			52 - 132	77%	SPK: 100		
118-79-6	2,4,6-Tribromophenol	131			44 - 137	87%	SPK: 150		
1718-51-0	Terphenyl-d14	80.4			42 - 152	80%	SPK: 100		
INTERNAL STANDARDS		**Area Count**							
3855-82-1	1,4-Dichlorobenzene-d4	156000							
1146-65-2	Naphthalene-d8	660000							
15067-26-2	Acenaphthene-d10	489000							
1517-22-2	Phenanthrene-d10	1080000							
1719-03-5	Chrysene-d12	1040000							
1520-96-3	Perylene-d12	886000							

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