

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM040215\
 Data File : BM000809.D
 Acq On : 02 Apr 2015 14:12
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC001

Quant Time: Apr 03 02:12:44 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	33557	5.00	ng	0.00
4) Naphthalene-d8	10.40	136	116558	5.00	ng	0.00
8) Acenaphthene-d10	14.28	164	57285	5.00	ng	0.00
13) Phenanthrene-d10	17.01	188	124654	5.00	ng	0.00
19) Chrysene-d12	21.21	240	103504	5.00	ng	0.00
25) Perylene-d12	23.40	264	93845	5.00	ng	0.00
System Monitoring Compounds						
5) Nitrobenzene-d5	8.77	82	5723	0.97	ng	0.00
9) 2-Fluorobiphenyl	12.89	172	18386	1.03	ng	0.00
21) Terphenyl-d14	19.66	244	12865	1.03	ng	0.00
Target Compounds						Qvalue
3) n-Nitrosodimethylamine	3.44	42	3147	0.97	ng	99
6) Naphthalene	10.45	128	27781	1.03	ng	99
7) 2-Methylnaphthalene	12.07	142	17598	1.03	ng	94
10) Acenaphthylene	13.99	152	25654	1.01	ng	100
11) Acenaphthene	14.33	154	17511	1.00	ng	100
12) Fluorene	15.32	166	20383	1.04	ng	100
14) Hexachlorobenzene	16.33	284	7086	0.97	ng	# 98
15) Pentachlorophenol	16.68	266	806	0.78	ng	97
16) Phenanthrene	17.05	178	34121	1.00	ng	99
17) Anthracene	17.13	178	27441m	0.96	ng	
18) Fluoranthene	19.08	202	32458	0.96	ng	100
20) Pyrene	19.44	202	33867	1.00	ng	100
22) Benzo(a)anthracene	21.19	228	29128	0.99	ng	# 88
23) Chrysene	21.24	228	31596	1.02	ng	99
24) Indeno(1,2,3-cd)pyrene	25.58	276	32654	1.02	ng	99
26) Benzo(b)fluoranthene	22.74	252	30619	1.04	ng	100
27) Benzo(k)fluoranthene	22.79	252	28238	1.00	ng	99
28) Benzo(a)pyrene	23.30	252	25556	1.01	ng	96
29) Dibenzo(a,h)anthracene	25.60	278	25658	1.02	ng	98
30) Benzo(g,h,i)perylene	26.25	276	27735	1.03	ng	98

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

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