

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060217\
 Data File : BM010412.D
 Acq On : 03 Jun 2017 05:59
 Operator : SJ/MA
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.419

Quant Time: Jun 05 04:04:11 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM053017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 05 04:03:07 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	728	0.40	ng/ul	0.00
2) Naphthalene-d8	10.72	136	2119	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.55	164	1177	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.30	188	2574	0.40	ng/ul	0.00
16) Chrysene-d12	21.46	240	2709	0.40	ng/ul	0.00
20) Perylene-d12	23.81	264	2756	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.30	152	1238	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	2859	0.42	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.77	128	2167	0.388	ng/ul	97
5) 2-Methylnaphthalene	12.37	142	1597	0.405	ng/ul	100
7) Acenaphthylene	14.26	152	2433	0.414	ng/ul	95
8) Acenaphthene	14.61	153	1733	0.390	ng/ul	99
9) Fluorene	15.60	166	2083	0.392	ng/ul	92
11) Pentachlorophenol	16.95	266	479	0.487	ng/ul	98
12) Phenanthrene	17.33	178	3301	0.399	ng/ul	99
13) Anthracene	17.42	178	3206	0.414	ng/ul	97
15) Fluoranthene	19.34	202	4256	0.415	ng/ul	96
17) Pyrene	19.71	202	4732	0.410	ng/ul#	92
18) Benzo(a)anthracene	21.44	228	4655	0.423	ng/ul	97
19) Chrysene	21.50	228	4006	0.376	ng/ul	96
21) Benzo(b)fluoranthene	23.08	252	4943	0.410	ng/ul	84
22) Benzo(k)fluoranthene	23.13	252	4407	0.384	ng/ul#	83
23) Benzo(a)pyrene	23.69	252	4781	0.419	ng/ul#	86
24) Indeno(1,2,3-cd)pyrene	26.20	276	5739	0.382	ng/ul#	93
25) Dibenzo(a,h)anthracene	26.21	278	4637	0.388	ng/ul	93
26) Benzo(g,h,i)perylene	26.94	276	4855	0.375	ng/ul	97

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

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