

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110924\
 Data File : BN034993.D
 Acq On : 10 Nov 2024 23:28
 Operator : RC/JU
 Sample : PB164818BS
 Misc :
 ALS Vial : 87 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS818

Quant Time: Nov 10 23:51:07 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	2816	0.400	ng/ul	0.00
4) Naphthalene-d8	10.328	136	6697	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.198	164	4820	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.945	188	11369	0.400	ng/ul	0.00
17) Chrysene-d12	21.139	240	11029	0.400	ng/ul	0.00
23) Perylene-d12	23.302	264	11674	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.149	96	1970	0.735	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.923	152	4377	0.427	ng/ul	-0.01
18) Fluoranthene-d10	18.981	212	12631	0.446	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.182	88	5380	1.832	ng/ul#	93
5) Naphthalene	10.377	128	6467	0.386	ng/ul	98
7) 2-Methylnaphthalene	12.000	142	4669	0.373	ng/ul	100
8) 1-Methylnaphthalene	12.220	142	5364	0.427	ng/ul	99
10) Acenaphthylene	13.916	152	7201	0.366	ng/ul	100
11) Acenaphthene	14.263	153	5523	0.370	ng/ul	99
12) Fluorene	15.253	166	6734	0.355	ng/ul	99
14) Pentachlorophenol	16.599	266	2003	0.762	ng/ul	98
15) Phenanthrene	16.987	178	11504	0.379	ng/ul	99
16) Anthracene	17.076	178	10664	0.374	ng/ul	98
19) Fluoranthene	19.009	202	14461	0.392	ng/ul#	97
20) Pyrene	19.372	202	15008	0.393	ng/ul	97
21) Benzo(a)anthracene	21.125	228	15109	0.378	ng/ul	99
22) Chrysene	21.174	228	15535	0.381	ng/ul	99
24) Benzo(b)fluoranthene	22.659	252	15499	0.378	ng/ul	99
25) Benzo(k)fluoranthene	22.703	252	16460	0.383	ng/ul	99
26) Benzo(a)pyrene	23.208	252	14014	0.382	ng/ul#	98
27) Indeno(1,2,3-cd)pyrene	25.453	276	19072	0.425	ng/ul#	94
28) Dibenzo(a,h)anthracene	25.470	278	15034	0.430	ng/ul	97
29) Benzo(g,h,i)perylene	26.107	276	14887	0.434	ng/ul	98

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

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