

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072924\
 Data File : BM046913.D
 Acq On : 30 Jul 2024 09:48
 Operator : MA/JU
 Sample : PB162176BS
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS176

Quant Time: Jul 30 10:57:09 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM072424.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 30 04:18:24 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.451	152	2706	0.400	ng/ul	0.00
4) Naphthalene-d8	10.211	136	7538	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.100	164	4437	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.858	188	8191	0.400	ng/ul	0.00
17) Chrysene-d12	21.076	240	5554	0.400	ng/ul	0.00
23) Perylene-d12	23.200	264	6355	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.088	96	1151	0.607	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.811	152	4534	0.378	ng/ul	0.00
18) Fluoranthene-d10	18.900	212	7135	0.394	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.117	88	3348	1.544	ng/ul#	86
5) Naphthalene	10.255	128	7095	0.372	ng/ul	98
7) 2-Methylnaphthalene	11.882	142	4965	0.378	ng/ul	99
8) 1-Methylnaphthalene	12.108	142	4982	0.368	ng/ul	100
10) Acenaphthylene	13.813	152	7513	0.370	ng/ul	100
11) Acenaphthene	14.164	153	5398	0.391	ng/ul	99
12) Fluorene	15.159	166	5961	0.346	ng/ul	100
14) Pentachlorophenol	16.521	266	2116	0.571	ng/ul	98
15) Phenanthrene	16.901	178	8396	0.357	ng/ul	98
16) Anthracene	16.989	178	7887	0.360	ng/ul	100
19) Fluoranthene	18.932	202	9072	0.372	ng/ul	98
20) Pyrene	19.295	202	9291	0.368	ng/ul	99
21) Benzo(a)anthracene	21.058	228	8075	0.349	ng/ul	99
22) Chrysene	21.111	228	7996	0.340	ng/ul	99
24) Benzo(b)fluoranthene	22.572	252	8624	0.309	ng/ul	96
25) Benzo(k)fluoranthene	22.612	252	8781	0.312	ng/ul	99
26) Benzo(a)pyrene	23.107	252	7992	0.399	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	25.299	276	11112	0.360	ng/ul#	97
28) Dibenzo(a,h)anthracene	25.313	278	8563	0.379	ng/ul	99
29) Benzo(g,h,i)perylene	25.936	276	8766	0.323	ng/ul	98

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

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