

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM020515\
 Data File : BM000426.D
 Acq On : 05 Feb 2015 12:17
 Operator : TP/IZ
 Sample : SSTD0.877
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.877

Quant Time: Feb 05 13:08:28 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM020515.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 05 12:58:49 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	2216	0.40	ng/ul	0.00
2) Naphthalene-d8	10.51	136	7126	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	4318	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.12	188	8793	0.40	ng/ul	0.00
16) Chrysene-d12	21.31	240	7736	0.40	ng/ul	0.00
20) Perylene-d12	23.55	264	6767	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	7508	0.77	ng/ul	0.00
14) Fluoranthene-d10	19.15	212	14730	0.77	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.56	128	14024	0.78	ng/ul	99
5) 2-Methylnaphthalene	12.18	142	9747	0.76	ng/ul	99
7) Acenaphthylene	14.10	152	14172	0.79	ng/ul	98
8) Acenaphthene	14.43	153	10244	0.77	ng/ul	100
9) Fluorene	15.43	166	12383	0.75	ng/ul	99
11) Pentachlorophenol	16.78	266	1532	1.37	ng/ul	95
12) Phenanthrene	17.16	178	19425	0.76	ng/ul	99
13) Anthracene	17.16	178	19425	0.83	ng/ul	99
15) Fluoranthene	19.19	202	22513	0.77	ng/ul	99
17) Pyrene	19.55	202	22760	0.72	ng/ul	98
18) Benzo(a)anthracene	21.29	228	19086	0.80	ng/ul	97
19) Chrysene	21.34	228	20536	0.76	ng/ul	99
21) Benzo(b)fluoranthene	22.88	252	19071	0.75	ng/ul	95
22) Benzo(k)fluoranthene	22.88	252	19071	0.72	ng/ul	92
23) Benzo(a)pyrene	23.45	252	17774	0.76	ng/ul	92
24) Indeno(1,2,3-cd)pyrene	25.82	276	20307	0.86	ng/ul	100
25) Dibenzo(a,h)anthracene	25.83	278	16430	0.87	ng/ul	96
26) Benzo(g,h,i)perylene	26.50	276	17696	0.86	ng/ul	98

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

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