

Report of Analysis

Client:	Aramark Uniforms		Date Collected:	10/16/25
Project:	Monthly 2025		Date Received:	10/16/25
Client Sample ID:	OUTSIDE-DUMPSTER		SDG No.:	Q3368
Lab Sample ID:	Q3368-06		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 19:17	PB170216
106-46-7	1,4-Dichlorobenzene	5.30	U	1	5.30	50.0	ug/L	10/22/25 19:17	PB170216
95-48-7	2-Methylphenol	11.2	U	1	11.2	50.0	ug/L	10/22/25 19:17	PB170216
65794-96-9	3+4-Methylphenols	11.0	U	1	11.0	100	ug/L	10/22/25 19:17	PB170216
67-72-1	Hexachloroethane	6.50	U	1	6.50	50.0	ug/L	10/22/25 19:17	PB170216
98-95-3	Nitrobenzene	7.60	U	1	7.60	50.0	ug/L	10/22/25 19:17	PB170216
88-06-2	2,4,6-Trichlorophenol	5.10	U	1	5.10	50.0	ug/L	10/22/25 19:17	PB170216
95-95-4	2,4,5-Trichlorophenol	6.20	U	1	6.20	50.0	ug/L	10/22/25 19:17	PB170216
121-14-2	2,4-Dinitrotoluene	12.2	U	1	12.2	50.0	ug/L	10/22/25 19:17	PB170216
118-74-1	Hexachlorobenzene	5.20	U	1	5.20	50.0	ug/L	10/22/25 19:17	PB170216
87-86-5	Pentachlorophenol	15.8	U	1	15.8	100	ug/L	10/22/25 19:17	PB170216
SURROGATES									
367-12-4	2-Fluorophenol	130			23 - 138	87%	SPK: 150		
13127-88-3	Phenol-d6	122			10 - 134	81%	SPK: 150		
4165-60-0	Nitrobenzene-d5	74.6			67 - 132	75%	SPK: 100		
321-60-8	2-Fluorobiphenyl	68.9			52 - 132	69%	SPK: 100		
118-79-6	2,4,6-Tribromophenol	129			44 - 137	86%	SPK: 150		
1718-51-0	Terphenyl-d14	55.3			42 - 152	55%	SPK: 100		
INTERNAL STANDARDS									
		Area Count							
3855-82-1	1,4-Dichlorobenzene-d4	151000							
1146-65-2	Naphthalene-d8	649000							
15067-26-2	Acenaphthene-d10	478000							
1517-22-2	Phenanthrene-d10	1050000							
1719-03-5	Chrysene-d12	1090000							
1520-96-3	Perylene-d12	970000							

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