

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM033115\
 Data File : BM000775.D
 Acq On : 01 Apr 2015 04:22
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC001

Quant Time: Apr 01 15:09:42 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	7.62	152	30279	5.00	ng	0.00
4) Naphthalene-d8	10.40	136	104396	5.00	ng	0.00
8) Acenaphthene-d10	14.28	164	51135	5.00	ng	0.00
13) Phenanthrene-d10	17.01	188	109087	5.00	ng	0.00
19) Chrysene-d12	21.21	240	91607	5.00	ng	0.00
25) Perylene-d12	23.40	264	83495	5.00	ng	0.00
System Monitoring Compounds						
5) Nitrobenzene-d5	8.77	82	5014	0.96	ng	0.00
9) 2-Fluorobiphenyl	12.89	172	16378	1.03	ng	0.00
21) Terphenyl-d14	19.66	244	11286	1.02	ng	0.00
Target Compounds						Qvalue
3) n-Nitrosodimethylamine	3.45	42	2859	0.97	ng	# 99
6) Naphthalene	10.45	128	24926	1.03	ng	99
7) 2-Methylnaphthalene	12.07	142	15619	1.02	ng	99
10) Acenaphthylene	13.99	152	22312	0.99	ng	100
11) Acenaphthene	14.33	154	15532	0.99	ng	100
12) Fluorene	15.32	166	18001	1.03	ng	99
14) Hexachlorobenzene	16.33	284	6437	1.01	ng	# 97
15) Pentachlorophenol	16.68	266	666	0.76	ng	99
16) Phenanthrene	17.05	178	29995	1.00	ng	100
17) Anthracene	17.15	178	23924	0.96	ng	99
18) Fluoranthene	19.08	202	28637	0.97	ng	100
20) Pyrene	19.44	202	30065	1.01	ng	100
22) Benzo(a)anthracene	21.19	228	25420	0.97	ng	# 89
23) Chrysene	21.24	228	28302	1.03	ng	100
24) Indeno(1,2,3-cd)pyrene	25.58	276	29158	1.03	ng	100
26) Benzo(b)fluoranthene	22.74	252	26620	1.01	ng	100
27) Benzo(k)fluoranthene	22.79	252	25690	1.02	ng	100
28) Benzo(a)pyrene	23.30	252	22369	0.99	ng	96
29) Dibenzo(a,h)anthracene	25.60	278	22889	1.03	ng	100
30) Benzo(g,h,i)perylene	26.26	276	24816	1.03	ng	100

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

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