

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010323\
 Data File : BM038355.D
 Acq On : 03 Jan 2023 11:10
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4059

Quant Time: Jan 04 00:42:29 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM123022.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 30 01:43:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.009	152	1379	0.400	ng/u1	0.00
4) Naphthalene-d8	10.818	136	3825	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.638	164	2073	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.379	188	4225	0.400	ng/u1	0.00
17) Chrysene-d12	21.551	240	4266	0.400	ng/u1	0.00
23) Perylene-d12	23.980	264	4910	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.372	96	811	0.395	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.401	152	2033	0.354	ng/u1	0.00
18) Fluoranthene-d10	19.399	212	4483	0.371	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.410	88	805	0.382	ng/u1	97
5) Naphthalene	10.867	128	3702	0.363	ng/u1	98
7) 2-Methylnaphthalene	12.478	142	2223	0.354	ng/u1	100
8) 1-Methylnaphthalene	12.693	142	2254	0.353	ng/u1	98
10) Acenaphthylene	14.361	152	3530	0.366	ng/u1	100
11) Acenaphthene	14.699	153	2559	0.357	ng/u1	97
12) Fluorene	15.684	166	2851	0.351	ng/u1	98
14) Pentachlorophenol	17.046	266	514	0.341	ng/u1	92
15) Phenanthrene	17.422	178	4340	0.348	ng/u1	99
16) Anthracene	17.515	178	4055	0.353	ng/u1	98
19) Fluoranthene	19.431	202	5462	0.370	ng/u1	100
20) Pyrene	19.794	202	5793	0.367	ng/u1	99
21) Benzo(a)anthracene	21.533	228	5130	0.361	ng/u1	99
22) Chrysene	21.589	228	5395	0.360	ng/u1	99
24) Benzo(b)fluoranthene	23.237	252	6314	0.355	ng/u1	99
25) Benzo(k)fluoranthene	23.284	252	6302	0.356	ng/u1	99
26) Benzo(a)pyrene	23.874	252	6048	0.360	ng/u1	99
27) Indeno(1,2,3-cd)pyrene	26.522	276	8750	0.356	ng/u1#	99
28) Dibenzo(a,h)anthracene	26.535	278	6902	0.358	ng/u1	98
29) Benzo(g,h,i)perylene	27.303	276	7861	0.360	ng/u1	98

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

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