

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072221\
 Data File : BM031087.D
 Acq On : 23 Jul 2021 12:08
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4061

Quant Time: Jul 23 12:39:12 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM072221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 22 11:33:47 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.632	152	1656	0.400	ng/ul	0.00
4) Naphthalene-d8	10.397	136	6680	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.262	164	4539	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.009	188	9900	0.400	ng/ul	0.00
17) Chrysene-d12	21.195	240	8868	0.400	ng/ul	0.00
23) Perylene-d12	23.370	264	7407	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.217	96	609	0.350	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.989	152	4391	0.374	ng/ul	0.00
18) Fluoranthene-d10	19.039	212	11152	0.412	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.248	88	563	0.331	ng/ul#	88
5) Naphthalene	10.446	128	7470	0.375	ng/ul	99
7) 2-Methylnaphthalene	12.065	142	5457	0.380	ng/ul	99
8) 1-Methylnaphthalene	12.286	142	5378	0.383	ng/ul	99
10) Acenaphthylene	13.981	152	7210	0.388	ng/ul	98
11) Acenaphthene	14.326	153	6158	0.375	ng/ul	98
12) Fluorene	15.316	166	7117	0.368	ng/ul	97
14) Pentachlorophenol	16.659	266	1210	0.353	ng/ul	98
15) Phenanthrene	17.051	178	11961	0.380	ng/ul	99
16) Anthracene	17.141	178	11063	0.376	ng/ul	98
19) Fluoranthene	19.069	202	14345	0.402	ng/ul#	96
20) Pyrene	19.432	202	14479	0.397	ng/ul	96
21) Benzo(a)anthracene	21.181	228	13644	0.385	ng/ul	99
22) Chrysene	21.231	228	13709	0.358	ng/ul	99
24) Benzo(b)fluoranthene	22.720	252	13207	0.379	ng/ul	97
25) Benzo(k)fluoranthene	22.764	252	13060	0.370	ng/ul	97
26) Benzo(a)pyrene	23.275	252	11057	0.367	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	25.524	276	11790	0.299	ng/ul#	95
28) Dibenzo(a,h)anthracene	25.543	278	9488	0.300	ng/ul	99
29) Benzo(g,h,i)perylene	26.183	276	9522	0.284	ng/ul	97

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

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