

Report of Analysis

Client:	Weston Solutions	Date Collected:	03/10/25
Project:	Ft Meade Tipton Airfield Parcel RI - PO 0111169	Date Received:	03/11/25
Client Sample ID:	TAPFTA-MW01I-031025-00-T2	SDG No.:	Q1539
Lab Sample ID:	Q1539-02	Matrix:	Water
Analytical Method:	SW8270	% Solid:	0
Sample Wt/Vol:	930 Units: mL	Final Vol:	1000 uL
Soil Aliquot Vol:	uL	Test:	SVOC-TCL BNA -20
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
BF141944.D	1	03/12/25 08:45	03/13/25 16:44	PB167097

CAS Number	Parameter	Conc.	Qualifier	MDL	LOD	LOQ / CRQL	Units
TARGETS							
100-52-7	Benzaldehyde	8.60	U	4.20	8.60	10.8	ug/L
108-95-2	Phenol	4.30	U	0.98	4.30	5.40	ug/L
111-44-4	bis(2-Chloroethyl)ether	4.30	U	0.87	4.30	5.40	ug/L
95-57-8	2-Chlorophenol	4.30	U	0.62	4.30	5.40	ug/L
95-48-7	2-Methylphenol	4.30	U	1.20	4.30	5.40	ug/L
108-60-1	2,2-oxybis(1-Chloropropane)	4.30	U	1.40	4.30	5.40	ug/L
98-86-2	Acetophenone	4.30	U	0.80	4.30	5.40	ug/L
65794-96-9	3+4-Methylphenols	8.60	U	1.20	8.60	10.8	ug/L
621-64-7	n-Nitroso-di-n-propylamine	2.70	U	1.50	2.70	2.70	ug/L
67-72-1	Hexachloroethane	4.30	U	0.70	4.30	5.40	ug/L
98-95-3	Nitrobenzene	4.30	U	0.82	4.30	5.40	ug/L
78-59-1	Isophorone	4.30	U	0.81	4.30	5.40	ug/L
88-75-5	2-Nitrophenol	4.30	U	1.90	4.30	5.40	ug/L
105-67-9	2,4-Dimethylphenol	4.30	U	2.00	4.30	5.40	ug/L
111-91-1	bis(2-Chloroethoxy)methane	4.30	U	0.73	4.30	5.40	ug/L
120-83-2	2,4-Dichlorophenol	4.30	U	0.56	4.30	5.40	ug/L
91-20-3	Naphthalene	4.30	U	0.54	4.30	5.40	ug/L
106-47-8	4-Chloroaniline	4.30	UQ	0.90	4.30	5.40	ug/L
87-68-3	Hexachlorobutadiene	4.30	U	0.58	4.30	5.40	ug/L
105-60-2	Caprolactam	8.60	U	1.20	8.60	10.8	ug/L
59-50-7	4-Chloro-3-methylphenol	4.30	U	0.63	4.30	5.40	ug/L
91-57-6	2-Methylnaphthalene	4.30	U	0.60	4.30	5.40	ug/L
77-47-4	Hexachlorocyclopentadiene	8.60	UQ	3.90	8.60	10.8	ug/L
88-06-2	2,4,6-Trichlorophenol	4.30	U	0.55	4.30	5.40	ug/L
95-95-4	2,4,5-Trichlorophenol	4.30	U	0.67	4.30	5.40	ug/L
92-52-4	1,1-Biphenyl	4.30	U	0.57	4.30	5.40	ug/L
91-58-7	2-Chloronaphthalene	4.30	U	0.66	4.30	5.40	ug/L
88-74-4	2-Nitroaniline	4.30	U	1.40	4.30	5.40	ug/L
131-11-3	Dimethylphthalate	4.30	U	0.66	4.30	5.40	ug/L

