

Report of Analysis

Client:	Aramark Uniforms		Date Collected:	08/06/25	
Project:	Monthly 2025		Date Received:	08/06/25	
Client Sample ID:	TP-4MSD		SDG No.:	Q2791	
Lab Sample ID:	Q2779-04MSD		Matrix:	TCLP	
Analytical Method:	8270E		% Solid:	0	
Sample Wt/Vol:	100	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 :					

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
BP025348.D	1	08/08/25 11:12	08/11/25 14:41	PB169180

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
110-86-1	Pyridine	320		12.8	50.0	ug/L
106-46-7	1,4-Dichlorobenzene	380		5.30	50.0	ug/L
95-48-7	2-Methylphenol	360		11.2	50.0	ug/L
65794-96-9	3+4-Methylphenols	350		11.0	100	ug/L
67-72-1	Hexachloroethane	390		6.50	50.0	ug/L
98-95-3	Nitrobenzene	420		7.60	50.0	ug/L
88-06-2	2,4,6-Trichlorophenol	480		5.10	50.0	ug/L
95-95-4	2,4,5-Trichlorophenol	460		6.20	50.0	ug/L
121-14-2	2,4-Dinitrotoluene	430		12.2	50.0	ug/L
118-74-1	Hexachlorobenzene	470		5.20	50.0	ug/L
87-86-5	Pentachlorophenol	980	E	15.8	100	ug/L
SURROGATES						
367-12-4	2-Fluorophenol	107		23 - 138	71%	SPK: 150
13127-88-3	Phenol-d6	91.7		10 - 134	61%	SPK: 150
4165-60-0	Nitrobenzene-d5	80.3		67 - 132	80%	SPK: 100
321-60-8	2-Fluorobiphenyl	75.8		52 - 132	76%	SPK: 100
118-79-6	2,4,6-Tribromophenol	135		44 - 137	90%	SPK: 150
1718-51-0	Terphenyl-d14	80.5		42 - 152	81%	SPK: 100
INTERNAL STANDARDS						
3855-82-1	1,4-Dichlorobenzene-d4	285000		7.796		
1146-65-2	Naphthalene-d8	1090000		10.572		
15067-26-2	Acenaphthene-d10	635000		14.419		
1517-22-2	Phenanthrene-d10	1140000		17.219		
1719-03-5	Chrysene-d12	1070000		21.66		
1520-96-3	Perylene-d12	1200000		25.048		

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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