

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM033020\
 Data File : BM025856.D
 Acq On : 30 Mar 2020 12:36
 Operator : CG/JU
 Sample : SSTD0.877
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.877

Quant Time: Mar 30 13:10:11 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	1443	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	5285	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.21	164	3233	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	7516	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	7393	0.40	ng/ul	0.00
20) Perylene-d12	23.29	264	7194	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.93	152	7627	0.77	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	18536	0.76	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.38	128	13015	0.818	ng/ul	99
5) 2-Methylnaphthalene	12.00	142	8894	0.768	ng/ul	99
7) Acenaphthylene	13.92	152	11394	0.753	ng/ul	97
8) Acenaphthene	14.27	153	9630	0.820	ng/ul	98
9) Fluorene	15.26	166	11699	0.795	ng/ul	99
11) Pentachlorophenol	16.61	266	1283	0.652	ng/ul	93
12) Phenanthrene	16.99	178	19806	0.817	ng/ul	99
13) Anthracene	17.08	178	16927	0.787	ng/ul	98
15) Fluoranthene	19.01	202	23445	0.773	ng/ul	99
17) Pyrene	19.37	202	23050	0.822	ng/ul	100
18) Benzo(a)anthracene	21.13	228	20070	0.739	ng/ul	100
19) Chrysene	21.18	228	23926	0.801	ng/ul	99
21) Benzo(b)fluoranthene	22.66	252	23195	0.806	ng/ul	95
22) Benzo(k)fluoranthene	22.70	252	23743	0.787	ng/ul	97
23) Benzo(a)pyrene	23.20	252	18680	0.773	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.41	276	23784	0.790	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.43	278	18626	0.746	ng/ul	95
26) Benzo(g,h,i)perylene	26.05	276	23062	0.902	ng/ul	98

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

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