

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM021115\
 Data File : BM000478.D
 Acq On : 11 Feb 2015 12:56
 Operator : TP/IZ
 Sample : SSTD0.499
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD0.499

Quant Time: Feb 11 13:54:28 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-SIM-BM021115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 11 13:45:17 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	2889	0.40	ng/ul	0.00
2) Naphthalene-d8	10.50	136	9994	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	6646	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	13837	0.40	ng/ul	0.00
16) Chrysene-d12	21.30	240	11970	0.40	ng/ul	0.00
20) Perylene-d12	23.53	264	9871	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.10	152	5246	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.14	212	11787	0.37	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.55	128	9309	0.34	ng/ul	100
5) 2-Methylnaphthalene	12.17	142	6565	0.35	ng/ul	100
7) Acenaphthylene	14.09	152	9572	0.35	ng/ul	100
8) Acenaphthene	14.41	153	7461	0.35	ng/ul	100
9) Fluorene	15.42	166	9188	0.36	ng/ul	100
11) Pentachlorophenol	16.76	266	484	0.48	ng/ul	100
12) Phenanthrene	17.15	178	15008	0.34	ng/ul	100
13) Anthracene	17.24	178	13528	0.34	ng/ul	100
15) Fluoranthene	19.18	202	16910	0.34	ng/ul	100
17) Pyrene	19.54	202	17191	0.33	ng/ul	100
18) Benzo(a)anthracene	21.28	228	13179	0.34	ng/ul	100
19) Chrysene	21.33	228	15863	0.34	ng/ul	100
21) Benzo(b)fluoranthene	22.86	252	13253	0.34	ng/ul	100
22) Benzo(k)fluoranthene	22.90	252	14221	0.36	ng/ul	100
23) Benzo(a)pyrene	23.43	252	12371	0.34	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.78	276	12535	0.31	ng/ul	100
25) Dibenzo(a,h)anthracene	25.79	278	9835	0.30	ng/ul	100
26) Benzo(g,h,i)perylene	26.47	276	10879	0.30	ng/ul	100

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

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