

Report of Analysis

Client:	Alliance Technical Group, LLC - Newark		Date Collected:	09/02/25
Project:	NJ Waste Water PT		Date Received:	09/04/25
Client Sample ID:	WP0925-RR-PAH-WPDL		SDG No.:	Q3015
Lab Sample ID:	Q3015-21DL		Matrix:	Water
Analytical Method:	SW8270ESIM		% Solid:	0
Sample Wt/Vol:	1000	Units: mL	Final Vol:	1000 uL
Soil Aliquot Vol:		uL	Test:	SVOCMS Group5
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
BN037876.D	10	09/16/25 09:39	09/24/25 15:05	PB169712

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
91-20-3	Naphthalene	2.70	D	0.40	1.00	ug/L
91-57-6	2-Methylnaphthalene	4.80	D	0.42	1.00	ug/L
208-96-8	Acenaphthylene	12.3	D	0.38	1.00	ug/L
83-32-9	Acenaphthene	3.00	D	0.38	1.00	ug/L
86-73-7	Fluorene	3.90	D	0.38	1.00	ug/L
85-01-8	Phenanthrene	1.60	D	0.34	1.00	ug/L
120-12-7	Anthracene	2.00	D	0.36	1.00	ug/L
206-44-0	Fluoranthene	0.99	JD	0.34	1.00	ug/L
129-00-0	Pyrene	4.40	D	0.31	1.00	ug/L
56-55-3	Benzo(a)anthracene	2.70	D	0.36	1.00	ug/L
218-01-9	Chrysene	4.60	D	0.37	1.00	ug/L
205-99-2	Benzo(b)fluoranthene	2.10	D	0.35	1.00	ug/L
207-08-9	Benzo(k)fluoranthene	2.10	D	0.50	1.00	ug/L
50-32-8	Benzo(a)pyrene	1.70	D	0.42	1.00	ug/L
193-39-5	Indeno(1,2,3-cd)pyrene	1.60	D	0.46	1.00	ug/L
53-70-3	Dibenzo(a,h)anthracene	3.60	D	0.47	1.00	ug/L
191-24-2	Benzo(g,h,i)perylene	3.50	D	0.42	1.00	ug/L
SURROGATES						
7297-45-2	2-Methylnaphthalene-d10	0.39		20 - 139	97%	SPK: 0.4
93951-69-0	Fluoranthene-d10	0.40		54 - 157	100%	SPK: 0.4
4165-60-0	Nitrobenzene-d5	0.99	*	27 - 154	247%	SPK: 0.4
321-60-8	2-Fluorobiphenyl	1.01	*	30 - 155	252%	SPK: 0.4
1718-51-0	Terphenyl-d14	1.08	*	54 - 175	270%	SPK: 0.4
INTERNAL STANDARDS						
3855-82-1	1,4-Dichlorobenzene-d4	5420		7.833		
1146-65-2	Naphthalene-d8	15300		10.637		
15067-26-2	Acenaphthene-d10	7910		14.484		
1517-22-2	Phenanthrene-d10	15600		17.223		
1719-03-5	Chrysene-d12	13300		21.411		

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1520-96-3	Perylene-d12	12100	23.741			

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