

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM033115\
 Data File : BM000777.D
 Acq On : 01 Apr 2015 06:09
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC001

Quant Time: Apr 01 15:09:56 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	30397	5.00	ng	0.00
4) Naphthalene-d8	10.40	136	104003	5.00	ng	0.00
8) Acenaphthene-d10	14.28	164	50801	5.00	ng	0.00
13) Phenanthrene-d10	17.01	188	108479	5.00	ng	0.00
19) Chrysene-d12	21.22	240	92417	5.00	ng	0.01
25) Perylene-d12	23.40	264	87075	5.00	ng	0.00
System Monitoring Compounds						
5) Nitrobenzene-d5	8.77	82	5188	0.99	ng	0.00
9) 2-Fluorobiphenyl	12.89	172	16350	1.04	ng	0.00
21) Terphenyl-d14	19.66	244	11421	1.03	ng	0.00
Target Compounds						Qvalue
3) n-Nitrosodimethylamine	3.45	42	2800	0.95	ng	98
6) Naphthalene	10.45	128	24921	1.03	ng	100
7) 2-Methylnaphthalene	12.07	142	15646	1.02	ng	98
10) Acenaphthylene	13.98	152	22599	1.01	ng	100
11) Acenaphthene	14.33	154	15760	1.01	ng	99
12) Fluorene	15.32	166	18105	1.04	ng	100
14) Hexachlorobenzene	16.33	284	6607	1.04	ng	# 96
15) Pentachlorophenol	16.68	266	833	0.86	ng	97
16) Phenanthrene	17.05	178	30452	1.02	ng	100
17) Anthracene	17.15	178	23781	0.95	ng	99
18) Fluoranthene	19.08	202	29349	1.00	ng	100
20) Pyrene	19.44	202	30556	1.01	ng	100
22) Benzo(a)anthracene	21.20	228	26266	1.00	ng	99
23) Chrysene	21.25	228	29031	1.05	ng	# 88
24) Indeno(1,2,3-cd)pyrene	25.59	276	30275	1.06	ng	99
26) Benzo(b)fluoranthene	22.74	252	27052	0.99	ng	99
27) Benzo(k)fluoranthene	22.79	252	27179	1.03	ng	99
28) Benzo(a)pyrene	23.31	252	23867	1.01	ng	99
29) Dibenzo(a,h)anthracene	25.60	278	23688	1.02	ng	100
30) Benzo(g,h,i)perylene	26.26	276	25604	1.02	ng	100

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

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