

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100325\
 Data File : BN037993.D
 Acq On : 03 Oct 2025 11:03
 Operator : RC/JU
 Sample : Q2899-02
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MDL-WATER-QT3-2025-06

Quant Time: Oct 03 11:31:48 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN100125.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 02 09:41:10 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.813	152	9016	0.400	ng/ul	0.00
4) Naphthalene-d8	10.612	136	25053	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.465	164	13131	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.213	188	25004	0.400	ng/ul	0.00
17) Chrysene-d12	21.394	240	16012	0.400	ng/ul	0.00
23) Perylene-d12	23.715	264	13614	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.256	96	5616	0.587	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.207	152	11016	0.311	ng/ul	0.00
18) Fluoranthene-d10	19.244	212	18742	0.387	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.294	88	778	0.078	ng/ul#	76
5) Naphthalene	10.661	128	5126	0.075	ng/ul	97
7) 2-Methylnaphthalene	12.284	142	3340	0.073	ng/ul	98
8) 1-Methylnaphthalene	12.504	142	3110	0.065	ng/ul	100
10) Acenaphthylene	14.183	152	4552	0.073	ng/ul	98
11) Acenaphthene	14.530	153	3472	0.076	ng/ul	97
12) Fluorene	15.520	166	3733	0.071	ng/ul	99
14) Pentachlorophenol	16.867	266	433	0.067	ng/ul	95
15) Phenanthrene	17.256	178	6062	0.075	ng/ul	97
16) Anthracene	17.348	178	5033	0.071	ng/ul	97
19) Fluoranthene	19.272	202	6132	0.091	ng/ul	98
20) Pyrene	19.635	202	6023	0.091	ng/ul	98
21) Benzo(a)anthracene	21.380	228	4486	0.077	ng/ul	98
22) Chrysene	21.429	228	4896	0.080	ng/ul	98
24) Benzo(b)fluoranthene	23.014	252	4813	0.081	ng/ul#	47
25) Benzo(k)fluoranthene	23.057	252	4758	0.080	ng/ul#	64
26) Benzo(a)pyrene	23.613	252	4161	0.085	ng/ul#	15
27) Indeno(1,2,3-cd)pyrene	26.081	276	5406	0.092	ng/ul#	77
28) Dibenzo(a,h)anthracene	26.101	278	4041	0.088	ng/ul#	77
29) Benzo(g,h,i)perylene	26.802	276	4136	0.085	ng/ul#	88

(#) = qualifier out of range (m) = manual integration (+) = signals summed

