

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062022\
 Data File : BM035578.D
 Acq On : 21 Jun 2022 18:35
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 57 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4105

Quant Time: Jun 22 01:33:07 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	7.879	152	3506	0.400	ng/ul	0.02
4) Naphthalene-d8	10.649	136	14275	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.473	164	7961	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.229	188	15883	0.400	ng/ul	0.00
17) Chrysene-d12	21.409	240	11691	0.400	ng/ul	0.00
23) Perylene-d12	23.759	264	10348	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.272	96	2206	0.447	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.255	152	8377	0.395	ng/ul	0.01
18) Fluoranthene-d10	19.252	212	16145	0.411	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.306	88	2148	0.437	ng/ul#	68
5) Naphthalene	10.699	128	15919	0.329	ng/ul	100
7) 2-Methylnaphthalene	12.332	142	9311	0.337	ng/ul	95
8) 1-Methylnaphthalene	12.525	142	9865	0.389	ng/ul	100
10) Acenaphthylene	14.200	152	15847	0.353	ng/ul	100
11) Acenaphthene	14.533	153	12129	0.365	ng/ul	98
12) Fluorene	15.542	166	13432	0.358	ng/ul	99
14) Pentachlorophenol	16.938	266	950	0.312	ng/ul	97
15) Phenanthrene	17.271	178	20273	0.348	ng/ul	99
16) Anthracene	17.381	178	17130	0.336	ng/ul	99
19) Fluoranthene	19.284	202	22066	0.369	ng/ul	100
20) Pyrene	19.647	202	23181	0.370	ng/ul	99
21) Benzo(a)anthracene	21.398	228	14544	0.329	ng/ul	98
22) Chrysene	21.444	228	18884	0.359	ng/ul	99
24) Benzo(b)fluoranthene	23.043	252	15863	0.350	ng/ul	99
25) Benzo(k)fluoranthene	23.093	252	16770	0.357	ng/ul	97
26) Benzo(a)pyrene	23.663	252	16828	0.358	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	26.202	276	19521	0.361	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.212	278	15904	0.360	ng/ul	97
29) Benzo(g,h,i)perylene	26.936	276	17653	0.364	ng/ul	98

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

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