

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM070823\
 Data File : BM040747.D
 Acq On : 08 Jul 2023 19:46
 Operator : MA/JU
 Sample : PB153845BS
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS845

Quant Time: Jul 08 23:32:09 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM062223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jul 08 22:50:20 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.895	152	3809	0.400	ng/ul	0.00
4) Naphthalene-d8	10.702	136	14815	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.541	164	6753	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.299	188	12016	0.400	ng/ul	0.00
17) Chrysene-d12	21.483	240	7313	0.400	ng/ul	0.00
23) Perylene-d12	23.880	264	5745	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.301	96	3549	0.593	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.291	152	6334	0.305	ng/ul	0.00
18) Fluoranthene-d10	19.324	212	8543	0.364	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.334	88	8502	1.396	ng/ul	88
5) Naphthalene	10.752	128	12338	0.302	ng/ul	98
7) 2-Methylnaphthalene	12.368	142	7214	0.301	ng/ul	97
8) 1-Methylnaphthalene	12.583	142	6809	0.284	ng/ul	99
10) Acenaphthylene	14.264	152	8893	0.295	ng/ul	96
11) Acenaphthene	14.606	153	7617	0.299	ng/ul	99
12) Fluorene	15.596	166	8115	0.298	ng/ul	98
14) Pentachlorophenol	16.957	266	549	0.237	ng/ul	97
15) Phenanthrene	17.337	178	10726	0.287	ng/ul	97
16) Anthracene	17.434	178	9558	0.305	ng/ul	95
19) Fluoranthene	19.357	202	10800	0.348	ng/ul	98
20) Pyrene	19.719	202	11445	0.344	ng/ul	97
21) Benzo(a)anthracene	21.469	228	7394	0.299	ng/ul	98
22) Chrysene	21.521	228	8644	0.306	ng/ul	99
24) Benzo(b)fluoranthene	23.149	252	7785	0.350	ng/ul	90
25) Benzo(k)fluoranthene	23.199	252	7524	0.343	ng/ul#	90
26) Benzo(a)pyrene	23.778	252	6341	0.317	ng/ul#	82
27) Indeno(1,2,3-cd)pyrene	26.371	276	8167	0.293	ng/ul#	95
28) Dibenzo(a,h)anthracene	26.384	278	6372	0.292	ng/ul#	91
29) Benzo(g,h,i)perylene	27.129	276	7335	0.300	ng/ul	95

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

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