

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM051715\
 Data File : BM001322.D
 Acq On : 15 May 2015 10:00
 Operator : TP/IZ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.481

Quant Time: May 16 02:25:36 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM051115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 14 01:20:45 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.43	152	1123	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.23	136	3891	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.12	164	1996	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.86	188	4054	0.40	ng/ul	0.00
16) Chrysene-d12	21.08	240	3336	0.40	ng/ul	0.00
20) Perylene-d12	23.18	264	3085	0.40	ng/ul	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	11.83	152	2009	0.41	ng/ul	0.00
14) Fluoranthene-d10	18.90	212	3495	0.39	ng/ul	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.26	128	3989	0.41	ng/ul	97
5) 2-Methylnaphthalene	11.90	142	2621	0.41	ng/ul	99
7) Acenaphthylene	13.82	152	3550	0.40	ng/ul	99
8) Acenaphthene	14.17	153	2811	0.40	ng/ul	94
9) Fluorene	15.18	166	3027	0.39	ng/ul	95
11) Pentachlorophenol	16.53	266	235	0.27	ng/ul	95
12) Phenanthrene	16.90	178	4670	0.41	ng/ul	97
13) Anthracene	17.00	178	4296	0.41	ng/ul	99
15) Fluoranthene	18.94	202	5132	0.40	ng/ul	98
17) Pyrene	19.30	202	5275	0.41	ng/ul	99
18) Benzo(a)anthracene	21.06	228	4258	0.42	ng/ul	98
19) Chrysene	21.11	228	4800	0.41	ng/ul	98
21) Benzo(b)fluoranthene	22.56	252	4683	0.42	ng/ul	99
22) Benzo(k)fluoranthene	22.60	252	4670	0.41	ng/ul	99
23) Benzo(a)pyrene	23.09	252	4374	0.41	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.27	276	4967	0.42	ng/ul	98
25) Dibenzo(a,h)anthracene	25.28	278	3921	0.42	ng/ul	97
26) Benzo(g,h,i)perylene	25.90	276	4466	0.43	ng/ul	98

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

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