

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072624\
 Data File : BM046858.D
 Acq On : 27 Jul 2024 08:35
 Operator : MA/JU
 Sample : PB162152BS
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS152

Quant Time: Jul 28 23:29:05 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	7.463	152	3324	0.400	ng/ul	0.00
4) Naphthalene-d8	10.222	136	8707	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.109	164	5066	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.867	188	9295	0.400	ng/ul	0.00
17) Chrysene-d12	21.079	240	6018	0.400	ng/ul	0.00
23) Perylene-d12	23.206	264	7123	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.096	96	1363	0.586	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.822	152	4705	0.340	ng/ul	0.00
18) Fluoranthene-d10	18.909	212	7143	0.364	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.126	88	3750	1.408	ng/ul#	82
5) Naphthalene	10.271	128	7460	0.339	ng/ul	98
7) 2-Methylnaphthalene	11.899	142	5212	0.344	ng/ul	99
8) 1-Methylnaphthalene	12.119	142	5199	0.332	ng/ul	99
10) Acenaphthylene	13.827	152	7820	0.337	ng/ul	99
11) Acenaphthene	14.174	153	5558	0.353	ng/ul	99
12) Fluorene	15.168	166	6165	0.313	ng/ul	99
14) Pentachlorophenol	16.525	266	1761	0.419	ng/ul	98
15) Phenanthrene	16.905	178	8662	0.324	ng/ul	99
16) Anthracene	16.998	178	8201	0.330	ng/ul	98
19) Fluoranthene	18.937	202	9188	0.348	ng/ul	99
20) Pyrene	19.299	202	9473	0.346	ng/ul	99
21) Benzo(a)anthracene	21.061	228	8120	0.324	ng/ul	98
22) Chrysene	21.114	228	8064	0.317	ng/ul	98
24) Benzo(b)fluoranthene	22.575	252	8757	0.280	ng/ul	96
25) Benzo(k)fluoranthene	22.615	252	9053	0.287	ng/ul	97
26) Benzo(a)pyrene	23.112	252	8084	0.360	ng/ul#	98
27) Indeno(1,2,3-cd)pyrene	25.306	276	11651	0.337	ng/ul#	97
28) Dibenzo(a,h)anthracene	25.319	278	8823	0.349	ng/ul	98
29) Benzo(g,h,i)perylene	25.946	276	9433	0.310	ng/ul	98

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

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