

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM040215\
 Data File : BM000806.D
 Acq On : 02 Apr 2015 10:22
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC001

Quant Time: Apr 03 02:07:17 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIMPAH-BM033115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 01 04:17:30 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	32988	5.00	ng	0.00
4) Naphthalene-d8	10.40	136	115766	5.00	ng	0.00
8) Acenaphthene-d10	14.26	164	56967	5.00	ng	-0.02
13) Phenanthrene-d10	17.01	188	128986	5.00	ng	0.00
19) Chrysene-d12	21.21	240	104195	5.00	ng	0.00
25) Perylene-d12	23.40	264	97125	5.00	ng	0.00
System Monitoring Compounds						
5) Nitrobenzene-d5	8.77	82	5815	0.99	ng	0.00
9) 2-Fluorobiphenyl	12.89	172	18412	1.04	ng	0.00
21) Terphenyl-d14	19.66	244	13069	1.04	ng	0.00
Target Compounds						Qvalue
3) n-Nitrosodimethylamine	3.44	42	3184	0.99	ng	# 99
6) Naphthalene	10.45	128	27778	1.03	ng	98
7) 2-Methylnaphthalene	12.07	142	17607	1.04	ng	94
10) Acenaphthylene	13.99	152	26143	1.04	ng	100
11) Acenaphthene	14.33	154	17482	1.00	ng	100
12) Fluorene	15.32	166	20782	1.06	ng	100
14) Hexachlorobenzene	16.33	284	7402	0.98	ng	# 96
15) Pentachlorophenol	16.68	266	1177	0.95	ng	96
16) Phenanthrene	17.05	178	34259	0.97	ng	99
17) Anthracene	17.13	178	28116m	0.95	ng	
18) Fluoranthene	19.08	202	33404	0.96	ng	100
20) Pyrene	19.44	202	35216	1.04	ng	99
22) Benzo(a)anthracene	21.19	228	30933	1.04	ng	# 91
23) Chrysene	21.24	228	32338	1.04	ng	97
24) Indeno(1,2,3-cd)pyrene	25.58	276	33637	1.04	ng	100
26) Benzo(b)fluoranthene	22.73	252	33402	1.09	ng	96
27) Benzo(k)fluoranthene	22.79	252	27818	0.95	ng	98
28) Benzo(a)pyrene	23.30	252	27542	1.05	ng	97
29) Dibenzo(a,h)anthracene	25.59	278	26340	1.02	ng	99
30) Benzo(g,h,i)perylene	26.25	276	28228	1.01	ng	100

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

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