

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071723\
 Data File : BM040926.D
 Acq On : 17 Jul 2023 17:27
 Operator : MA/JU
 Sample : PB153935BS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS935

Quant Time: Jul 18 02:47:01 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 15 03:58:02 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.857	152	4622	0.400	ng/ul	0.00
4) Naphthalene-d8	10.660	136	14818	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.505	164	5915	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.258	188	9802	0.400	ng/ul	0.00
17) Chrysene-d12	21.449	240	6914	0.400	ng/ul	0.00
23) Perylene-d12	23.832	264	7273	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.280	96	4828	0.665	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.254	152	6162	0.297	ng/ul	0.00
18) Fluoranthene-d10	19.288	212	7228	0.326	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.314	88	13356	1.807	ng/ul	87
5) Naphthalene	10.709	128	14072	0.344	ng/ul	99
7) 2-Methylnaphthalene	12.326	142	7462	0.311	ng/ul	98
8) 1-Methylnaphthalene	12.546	142	7470	0.311	ng/ul	99
10) Acenaphthylene	14.227	152	8233	0.312	ng/ul	99
11) Acenaphthene	14.565	153	7241	0.324	ng/ul	97
12) Fluorene	15.560	166	6995	0.294	ng/ul	99
14) Pentachlorophenol	16.916	266	1710	0.906	ng/ul	92
15) Phenanthrene	17.300	178	9729	0.320	ng/ul	99
16) Anthracene	17.397	178	7643	0.299	ng/ul	99
19) Fluoranthene	19.321	202	9668	0.329	ng/ul	99
20) Pyrene	19.683	202	10147	0.322	ng/ul	98
21) Benzo(a)anthracene	21.435	228	7633	0.326	ng/ul	99
22) Chrysene	21.487	228	9057	0.340	ng/ul	100
24) Benzo(b)fluoranthene	23.104	252	10055	0.357	ng/ul	98
25) Benzo(k)fluoranthene	23.153	252	9630	0.347	ng/ul	99
26) Benzo(a)pyrene	23.723	252	8722	0.344	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.288	276	13131	0.372	ng/ul#	94
28) Dibenzo(a,h)anthracene	26.309	278	10482	0.379	ng/ul	97
29) Benzo(g,h,i)perylene	27.040	276	11671	0.377	ng/ul	98

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

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