

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071524\
 Data File : BM046649.D
 Acq On : 17 Jul 2024 01:01
 Operator : MA/JU
 Sample : PB161884BS
 Misc :
 ALS Vial : 61 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS884

Quant Time: Jul 17 01:37:43 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM071024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 15 08:57:27 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.502	152	2275	0.400	ng/ul	0.00
4) Naphthalene-d8	10.259	136	6227	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.145	164	4284	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.895	188	9590	0.400	ng/ul	0.00
17) Chrysene-d12	21.102	240	7626	0.400	ng/ul	0.00
23) Perylene-d12	23.239	264	8823	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.113	96	1036	0.554	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.859	152	3840	0.382	ng/ul	0.00
18) Fluoranthene-d10	18.936	212	9391	0.405	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.143	88	2826	1.417	ng/ul#	74
5) Naphthalene	10.308	128	5901	0.373	ng/ul	99
7) 2-Methylnaphthalene	11.931	142	4149	0.373	ng/ul	99
8) 1-Methylnaphthalene	12.156	142	4200	0.370	ng/ul	99
10) Acenaphthylene	13.858	152	6852	0.351	ng/ul	98
11) Acenaphthene	14.205	153	4828	0.370	ng/ul	97
12) Fluorene	15.199	166	5741	0.355	ng/ul	96
14) Pentachlorophenol	16.553	266	2684	0.552	ng/ul	99
15) Phenanthrene	16.937	178	9467	0.346	ng/ul	100
16) Anthracene	17.026	178	9298	0.348	ng/ul	100
19) Fluoranthene	18.964	202	11828	0.378	ng/ul	98
20) Pyrene	19.326	202	12482	0.368	ng/ul#	97
21) Benzo(a)anthracene	21.085	228	11733	0.348	ng/ul	99
22) Chrysene	21.138	228	11192	0.343	ng/ul	100
24) Benzo(b)fluoranthene	22.604	252	12114	0.344	ng/ul	95
25) Benzo(k)fluoranthene	22.645	252	12228	0.353	ng/ul#	95
26) Benzo(a)pyrene	23.145	252	11169	0.385	ng/ul#	95
27) Indeno(1,2,3-cd)pyrene	25.357	276	14713	0.333	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.374	278	11508	0.329	ng/ul#	96
29) Benzo(g,h,i)perylene	26.001	276	11510	0.327	ng/ul	96

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

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