

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030620\
 Data File : BM025315.D
 Acq On : 06 Mar 2020 14:44
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.413

Quant Time: Mar 06 16:44:44 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 04 08:02:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	3491	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	14779	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	10426	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	24971	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	22203	0.40	ng/ul	0.00
20) Perylene-d12	23.42	264	18259	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	10004	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.07	212	27613	0.34	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.49	128	15893	0.357	ng/ul#	82
5) 2-Methylnaphthalene	12.11	142	11657	0.360	ng/ul	99
7) Acenaphthylene	14.02	152	16866	0.345	ng/ul	93
8) Acenaphthene	14.36	153	13645	0.360	ng/ul	99
9) Fluorene	15.35	166	16878	0.356	ng/ul	95
11) Pentachlorophenol	16.70	266	1928	0.295	ng/ul	95
12) Phenanthrene	17.08	178	28860	0.358	ng/ul	95
13) Anthracene	17.17	178	25354	0.355	ng/ul	95
15) Fluoranthene	19.10	202	34141	0.339	ng/ul	95
17) Pyrene	19.46	202	33626	0.399	ng/ul	96
18) Benzo(a)anthracene	21.21	228	29775	0.365	ng/ul	98
19) Chrysene	21.27	228	32621	0.364	ng/ul	98
21) Benzo(b)fluoranthene	22.77	252	27121	0.371	ng/ul	95
22) Benzo(k)fluoranthene	22.81	252	28934	0.378	ng/ul#	96
23) Benzo(a)pyrene	23.33	252	22194	0.362	ng/ul#	94
24) Indeno(1,2,3-cd)pyrene	25.60	276	26018	0.340	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.62	278	21701	0.343	ng/ul#	94
26) Benzo(g,h,i)perylene	26.27	276	21475	0.331	ng/ul#	91

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

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