

Data Path : Z:\HPCHEM1\BNA_M\Data\BM060817\
 Data File : BM010444.D
 Acq On : 07 Jun 2017 23:22
 Operator : SJ/MA
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.423

Quant Time: Jun 08 07:10:50 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM053017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 08 07:10:23 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	697	0.40	ng/ul	0.00
2) Naphthalene-d8	10.72	136	2065	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	1271	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.28	188	2907	0.40	ng/ul	-0.02
16) Chrysene-d12	21.46	240	2252	0.40	ng/ul	0.00
20) Perylene-d12	23.80	264	2247	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	1302	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.31	212	2957	0.38	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.77	128	2108	0.387	ng/ul	98
5) 2-Methylnaphthalene	12.36	142	1593	0.415	ng/ul	100
7) Acenaphthylene	14.26	152	2511	0.396	ng/ul	100
8) Acenaphthene	14.61	153	1833	0.382	ng/ul	92
9) Fluorene	15.60	166	2221	0.387	ng/ul	99
11) Pentachlorophenol	16.94	266	449	0.404	ng/ul	96
12) Phenanthrene	17.33	178	3581	0.383	ng/ul	95
13) Anthracene	17.42	178	3363	0.384	ng/ul	98
15) Fluoranthene	19.34	202	4323	0.373	ng/ul	95
17) Pyrene	19.70	202	4394	0.458	ng/ul	96
18) Benzo(a)anthracene	21.44	228	3689	0.403	ng/ul	98
19) Chrysene	21.49	228	3544	0.400	ng/ul	96
21) Benzo(b)fluoranthene	23.08	252	3723	0.379	ng/ul	86
22) Benzo(k)fluoranthene	23.13	252	3790	0.405	ng/ul#	82
23) Benzo(a)pyrene	23.69	252	3699	0.397	ng/ul#	83
24) Indeno(1,2,3-cd)pyrene	26.18	276	4754	0.389	ng/ul#	89
25) Dibenzo(a,h)anthracene	26.19	278	3783	0.388	ng/ul#	90
26) Benzo(g,h,i)perylene	26.93	276	4046	0.384	ng/ul#	91

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

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