

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071124\
 Data File : BM046524.D
 Acq On : 11 Jul 2024 19:06
 Operator : MA/JU
 Sample : PB161731BS
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS731

Quant Time: Jul 12 04:22:58 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 04:18:37 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.514	152	2404	0.400	ng/ul	0.00
4) Naphthalene-d8	10.275	136	6014	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.158	164	3832	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.908	188	8061	0.400	ng/ul	0.00
17) Chrysene-d12	21.114	240	6192	0.400	ng/ul	0.00
23) Perylene-d12	23.256	264	7188	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.122	96	1211	0.613	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.876	152	3292	0.339	ng/ul	0.00
18) Fluoranthene-d10	18.945	212	6912	0.367	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.151	88	3178	1.508	ng/ul#	70
5) Naphthalene	10.325	128	4981	0.326	ng/ul	98
7) 2-Methylnaphthalene	11.953	142	3502	0.326	ng/ul	99
8) 1-Methylnaphthalene	12.172	142	3481	0.317	ng/ul	97
10) Acenaphthylene	13.872	152	5479	0.314	ng/ul	98
11) Acenaphthene	14.219	153	3886	0.333	ng/ul	96
12) Fluorene	15.213	166	4499	0.311	ng/ul	96
14) Pentachlorophenol	16.562	266	2023	0.495	ng/ul	99
15) Phenanthrene	16.950	178	7159	0.311	ng/ul	99
16) Anthracene	17.043	178	6921	0.308	ng/ul	98
19) Fluoranthene	18.978	202	8583	0.337	ng/ul	99
20) Pyrene	19.340	202	9027	0.328	ng/ul	99
21) Benzo(a)anthracene	21.096	228	8550	0.312	ng/ul	99
22) Chrysene	21.149	228	8358	0.315	ng/ul	100
24) Benzo(b)fluoranthene	22.622	252	9128	0.318	ng/ul	99
25) Benzo(k)fluoranthene	22.663	252	9108	0.323	ng/ul	98
26) Benzo(a)pyrene	23.163	252	8376	0.354	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.380	276	11351	0.316	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.407	278	8909	0.313	ng/ul	98
29) Benzo(g,h,i)perylene	26.027	276	8952	0.312	ng/ul	99

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

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