

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM100820\
 Data File : BM027645.D
 Acq On : 08 Oct 2020 12:30
 Operator : JU/CG
 Sample : SST00.203
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST00.2103

Quant Time: Oct 08 13:42:28 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SFAM-EPA-SIM-BM100820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 08 13:40:34 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.79	152	853	0.40	ng/ul	0.00
2) Naphthalene-d8	10.57	136	2352	0.40	ng/ul	0.00
7) Acenaphthene-d10	14.42	164	1575	0.40	ng/ul	0.00
11) Phenanthrene-d10	17.16	188	3934	0.40	ng/ul	0.00
17) Chrysene-d12	21.34	240	4068	0.40	ng/ul	0.00
21) Perylene-d12	23.59	264	4179	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.16	152	850	0.18	ng/ul	0.00
15) Fluoranthene-d10	19.18	212	2479	0.18	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.62	128	1320	0.202	ng/ul#	93
5) 2-Methylnaphthalene	12.24	142	934	0.185	ng/ul	95
6) 1-Methylnaphthalene	12.45	142	1073	0.186	ng/ul	99
8) Acenaphthylene	14.13	152	1322	0.186	ng/ul	95
9) Acenaphthene	14.48	153	1063	0.201	ng/ul	100
10) Fluorene	15.47	166	1346	0.193	ng/ul	97
12) Pentachlorophenol	16.81	266	313	0.203	ng/ul	99
13) Phenanthrene	17.20	178	2327	0.199	ng/ul	95
14) Anthracene	17.29	178	2200	0.195	ng/ul	95
16) Fluoranthene	19.21	202	3076	0.186	ng/ul	95
18) Pyrene	19.58	202	2381	0.154	ng/ul#	95
19) Benzo(a)anthracene	21.32	228	3098	0.193	ng/ul	97
20) Chrysene	21.37	228	3315	0.202	ng/ul	99
22) Benzo(b)fluoranthene	22.91	252	3505	0.204	ng/ul	85
23) Benzo(k)fluoranthene	22.96	252	3359	0.198	ng/ul#	84
24) Benzo(a)pyrene	23.49	252	2944	0.191	ng/ul#	79
25) Indeno(1,2,3-cd)pyrene	25.86	276	4070	0.199	ng/ul#	99
26) Dibenzo(a,h)anthracene	25.87	278	3329	0.196	ng/ul#	86
27) Benzo(a,h,i)perylene	26.55	276	3615	0.206	ng/ul#	96

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

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