

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM021215\
 Data File : BM000504.D
 Acq On : 12 Feb 2015 16:04
 Operator : TP/IZ
 Sample : SSTD00102
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD00102

Quant Time: Feb 12 16:50:01 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 16:25:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.71	152	3365	0.40	ng/ul	0.00
2) Naphthalene-d8	10.49	136	11147	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.35	164	6405	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.09	188	13473	0.40	ng/ul	0.00
16) Chrysene-d12	21.29	240	11611	0.40	ng/ul	0.00
20) Perylene-d12	23.52	264	9993	0.40	ng/ul	0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.09	152	13976	0.94	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	27711	0.93	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.54	128	25916	0.95	ng/ul	99
5) 2-Methylnaphthalene	12.16	142	17667	0.93	ng/ul	99
7) Acenaphthylene	14.07	152	25179	0.98	ng/ul	99
8) Acenaphthene	14.41	153	19096	0.97	ng/ul	98
9) Fluorene	15.40	166	23015	0.96	ng/ul	99
11) Pentachlorophenol	16.75	266	1878	1.40	ng/ul	95
12) Phenanthrene	17.13	178	35941	0.95	ng/ul	99
13) Anthracene	17.23	178	32569	0.96	ng/ul	99
15) Fluoranthene	19.16	202	40410	0.92	ng/ul	99
17) Pyrene	19.53	202	40894	0.97	ng/ul	99
18) Benzo(a)anthracene	21.27	228	32943	0.96	ng/ul	97
19) Chrysene	21.32	228	35985	0.92	ng/ul	96
21) Benzo(b)fluoranthene	22.84	252	32833	0.95	ng/ul	94
22) Benzo(k)fluoranthene	22.89	252	34132	0.95	ng/ul	95
23) Benzo(a)pyrene	23.42	252	31240	0.97	ng/ul	92
24) Indeno(1,2,3-cd)pyrene	25.76	276	35313	1.02	ng/ul	100
25) Dibenzo(a,h)anthracene	25.77	278	28239	1.03	ng/ul	95
26) Benzo(g,h,i)perylene	26.45	276	30764	1.02	ng/ul	98

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

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