

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071224\
 Data File : BM046553.D
 Acq On : 12 Jul 2024 20:05
 Operator : MA/JU
 Sample : P3059-04MS
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 E1GM5MS

Quant Time: Jul 13 00:35:48 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 : Fri Jul 12 05:09:19 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.510	152	1927	0.400	ng/ul	0.00
4) Naphthalene-d8	10.270	136	4721	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.154	164	3078	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.903	188	6701	0.400	ng/ul	0.00
17) Chrysene-d12	21.111	240	5926	0.400	ng/ul #	0.00
23) Perylene-d12	23.247	264	8077	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.122	96	469	0.296	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.870	152	2651	0.348	ng/ul	0.00
18) Fluoranthene-d10	18.940	212	7142	0.396	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.151	88	1287	0.762	ng/ul#	74
5) Naphthalene	10.319	128	4061	0.338	ng/ul	99
7) 2-Methylnaphthalene	11.942	142	2936	0.348	ng/ul	100
8) 1-Methylnaphthalene	12.167	142	2912	0.338	ng/ul	99
10) Acenaphthylene	13.867	152	4680	0.334	ng/ul	99
11) Acenaphthene	14.214	153	3414	0.364	ng/ul	97
12) Fluorene	15.209	166	5002	0.431	ng/ul	99
14) Pentachlorophenol	16.557	266	2077	0.612	ng/ul	98
15) Phenanthrene	16.946	178	7191	0.376	ng/ul	99
16) Anthracene	17.034	178	6544	0.351	ng/ul	100
19) Fluoranthene	18.973	202	9014	0.370	ng/ul	98
20) Pyrene	19.335	202	9432	0.358	ng/ul#	96
21) Benzo(a)anthracene	21.094	228	9400	0.359	ng/ul	100
22) Chrysene	21.143	228	9279	0.366	ng/ul	99
24) Benzo(b)fluoranthene	22.616	252	11319	0.351	ng/ul	98
25) Benzo(k)fluoranthene	22.657	252	11240	0.355	ng/ul	98
26) Benzo(a)pyrene	23.157	252	10376	0.391	ng/ul	95
27) Indeno(1,2,3-cd)pyrene	25.370	276	17405	0.431	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.387	278	13785	0.431	ng/ul#	94
29) Benzo(g,h,i)perylene	26.011	276	14210	0.441	ng/ul	97

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

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