

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

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	5.000	5.000	0.0	104	0.00
2	1,4-Dioxane	1.000	0.000	100.0#	0	-2.47#
3	n-Nitrosodimethylamine	1.000	0.966	3.4	101	0.00
4 I	Naphthalene-d8	5.000	5.000	0.0	103	0.00
5 S	Nitrobenzene-d5	1.000	0.975	2.5	104	0.00
6	Naphthalene	1.000	1.026	-2.6	104	0.00
7	2-Methylnaphthalene	1.000	1.028	-2.8	104	0.00
8 I	Acenaphthene-d10	5.000	5.000	0.0	104	0.00
9 S	2-Fluorobiphenyl	1.000	1.035	-3.5	105	0.00
10	Acenaphthylene	1.000	1.012	-1.2	106	0.00
11	Acenaphthene	1.000	0.997	0.3	105	0.00
12	Fluorene	1.000	1.037	-3.7	107	0.00
13 I	Phenanthrene-d10	5.000	5.000	0.0	111	0.00
14	Hexachlorobenzene	1.000	0.970	3.0	105	0.00
15	Pentachlorophenol	1.000	0.782	21.8	78	0.00
16	Phenanthrene	1.000	0.999	0.1	109	0.00
17	Anthracene	1.000	0.959	4.1	105	-0.02
18	Fluoranthene	1.000	0.962	3.8	107	0.00
19 I	Chrysene-d12	5.000	5.000	0.0	108	0.00
20	Pyrene	1.000	1.003	-0.3	107	0.00
21 S	Terphenyl-d14	1.000	1.033	-3.3	109	0.00
22	Benzo(a)anthracene	1.000	0.985	1.5	110	-0.01
23	Chrysene	1.000	1.020	-2.0	108	0.00
24	Indeno(1,2,3-cd)pyrene	1.000	1.018	-1.8	111	-0.01
25 I	Perylene-d12	5.000	5.000	0.0	110	0.00
26	Benzo(b)fluoranthene	1.000	1.036	-3.6	113	0.00
27	Benzo(k)fluoranthene	1.000	0.996	0.4	107	0.00
28 C	Benzo(a)pyrene	1.000	1.007	-0.7	111	-0.01
29	Dibenzo(a,h)anthracene	1.000	1.024	-2.4	111	0.00
30	Benzo(g,h,i)perylene	1.000	1.025	-2.5	111	-0.01

(#) = Out of Range

SPCC's out = 0 CCC's out = 0