

Data Path : Z:\HPCHEM1\BNA M\Data\BM042215\
 Data File : BM001107.D
 Acq On : 22 Apr 2015 14:19
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC001

Quant Time: Apr 23 11:10:30 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIMPAH-BM041815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 20 08:05:22 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	140	-0.01
2	1,4-Dioxane	1.000	0.925	7.5	127	-0.01
3	n-Nitrosodimethylamine	1.000	1.213	-21.3	172	-0.01
4 I	Naphthalene-d8	5.000	5.000	0.0	138	-0.01
5 S	Nitrobenzene-d5	1.000	1.217	-21.7	167	0.00
6	Naphthalene	1.000	0.975	2.5	136	-0.01
7	2-Methylnaphthalene	1.000	0.967	3.3	133	-0.01
8 I	Acenaphthene-d10	5.000	5.000	0.0	126	-0.02
9 S	2-Fluorobiphenyl	1.000	1.000	0.0	126	-0.01
10	Acenaphthylene	1.000	1.093	-9.3	136	-0.02
11	Acenaphthene	1.000	0.955	4.5	124	0.00
12	Fluorene	1.000	0.989	1.1	124	-0.02
13 I	Phenanthrene-d10	5.000	5.000	0.0	115	-0.02
14	Hexachlorobenzene	1.000	1.031	-3.1	121	-0.02
15	Pentachlorophenol	1.000	1.340	-34.0#	182	0.00
16	Phenanthrene	1.000	1.021	-2.1	120	-0.02
17	Anthracene	1.000	1.030	-3.0	117	0.00
18	Fluoranthene	1.000	1.013	-1.3	116	-0.01
19 I	Chrysene-d12	5.000	5.000	0.0	112	-0.01
20	Pvrene	1.000	1.036	-3.6	118	0.00
21 S	Terphenyl-d14	1.000	1.008	-0.8	113	-0.01
22	Benzo(a)anthracene	1.000	1.055	-5.5	122	0.00
23	Chrysene	1.000	0.959	4.1	109	0.00
24	Indeno(1,2,3-cd)pyrene	1.000	1.144	-14.4	130	0.00
25 I	Perylene-d12	5.000	5.000	0.0	120	-0.01
26	Benzo(b)fluoranthene	1.000	1.007	-0.7	121	0.00
27	Benzo(k)fluoranthene	1.000	0.982	1.8	118	0.00
28 C	Benzo(a)pyrene	1.000	1.067	-6.7	128	0.00
29	Dibenzo(a,h)anthracene	1.000	1.061	-6.1	127	0.00
30	Benzo(g,h,i)perylene	1.000	1.058	-5.8	129	0.00

(#) = Out of Range

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