

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM033020\
 Data File : BM025855.D
 Acq On : 30 Mar 2020 12:00
 Operator : CG/JU
 Sample : SST0.476
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.476

Quant Time: Mar 30 12:33:17 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM033020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 30 12:31:41 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	1804	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	6952	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	4295	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	9687	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	8842	0.40	ng/ul	0.00
20) Perylene-d12	23.29	264	8801	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.93	152	3999	0.31	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	8995	0.29	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.38	128	6899	0.330	ng/ul	100
5) 2-Methylnaphthalene	12.00	142	4680	0.307	ng/ul	100
7) Acenaphthylene	13.92	152	5949	0.296	ng/ul	100
8) Acenaphthene	14.27	153	4976	0.319	ng/ul	100
9) Fluorene	15.26	166	5826	0.298	ng/ul	100
11) Pentachlorophenol	16.62	266	542	0.214	ng/ul	100
12) Phenanthrene	16.99	178	9842	0.315	ng/ul	100
13) Anthracene	17.08	178	8270	0.298	ng/ul	100
15) Fluoranthene	19.01	202	11210	0.287	ng/ul	100
17) Pyrene	19.38	202	11133	0.332	ng/ul	100
18) Benzo(a)anthracene	21.13	228	9358	0.288	ng/ul	100
19) Chrysene	21.18	228	11337	0.317	ng/ul	100
21) Benzo(b)fluoranthene	22.66	252	10884	0.309	ng/ul	100
22) Benzo(k)fluoranthene	22.70	252	11185	0.303	ng/ul	100
23) Benzo(a)pyrene	23.20	252	8840	0.299	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.41	276	11105	0.301	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.43	278	8628	0.283	ng/ul	100
26) Benzo(g,h,i)perylene	26.05	276	10923	0.349	ng/ul	100

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

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