

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110924\
 Data File : BN034974.D
 Acq On : 10 Nov 2024 11:33
 Operator : RC/JU
 Sample : PB164816BS
 Misc :
 ALS Vial : 67 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS816

Quant Time: Nov 10 21:51:48 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	7.566	152	3158	0.400	ng/ul	0.00
4) Naphthalene-d8	10.328	136	7625	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.203	164	5514	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.949	188	12666	0.400	ng/ul	0.00
17) Chrysene-d12	21.142	240	11677	0.400	ng/ul	0.00
23) Perylene-d12	23.305	264	12225	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.149	96	2177	0.724	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.928	152	5103	0.437	ng/ul	0.00
18) Fluoranthene-d10	18.981	212	13739	0.458	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.182	88	5866	1.781	ng/ul#	91
5) Naphthalene	10.377	128	7489	0.392	ng/ul	98
7) 2-Methylnaphthalene	12.000	142	5398	0.379	ng/ul	100
8) 1-Methylnaphthalene	12.225	142	6257	0.438	ng/ul	99
10) Acenaphthylene	13.916	152	8421	0.374	ng/ul	100
11) Acenaphthene	14.263	153	6403	0.375	ng/ul	99
12) Fluorene	15.258	166	7694	0.354	ng/ul	100
14) Pentachlorophenol	16.603	266	2194	0.749	ng/ul	97
15) Phenanthrene	16.992	178	12979	0.383	ng/ul	99
16) Anthracene	17.080	178	11882	0.374	ng/ul	99
19) Fluoranthene	19.009	202	15781	0.404	ng/ul#	96
20) Pyrene	19.372	202	16218	0.401	ng/ul#	96
21) Benzo(a)anthracene	21.125	228	16005	0.378	ng/ul	99
22) Chrysene	21.177	228	16378	0.380	ng/ul	99
24) Benzo(b)fluoranthene	22.659	252	16054	0.373	ng/ul	99
25) Benzo(k)fluoranthene	22.706	252	17218	0.383	ng/ul	99
26) Benzo(a)pyrene	23.211	252	14708	0.383	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.456	276	19411	0.413	ng/ul#	93
28) Dibenzo(a,h)anthracene	25.473	278	15364	0.420	ng/ul	98
29) Benzo(g,h,i)perylene	26.110	276	14995	0.418	ng/ul	98

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

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