

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062022\
 Data File : BM035598.D
 Acq On : 22 Jun 2022 07:52
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4107

Quant Time: Jun 22 08:24:01 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM061822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 20 01:56:13 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.871	152	4395	0.400	ng/ul	0.01
4) Naphthalene-d8	10.638	136	16796	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.465	164	8916	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.225	188	16360	0.400	ng/ul	0.00
17) Chrysene-d12	21.405	240	13073	0.400	ng/ul	0.00
23) Perylene-d12	23.750	264	12287	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.268	96	2808	0.454	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.244	152	9667	0.387	ng/ul	0.00
18) Fluoranthene-d10	19.248	212	17090	0.389	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.302	88	2678	0.435	ng/ul#	69
5) Naphthalene	10.693	128	18871	0.332	ng/ul	99
7) 2-Methylnaphthalene	12.316	142	11132	0.343	ng/ul	96
8) 1-Methylnaphthalene	12.519	142	11478	0.385	ng/ul	99
10) Acenaphthylene	14.196	152	17679	0.352	ng/ul	99
11) Acenaphthene	14.530	153	13613	0.366	ng/ul	99
12) Fluorene	15.538	166	14480	0.344	ng/ul	99
14) Pentachlorophenol	16.926	266	1128	0.359	ng/ul	96
15) Phenanthrene	17.268	178	21251	0.354	ng/ul	100
16) Anthracene	17.377	178	17657	0.336	ng/ul	100
19) Fluoranthene	19.281	202	23408	0.350	ng/ul	99
20) Pyrene	19.638	202	26800	0.383	ng/ul	99
21) Benzo(a)anthracene	21.394	228	16690	0.338	ng/ul	100
22) Chrysene	21.443	228	21107	0.359	ng/ul	100
24) Benzo(b)fluoranthene	23.036	252	18849	0.350	ng/ul	98
25) Benzo(k)fluoranthene	23.083	252	19473	0.349	ng/ul	98
26) Benzo(a)pyrene	23.653	252	19723	0.354	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	26.191	276	23556	0.367	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.194	278	19164	0.365	ng/ul	99
29) Benzo(g,h,i)perylene	26.922	276	21259	0.370	ng/ul	99

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

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