

Data Path : Z:\HPCHEM1\BNA_M\Data\BM021715\
 Data File : BM000537.D
 Acq On : 17 Feb 2015 12:13
 Operator : TP/IZ
 Sample : SSTD0.415
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC0.4

Quant Time: Feb 18 03:50:53 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-SIM-BM021215.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 12 16:45:55 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	2940	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.47	136	9614	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.34	164	5448	0.40	ng/ul	-0.02
10) Phenanthrene-d10	17.07	188	10557	0.40	ng/ul	-0.02
16) Chrysene-d12	21.27	240	8936	0.40	ng/ul	-0.02
20) Perylene-d12	23.49	264	7803	0.40	ng/ul	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.07	152	4729	0.39	ng/ul	-0.02
14) Fluoranthene-d10	19.12	212	8744	0.39	ng/ul	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.51	128	8960	0.40	ng/ul	98
5) 2-Methylnaphthalene	12.14	142	5898	0.39	ng/ul	100
7) Acenaphthylene	14.05	152	7944	0.39	ng/ul	99
8) Acenaphthene	14.38	153	6172	0.39	ng/ul#	80
9) Fluorene	15.39	166	7233	0.38	ng/ul	97
11) Pentachlorophenol	16.73	266	342	0.25	ng/ul#	91
12) Phenanthrene	17.12	178	11233	0.39	ng/ul	98
13) Anthracene	17.21	178	10036	0.39	ng/ul	97
15) Fluoranthene	19.14	202	12345	0.39	ng/ul	95
17) Pyrene	19.50	202	12728	0.40	ng/ul	95
18) Benzo(a)anthracene	21.25	228	9817	0.39	ng/ul	95
19) Chrysene	21.30	228	11459	0.39	ng/ul	94
21) Benzo(b)fluoranthene	22.82	252	10784	0.41	ng/ul	98
22) Benzo(k)fluoranthene	22.87	252	10070	0.38	ng/ul	99
23) Benzo(a)pyrene	23.39	252	9719	0.39	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.72	276	10900	0.40	ng/ul	99
25) Dibenzo(a,h)anthracene	25.73	278	8728	0.40	ng/ul	97
26) Benzo(g,h,i)perylene	26.40	276	9614	0.40	ng/ul	100

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

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