

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081221\
 Data File : BM031712.D
 Acq On : 13 Aug 2021 13:17
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4116

Quant Time: Aug 13 13:54:39 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	7.570	152	1290	0.400	ng/ul	0.00
4) Naphthalene-d8	10.329	136	4767	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.205	164	2522	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.961	188	4828	0.400	ng/ul	0.00
17) Chrysene-d12	21.154	240	3948	0.400	ng/ul	# 0.00
23) Perylene-d12	23.314	264	3966	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.186	96	548	0.382	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.926	152	2850	0.355	ng/ul	0.00
18) Fluoranthene-d10	18.989	212	4972	0.377	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.217	88	534	0.369	ng/ul#	84
5) Naphthalene	10.379	128	5341	0.376	ng/ul	99
7) 2-Methylnaphthalene	12.002	142	3456	0.356	ng/ul	97
8) 1-Methylnaphthalene	12.223	142	3360	0.351	ng/ul	100
10) Acenaphthylene	13.923	152	4485	0.418	ng/ul	100
11) Acenaphthene	14.265	153	3442	0.381	ng/ul	100
12) Fluorene	15.263	166	3702	0.356	ng/ul	97
14) Pentachlorophenol	16.621	266	503	0.325	ng/ul	99
15) Phenanthrene	16.999	178	5670	0.370	ng/ul	99
16) Anthracene	17.093	178	5414	0.377	ng/ul	98
19) Fluoranthene	19.020	202	6378	0.378	ng/ul#	95
20) Pyrene	19.386	202	6599	0.386	ng/ul	96
21) Benzo(a)anthracene	21.142	228	5975	0.375	ng/ul	98
22) Chrysene	21.193	228	6076	0.368	ng/ul	99
24) Benzo(b)fluoranthene	22.672	252	6166	0.345	ng/ul	91
25) Benzo(k)fluoranthene	22.716	252	6398	0.347	ng/ul#	95
26) Benzo(a)pyrene	23.222	252	5576	0.356	ng/ul#	89
27) Indeno(1,2,3-cd)pyrene	25.449	276	7112	0.352	ng/ul#	96
28) Dibenzo(a,h)anthracene	25.461	278	5686	0.349	ng/ul	95
29) Benzo(g,h,i)perylene	26.098	276	6095	0.353	ng/ul	95

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

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