

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM031215\
 Data File : BM000700.D
 Acq On : 12 Mar 2015 15:41
 Operator : TP/IZ
 Sample : SSTD0.874
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.874

Quant Time: Mar 13 09:02:17 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM021815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 13 08:51:18 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	3569	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	12829	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	1753	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	16489	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	16413	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	15706	0.40	ng/ul	0.01

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	13841	0.79	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	31223	0.86	ng/ul	0.00

Target Compounds						Qvalue
3) Naphthalene	10.48	128	27735	0.83	ng/ul	100
5) 2-Methylnaphthalene	12.10	142	19036	0.81	ng/ul	100
7) Acenaphthylene	14.02	152	26784	2.25	ng/ul	99
8) Acenaphthene	14.36	153	20618	2.25	ng/ul	98
9) Fluorene	15.35	166	25012	2.22	ng/ul	99
11) Pentachlorophenol	16.71	266	1634	0.38	ng/ul	98
12) Phenanthrene	17.08	178	40147	0.83	ng/ul	99
13) Anthracene	17.17	178	36767	0.81	ng/ul	100
15) Fluoranthene	19.11	202	47612	0.86	ng/ul	100
17) Pyrene	19.48	202	48667	0.66	ng/ul	99
18) Benzo(a)anthracene	21.23	228	42060	0.80	ng/ul	99
19) Chrysene	21.28	228	48678	0.84	ng/ul	99
21) Benzo(b)fluoranthene	22.78	252	48100	0.80	ng/ul	97
22) Benzo(k)fluoranthene	22.82	252	46940	0.79	ng/ul	98
23) Benzo(a)pyrene	23.34	252	44365	0.82	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.65	276	51484	0.99	ng/ul	100
25) Dibenzo(a,h)anthracene	25.66	278	41315	1.00	ng/ul	98
26) Benzo(g,h,i)perylene	26.31	276	45094	0.99	ng/ul	99

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

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