

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072524\
 Data File : BM046802.D
 Acq On : 25 Jul 2024 16:33
 Operator : MA/JU
 Sample : PB162099BS
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS099

Quant Time: Jul 25 17:07:05 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 : Wed Jul 24 15:10:37 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.472	152	3153	0.400	ng/ul	0.00
4) Naphthalene-d8	10.227	136	9023	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.118	164	5854	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.871	188	12525	0.400	ng/ul	0.00
17) Chrysene-d12	21.081	240	8131	0.400	ng/ul	0.00
23) Perylene-d12	23.209	264	7115	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.101	96	1324	0.600	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.827	152	4768	0.332	ng/ul	0.00
18) Fluoranthene-d10	18.909	212	10594	0.400	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.134	88	3461	1.370	ng/ul#	85
5) Naphthalene	10.277	128	7218	0.316	ng/ul	100
7) 2-Methylnaphthalene	11.904	142	5178	0.330	ng/ul	99
8) 1-Methylnaphthalene	12.130	142	5260	0.324	ng/ul	99
10) Acenaphthylene	13.831	152	8581	0.320	ng/ul	98
11) Acenaphthene	14.178	153	6160	0.338	ng/ul	98
12) Fluorene	15.173	166	7125	0.313	ng/ul	99
14) Pentachlorophenol	16.529	266	3017	0.532	ng/ul	99
15) Phenanthrene	16.913	178	11289	0.314	ng/ul	99
16) Anthracene	17.002	178	10664	0.318	ng/ul	99
19) Fluoranthene	18.942	202	13330	0.373	ng/ul	100
20) Pyrene	19.304	202	13859	0.375	ng/ul	99
21) Benzo(a)anthracene	21.064	228	11098	0.328	ng/ul	99
22) Chrysene	21.116	228	10614	0.308	ng/ul	99
24) Benzo(b)fluoranthene	22.577	252	9345	0.299	ng/ul	92
25) Benzo(k)fluoranthene	22.618	252	9345	0.297	ng/ul	96
26) Benzo(a)pyrene	23.112	252	8325	0.371	ng/ul#	95
27) Indeno(1,2,3-cd)pyrene	25.309	276	9763	0.283	ng/ul	95
28) Dibenzo(a,h)anthracene	25.326	278	7426	0.294	ng/ul#	93
29) Benzo(g,h,i)perylene	25.946	276	7850	0.258	ng/ul#	95

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

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