

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM040815\
 Data File : BM000902.D
 Acq On : 08 Apr 2015 10:33
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC001

Quant Time: Apr 09 03:33:30 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	29284	5.00	ng	-0.01
4) Naphthalene-d8	10.39	136	101700	5.00	ng	-0.01
8) Acenaphthene-d10	14.26	164	52130	5.00	ng	-0.02
13) Phenanthrene-d10	17.01	188	109283	5.00	ng	0.00
19) Chrysene-d12	21.21	240	94676	5.00	ng	0.00
25) Perylene-d12	23.39	264	85620	5.00	ng	-0.01

System Monitoring Compounds						
5) Nitrobenzene-d5	8.76	82	4735	0.93	ng	-0.01
9) 2-Fluorobiphenyl	12.88	172	15797	0.98	ng	-0.01
21) Terphenyl-d14	19.65	244	11825	1.04	ng	-0.01

Target Compounds				Qvalue		
3) n-Nitrosodimethylamine	3.44	42	2727	0.96	ng	95
6) Naphthalene	10.44	128	23312	0.99	ng	100
7) 2-Methylnaphthalene	12.06	142	15176	1.02	ng	98
10) Acenaphthylene	13.97	152	22047	0.96	ng	100
11) Acenaphthene	14.33	154	15183	0.95	ng	99
12) Fluorene	15.32	166	17788	0.99	ng	99
14) Hexachlorobenzene	16.33	284	5925	0.93	ng	# 43
15) Pentachlorophenol	16.66	266	1145	1.03	ng	88
16) Phenanthrene	17.05	178	27874	0.93	ng	99
17) Anthracene	17.13	178	26459	1.05	ng	99
18) Fluoranthene	19.08	202	29770	1.01	ng	99
20) Pyrene	19.44	202	31241	1.01	ng	100
22) Benzo(a)anthracene	21.19	228	26053	0.96	ng	96
23) Chrysene	21.24	228	28403	1.00	ng	91
24) Indeno(1,2,3-cd)pyrene	25.57	276	27621	0.94	ng	98
26) Benzo(b)fluoranthene	22.73	252	26551	0.98	ng	99
27) Benzo(k)fluoranthene	22.78	252	25553	0.99	ng	99
28) Benzo(a)pyrene	23.29	252	22535	0.97	ng	96
29) Dibenzo(a,h)anthracene	25.58	278	20980	0.92	ng	99
30) Benzo(g,h,i)perylene	26.23	276	24007	0.97	ng	99

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

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