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208-96-8	Acenaphthylene	4.30	U	0.81	4.30	5.40	ug/L
606-20-2	2,6-Dinitrotoluene	4.30	U	0.99	4.30	5.40	ug/L
99-09-2	3-Nitroaniline	4.30	U	1.10	4.30	5.40	ug/L
83-32-9	Acenaphthene	4.30	U	0.59	4.30	5.40	ug/L
51-28-5	2,4-Dinitrophenol	8.60	U	6.40	8.60	10.8	ug/L
100-02-7	4-Nitrophenol	8.60	U	2.60	8.60	10.8	ug/L
132-64-9	Dibenzofuran	4.30	U	0.66	4.30	5.40	ug/L
121-14-2	2,4-Dinitrotoluene	4.30	U	1.30	4.30	5.40	ug/L
84-66-2	Diethylphthalate	4.30	U	0.74	4.30	5.40	ug/L
7005-72-3	4-Chlorophenyl-phenylether	4.30	U	0.73	4.30	5.40	ug/L
86-73-7	Fluorene	4.30	U	0.68	4.30	5.40	ug/L
100-01-6	4-Nitroaniline	4.30	U	1.60	4.30	5.40	ug/L
534-52-1	4,6-Dinitro-2-methylphenol	8.60	U	3.10	8.60	10.8	ug/L
86-30-6	n-Nitrosodiphenylamine	4.30	U	0.62	4.30	5.40	ug/L
101-55-3	4-Bromophenyl-phenylether	4.30	U	0.43	4.30	5.40	ug/L
118-74-1	Hexachlorobenzene	4.30	U	0.56	4.30	5.40	ug/L
1912-24-9	Atrazine	4.30	U	1.10	4.30	5.40	ug/L
87-86-5	Pentachlorophenol	8.60	U	1.70	8.60	10.8	ug/L
85-01-8	Phenanthrene	4.30	U	0.54	4.30	5.40	ug/L
120-12-7	Anthracene	4.30	U	0.66	4.30	5.40	ug/L
86-74-8	Carbazole	4.30	U	0.77	4.30	5.40	ug/L
84-74-2	Di-n-butylphthalate	4.30	U	1.30	4.30	5.40	ug/L
206-44-0	Fluoranthene	4.30	U	0.88	4.30	5.40	ug/L
129-00-0	Pyrene	4.30	U	0.54	4.30	5.40	ug/L
85-68-7	Butylbenzylphthalate	4.30	U	2.10	4.30	5.40	ug/L
91-94-1	3,3-Dichlorobenzidine	8.60	U	1.00	8.60	10.8	ug/L
56-55-3	Benzo(a)anthracene	4.30	U	0.48	4.30	5.40	ug/L
218-01-9	Chrysene	4.30	U	0.47	4.30	5.40	ug/L
117-81-7	Bis(2-ethylhexyl)phthalate	4.30	U	1.70	4.30	5.40	ug/L
117-84-0	Di-n-octyl phthalate	8.60	U	2.50	8.60	10.8	ug/L
205-99-2	Benzo(b)fluoranthene	4.30	U	0.53	4.30	5.40	ug/L

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207-08-9	Benzo(k)fluoranthene	4.30	U	0.52	4.30	5.40	ug/L
50-32-8	Benzo(a)pyrene	4.30	U	0.59	4.30	5.40	ug/L
193-39-5	Indeno(1,2,3-cd)pyrene	4.30	U	0.63	4.30	5.40	ug/L
53-70-3	Dibenzo(a,h)anthracene	4.30	U	0.72	4.30	5.40	ug/L
191-24-2	Benzo(g,h,i)perylene	4.30	U	0.74	4.30	5.40	ug/L
95-94-3	1,2,4,5-Tetrachlorobenzene	4.30	U	0.56	4.30	5.40	ug/L
123-91-1	1,4-Dioxane	4.30	U	1.10	4.30	5.40	ug/L
58-90-2	2,3,4,6-Tetrachlorophenol	4.30	U	0.77	4.30	5.40	ug/L
SURROGATES							
367-12-4	2-Fluorophenol	73.5		19 - 119		49%	SPK: 150
13127-88-3	Phenol-d6	45.4		10 - 130		30%	SPK: 150
4165-60-0	Nitrobenzene-d5	104		44 - 120		104%	SPK: 100
321-60-8	2-Fluorobiphenyl	96.7		44 - 119		97%	SPK: 100
118-79-6	2,4,6-Tribromophenol	176		43 - 140		117%	SPK: 150
1718-51-0	Terphenyl-d14	86.4		50 - 134		86%	SPK: 100
INTERNAL STANDARDS							
3855-82-1	1,4-Dichlorobenzene-d4	169000	6.875				
1146-65-2	Naphthalene-d8	663000	8.157				
15067-26-2	Acenaphthene-d10	386000	9.91				
1517-22-2	Phenanthrene-d10	719000	11.392				
1719-03-5	Chrysene-d12	525000	14.027				
1520-96-3	Perylene-d12	378000	15.498				
TENTATIVE IDENTIFIED COMPOUNDS							
000994-05-8	Butane, 2-methoxy-2-methyl-	130	JB			2.21	ug/L
000111-76-2	Ethanol, 2-butoxy-	34.8	J			5.83	ug/L
002516-93-0	Butoxyacetic acid	3.00	J			7.36	ug/L

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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