

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081221\
 Data File : BM031727.D
 Acq On : 13 Aug 2021 22:44
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4117

Quant Time: Aug 13 23:17:49 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM072821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 12 17:28:09 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.569	152	1332	0.400	ng/ul	0.00
4) Naphthalene-d8	10.329	136	4863	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.201	164	2665	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.957	188	5608	0.400	ng/ul	0.00
17) Chrysene-d12	21.151	240	4838	0.400	ng/ul	0.00
23) Perylene-d12	23.309	264	4998	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.185	96	586	0.396	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.926	152	2994	0.365	ng/ul	0.00
18) Fluoranthene-d10	18.989	212	5936	0.368	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.217	88	558	0.373	ng/ul#	73
5) Naphthalene	10.379	128	5489	0.379	ng/ul	99
7) 2-Methylnaphthalene	12.002	142	3607	0.364	ng/ul	99
8) 1-Methylnaphthalene	12.218	142	3473	0.355	ng/ul	99
10) Acenaphthylene	13.923	152	4619	0.407	ng/ul	99
11) Acenaphthene	14.265	153	3569	0.374	ng/ul	98
12) Fluorene	15.262	166	4083	0.371	ng/ul	98
14) Pentachlorophenol	16.617	266	628	0.350	ng/ul	99
15) Phenanthrene	16.999	178	6561	0.368	ng/ul	99
16) Anthracene	17.089	178	6246	0.375	ng/ul	99
19) Fluoranthene	19.020	202	7580	0.367	ng/ul	97
20) Pyrene	19.382	202	7781	0.371	ng/ul	97
21) Benzo(a)anthracene	21.134	228	7244	0.371	ng/ul	98
22) Chrysene	21.188	228	7359	0.364	ng/ul	99
24) Benzo(b)fluoranthene	22.665	252	7815	0.347	ng/ul	93
25) Benzo(k)fluoranthene	22.708	252	7648	0.329	ng/ul#	97
26) Benzo(a)pyrene	23.212	252	6893	0.350	ng/ul#	91
27) Indeno(1,2,3-cd)pyrene	25.439	276	9065	0.356	ng/ul#	96
28) Dibenzo(a,h)anthracene	25.451	278	7330	0.357	ng/ul	99
29) Benzo(g,h,i)perylene	26.088	276	7764	0.357	ng/ul	94

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

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