

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082824\
 Data File : BN033663.D
 Acq On : 29 Aug 2024 02:12
 Operator : MA/JU
 Sample : PB162970BS
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS970

Quant Time: Aug 29 02:41:18 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN082724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 27 23:48:09 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.536	152	5239	0.400	ng/ul	0.00
4) Naphthalene-d8	10.294	136	14225	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.174	164	7080	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.923	188	14906	0.400	ng/ul	0.00
17) Chrysene-d12	21.130	240	12049	0.400	ng/ul	0.00
23) Perylene-d12	23.292	264	12600	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.143	96	4619	0.858	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.894	152	8589	0.421	ng/ul	0.00
18) Fluoranthene-d10	18.962	212	16640	0.463	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.177	88	12116	2.027	ng/ul#	80
5) Naphthalene	10.343	128	14528	0.382	ng/ul	100
7) 2-Methylnaphthalene	11.965	142	9224	0.383	ng/ul	99
8) 1-Methylnaphthalene	12.191	142	10721	0.432	ng/ul	99
10) Acenaphthylene	13.887	152	11766	0.366	ng/ul	99
11) Acenaphthene	14.234	153	9487	0.409	ng/ul	98
12) Fluorene	15.229	166	10247	0.395	ng/ul	99
14) Pentachlorophenol	16.581	266	3021	0.644	ng/ul	99
15) Phenanthrene	16.965	178	16518	0.376	ng/ul	100
16) Anthracene	17.054	178	14610	0.370	ng/ul	100
19) Fluoranthene	18.989	202	18623	0.392	ng/ul	98
20) Pyrene	19.357	202	19417	0.385	ng/ul	98
21) Benzo(a)anthracene	21.115	228	17806	0.375	ng/ul	99
22) Chrysene	21.165	228	18654	0.389	ng/ul	100
24) Benzo(b)fluoranthene	22.649	252	18624	0.381	ng/ul	99
25) Benzo(k)fluoranthene	22.690	252	19047	0.389	ng/ul	99
26) Benzo(a)pyrene	23.199	252	16587	0.422	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.438	276	22284	0.378	ng/ul	99
28) Dibenzo(a,h)anthracene	25.451	278	17329	0.378	ng/ul	98
29) Benzo(g,h,i)perylene	26.085	276	17593	0.382	ng/ul	99

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

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