

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM021215\
 Data File : BM000502.D
 Acq On : 12 Feb 2015 14:44
 Operator : TP/IZ
 Sample : SSTD0.499
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.499

Quant Time: Feb 12 16:34:46 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	2941	0.40	ng/ul	0.00
2) Naphthalene-d8	10.49	136	9650	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.35	164	5776	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.09	188	11895	0.40	ng/ul	0.00
16) Chrysene-d12	21.29	240	10943	0.40	ng/ul	0.00
20) Perylene-d12	23.51	264	9902	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.09	152	4927	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	10109	0.38	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.54	128	9107	0.39	ng/ul	100
5) 2-Methylnaphthalene	12.16	142	6130	0.37	ng/ul	100
7) Acenaphthylene	14.08	152	8445	0.37	ng/ul	100
8) Acenaphthene	14.41	153	6693	0.38	ng/ul	100
9) Fluorene	15.40	166	7928	0.37	ng/ul	100
11) Pentachlorophenol	16.76	266	558	0.47	ng/ul	100
12) Phenanthrene	17.13	178	12653	0.38	ng/ul	100
13) Anthracene	17.23	178	11468	0.38	ng/ul	100
15) Fluoranthene	19.16	202	14709	0.38	ng/ul	100
17) Pyrene	19.53	202	15052	0.38	ng/ul	100
18) Benzo(a)anthracene	21.27	228	12162	0.38	ng/ul	100
19) Chrysene	21.32	228	14451	0.39	ng/ul	100
21) Benzo(b)fluoranthene	22.85	252	13417	0.39	ng/ul	100
22) Benzo(k)fluoranthene	22.89	252	12945	0.36	ng/ul	100
23) Benzo(a)pyrene	23.42	252	12363	0.39	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.75	276	13941	0.40	ng/ul	100
25) Dibenzo(a,h)anthracene	25.77	278	10965	0.40	ng/ul	100
26) Benzo(g,h,i)perylene	26.44	276	12292	0.41	ng/ul	100

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

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