

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM070323\
 Data File : BM040584.D
 Acq On : 03 Jul 2023 20:39
 Operator : MA/JU
 Sample : PB153508BS
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS508

Quant Time: Jul 04 01:30:07 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 : Mon Jul 03 13:40:38 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.923	152	5365	0.400	ng/ul	0.00
4) Naphthalene-d8	10.734	136	20083	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.568	164	8990	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.319	188	14754	0.400	ng/ul	0.00
17) Chrysene-d12	21.503	240	9461	0.400	ng/ul	0.00
23) Perylene-d12	23.914	264	10036	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.321	96	5057	0.600	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.323	152	9413	0.334	ng/ul	0.00
18) Fluoranthene-d10	19.347	212	10707	0.353	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.354	88	12207	1.423	ng/ul	93
5) Naphthalene	10.783	128	17213	0.311	ng/ul	99
7) 2-Methylnaphthalene	12.395	142	9992	0.307	ng/ul	98
8) 1-Methylnaphthalene	12.615	142	9547	0.294	ng/ul	99
10) Acenaphthylene	14.290	152	12868	0.320	ng/ul	98
11) Acenaphthene	14.628	153	10096	0.298	ng/ul	98
12) Fluorene	15.618	166	10379	0.287	ng/ul	99
14) Pentachlorophenol	16.977	266	732	0.258	ng/ul	92
15) Phenanthrene	17.362	178	13316	0.291	ng/ul	98
16) Anthracene	17.455	178	11765	0.305	ng/ul	97
19) Fluoranthene	19.375	202	12674	0.315	ng/ul	100
20) Pyrene	19.737	202	13357	0.310	ng/ul	97
21) Benzo(a)anthracene	21.485	228	11035	0.345	ng/ul	98
22) Chrysene	21.538	228	11453	0.314	ng/ul	99
24) Benzo(b)fluoranthene	23.178	252	11723	0.302	ng/ul	86
25) Benzo(k)fluoranthene	23.224	252	11905	0.311	ng/ul#	84
26) Benzo(a)pyrene	23.803	252	10401	0.297	ng/ul#	81
27) Indeno(1,2,3-cd)pyrene	26.414	276	14612	0.300	ng/ul#	92
28) Dibenzo(a,h)anthracene	26.430	278	11486	0.301	ng/ul	92
29) Benzo(g,h,i)perylene	27.185	276	12352	0.289	ng/ul	96

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

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