Alliance TECHNICAL GROUP
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Report of Analysis

Client:	Aramark Uniforms			Date Collected:	10/22/25
Project:	Monthly 2025			Date Received:	10/22/25
Client Sample ID:	PB170172TB			SDG No.:	Q3368
Lab Sample ID:	PB170172TB			Matrix:	TCLP
Analytical Method:	8270E	Level :	LOW	% Solid:	0
Sample Wt/Vol:	100 mL	Final Vol:	1000 uL	Test:	TCLP BNA Group1
Prep Method :	SW3541	Prep Date:	10/22/25		

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	Prep BatchID
TARGETS									
110-86-1	Pyridine	12.8	U	1	12.8	50.0	ug/L	10/22/25 18:36	PB170216
106-46-7	1,4-Dichlorobenzene	5.30	U	1	5.30	50.0	ug/L	10/22/25 18:36	PB170216
95-48-7	2-Methylphenol	11.2	U	1	11.2	50.0	ug/L	10/22/25 18:36	PB170216
65794-96-9	3+4-Methylphenols	11.0	U	1	11.0	100	ug/L	10/22/25 18:36	PB170216
67-72-1	Hexachloroethane	6.50	U	1	6.50	50.0	ug/L	10/22/25 18:36	PB170216
98-95-3	Nitrobenzene	7.60	U	1	7.60	50.0	ug/L	10/22/25 18:36	PB170216
88-06-2	2,4,6-Trichlorophenol	5.10	U	1	5.10	50.0	ug/L	10/22/25 18:36	PB170216
95-95-4	2,4,5-Trichlorophenol	6.20	U	1	6.20	50.0	ug/L	10/22/25 18:36	PB170216
121-14-2	2,4-Dinitrotoluene	12.2	U	1	12.2	50.0	ug/L	10/22/25 18:36	PB170216
118-74-1	Hexachlorobenzene	5.20	U	1	5.20	50.0	ug/L	10/22/25 18:36	PB170216
87-86-5	Pentachlorophenol	15.8	U	1	15.8	100	ug/L	10/22/25 18:36	PB170216
SURROGATES									
367-12-4	2-Fluorophenol	146			23 - 138	98%	SPK: 150		
13127-88-3	Phenol-d6	136			10 - 134	91%	SPK: 150		
4165-60-0	Nitrobenzene-d5	80.4			67 - 132	80%	SPK: 100		
321-60-8	2-Fluorobiphenyl	77.1			52 - 132	77%	SPK: 100		
118-79-6	2,4,6-Tribromophenol	129			44 - 137	86%	SPK: 150		
1718-51-0	Terphenyl-d14	89.0			42 - 152	89%	SPK: 100		
INTERNAL STANDARDS		**Area Count**							
3855-82-1	1,4-Dichlorobenzene-d4	158000							
1146-65-2	Naphthalene-d8	655000							
15067-26-2	Acenaphthene-d10	478000							
1517-22-2	Phenanthrene-d10	1020000							
1719-03-5	Chrysene-d12	810000							
1520-96-3	Perylene-d12	659000							

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