

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM071415\
 Data File : BM001908.D
 Acq On : 13 Jul 2015 22:10
 Operator : TP/UM
 Sample : SSTD0.845
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.845

Quant Time: Jul 14 04:21:10 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM071415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 14 04:19:54 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	1213	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	4658	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	2471	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	5259	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	4772	0.40	ng/ul	0.00
20) Perylene-d12	23.47	264	4162	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.04	152	5814	0.96	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	11633	0.98	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.49	128	11222	0.88	ng/ul	99
5) 2-Methylnaphthalene	12.11	142	7531	0.93	ng/ul	99
7) Acenaphthylene	14.01	152	11567	0.93	ng/ul	99
8) Acenaphthene	14.37	153	8153	0.87	ng/ul	99
9) Fluorene	15.37	166	9390	0.89	ng/ul	97
11) Pentachlorophenol	16.70	266	1080	0.92	ng/ul	99
12) Phenanthrene	17.09	178	14328	0.87	ng/ul	99
13) Anthracene	17.18	178	14228	0.94	ng/ul	99
15) Fluoranthene	19.12	202	16368	0.93	ng/ul	98
17) Pyrene	19.48	202	16732	0.86	ng/ul	98
18) Benzo(a)anthracene	21.23	228	14453	0.90	ng/ul	99
19) Chrysene	21.28	228	14153	0.84	ng/ul	97
21) Benzo(b)fluoranthene	22.80	252	15158	0.91	ng/ul	96
22) Benzo(k)fluoranthene	22.85	252	12022	0.78	ng/ul	98
23) Benzo(a)pyrene	23.37	252	12844	0.87	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.71	276	13346	0.79	ng/ul	100
25) Dibenzo(a,h)anthracene	25.73	278	10618	0.77	ng/ul	98
26) Benzo(g,h,i)perylene	26.40	276	11328	0.77	ng/ul	98

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

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