

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM070523\
 Data File : BM040708.D
 Acq On : 07 Jul 2023 11:18
 Operator : MA/JU
 Sample : PB153582BS
 Misc :
 ALS Vial : 77 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS582

Quant Time: Jul 07 13:35:44 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 : Fri Jul 07 10:22:24 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.899	152	5086	0.400	ng/ul	0.00
4) Naphthalene-d8	10.707	136	18698	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.546	164	8228	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.303	188	13287	0.400	ng/ul	0.00
17) Chrysene-d12	21.492	240	8196	0.400	ng/ul	0.00
23) Perylene-d12	23.892	264	7873	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.309	96	4488	0.562	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.302	152	7909	0.302	ng/ul	0.00
18) Fluoranthene-d10	19.334	212	8906	0.339	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.343	88	11272	1.386	ng/ul	90
5) Naphthalene	10.757	128	15965	0.310	ng/ul	99
7) 2-Methylnaphthalene	12.374	142	9119	0.301	ng/ul	98
8) 1-Methylnaphthalene	12.588	142	8599	0.284	ng/ul	98
10) Acenaphthylene	14.268	152	10802	0.294	ng/ul	98
11) Acenaphthene	14.610	153	9216	0.297	ng/ul	98
12) Fluorene	15.600	166	9489	0.286	ng/ul	99
14) Pentachlorophenol	16.961	266	494	0.193	ng/ul	96
15) Phenanthrene	17.345	178	12046	0.292	ng/ul	97
16) Anthracene	17.438	178	10217	0.295	ng/ul	97
19) Fluoranthene	19.361	202	11244	0.323	ng/ul	99
20) Pyrene	19.724	202	11587	0.310	ng/ul	98
21) Benzo(a)anthracene	21.474	228	8428	0.304	ng/ul	98
22) Chrysene	21.527	228	9444	0.299	ng/ul	100
24) Benzo(b)fluoranthene	23.158	252	9464	0.311	ng/ul	89
25) Benzo(k)fluoranthene	23.202	252	9571	0.319	ng/ul#	87
26) Benzo(a)pyrene	23.783	252	8274	0.301	ng/ul#	81
27) Indeno(1,2,3-cd)pyrene	26.381	276	11538	0.302	ng/ul#	95
28) Dibenzo(a,h)anthracene	26.397	278	9054	0.303	ng/ul	93
29) Benzo(g,h,i)perylene	27.145	276	10378	0.310	ng/ul	98

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

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