

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110924\
 Data File : BN034944.D
 Acq On : 09 Nov 2024 16:26
 Operator : RC/JU
 Sample : PB164777BS
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS777

Quant Time: Nov 09 22:52:54 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN110924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Nov 09 00:09:12 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.571	152	2903	0.400	ng/ul	0.00
4) Naphthalene-d8	10.333	136	7356	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.203	164	5647	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.950	188	14140	0.400	ng/ul	0.00
17) Chrysene-d12	21.142	240	13818	0.400	ng/ul	0.00
23) Perylene-d12	23.308	264	12998	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.153	96	1690	0.612	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.928	152	4386	0.389	ng/ul	0.00
18) Fluoranthene-d10	18.981	212	14230	0.401	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.182	88	2290	0.756	ng/ul	94
5) Naphthalene	10.383	128	3188	0.173	ng/ul	100
7) 2-Methylnaphthalene	12.005	142	2326	0.169	ng/ul	99
8) 1-Methylnaphthalene	12.225	142	2389	0.173	ng/ul	99
10) Acenaphthylene	13.921	152	3804	0.165	ng/ul	99
11) Acenaphthene	14.268	153	2956	0.169	ng/ul	98
12) Fluorene	15.258	166	3571	0.161	ng/ul	97
14) Pentachlorophenol	16.603	266	1067	0.326	ng/ul	95
15) Phenanthrene	16.992	178	6343	0.168	ng/ul	99
16) Anthracene	17.080	178	5731	0.161	ng/ul	100
19) Fluoranthene	19.014	202	8225	0.178	ng/ul#	96
20) Pyrene	19.376	202	8518	0.178	ng/ul	97
21) Benzo(a)anthracene	21.128	228	8501	0.170	ng/ul	100
22) Chrysene	21.177	228	8876	0.174	ng/ul	99
24) Benzo(b)fluoranthene	22.665	252	8246	0.180	ng/ul	93
25) Benzo(k)fluoranthene	22.706	252	8509	0.178	ng/ul#	92
26) Benzo(a)pyrene	23.214	252	7007	0.172	ng/ul#	89
27) Indeno(1,2,3-cd)pyrene	25.467	276	8111	0.162	ng/ul#	95
28) Dibenzo(a,h)anthracene	25.480	278	6281	0.161	ng/ul	94
29) Benzo(g,h,i)perylene	26.121	276	6191	0.162	ng/ul#	94

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

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