

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM021215\
 Data File : BM000503.D
 Acq On : 12 Feb 2015 15:19
 Operator : TP/IZ
 Sample : SSTD0.801
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.801

Quant Time: Feb 12 16:34:50 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	3568	0.40	ng/ul	0.00
2) Naphthalene-d8	10.49	136	11957	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.35	164	7397	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.09	188	14646	0.40	ng/ul	0.00
16) Chrysene-d12	21.29	240	11537	0.40	ng/ul	0.00
20) Perylene-d12	23.51	264	9700	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.09	152	12552	0.79	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	24095	0.74	ng/ul	0.00

Target Compounds						Qvalue
3) Naphthalene	10.54	128	22785	0.78	ng/ul	99
5) 2-Methylnaphthalene	12.16	142	15687	0.77	ng/ul	99
7) Acenaphthylene	14.07	152	22982	0.78	ng/ul	99
8) Acenaphthene	14.41	153	17631	0.77	ng/ul	99
9) Fluorene	15.40	166	20792	0.75	ng/ul	98
11) Pentachlorophenol	16.75	266	1581	1.08	ng/ul	98
12) Phenanthrene	17.13	178	32147	0.78	ng/ul	99
13) Anthracene	17.23	178	29597	0.80	ng/ul	99
15) Fluoranthene	19.16	202	35382	0.74	ng/ul	99
17) Pyrene	19.53	202	35376	0.84	ng/ul	99
18) Benzo(a)anthracene	21.27	228	26177	0.77	ng/ul	99
19) Chrysene	21.32	228	30610	0.78	ng/ul	100
21) Benzo(b)fluoranthene	22.84	252	27311	0.82	ng/ul	97
22) Benzo(k)fluoranthene	22.89	252	26102	0.75	ng/ul	97
23) Benzo(a)pyrene	23.42	252	24727	0.79	ng/ul	94
24) Indeno(1,2,3-cd)pyrene	25.75	276	27270	0.81	ng/ul	100
25) Dibenzo(a,h)anthracene	25.77	278	21774	0.82	ng/ul	95
26) Benzo(g,h,i)perylene	26.44	276	23840	0.81	ng/ul	99

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

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