

Data Path : Z:\HPCHEM1\BNA M\Data\BM042215\
 Data File : BM001110.D
 Acq On : 22 Apr 2015 18:10
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC001

Quant Time: Apr 23 03:45:54 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	124	0.00
2	1,4-Dioxane	1.000	0.886	11.4	108	0.00
3	n-Nitrosodimethylamine	1.000	1.234	-23.4	154	-0.01
4 I	Naphthalene-d8	5.000	5.000	0.0	117	-0.01
5 S	Nitrobenzene-d5	1.000	1.203	-20.3	140	0.00
6	Naphthalene	1.000	0.992	0.8	118	-0.01
7	2-Methylnaphthalene	1.000	0.953	4.7	111	-0.01
8 I	Acenaphthene-d10	5.000	5.000	0.0	104	-0.02
9 S	2-Fluorobiphenyl	1.000	1.003	-0.3	105	-0.01
10	Acenaphthylene	1.000	1.082	-8.2	112	-0.02
11	Acenaphthene	1.000	0.964	3.6	104	0.00
12	Fluorene	1.000	0.980	2.0	102	-0.02
13 I	Phenanthrene-d10	5.000	5.000	0.0	95	-0.02
14	Hexachlorobenzene	1.000	1.054	-5.4	103	-0.02
15	Pentachlorophenol	1.000	1.253	-25.3#	139	0.00
16	Phenanthrene	1.000	1.049	-4.9	102	-0.02
17	Anthracene	1.000	1.028	-2.8	97	0.00
18	Fluoranthene	1.000	1.062	-6.2	101	-0.01
19 I	Chrysene-d12	5.000	5.000	0.0	105	-0.01
20	Pvrene	1.000	0.984	1.6	105	-0.01
21 S	Terphenyl-d14	1.000	0.941	5.9	99	-0.01
22	Benzo(a)anthracene	1.000	1.053	-5.3	114	0.00
23	Chrysene	1.000	0.970	3.0	103	0.00
24	Indeno(1,2,3-cd)pyrene	1.000	1.162	-16.2	123	-0.01
25 I	Perylene-d12	5.000	5.000	0.0	114	-0.01
26	Benzo(b)fluoranthene	1.000	1.066	-6.6	122	0.00
27	Benzo(k)fluoranthene	1.000	0.923	7.7	105	0.00
28 C	Benzo(a)pyrene	1.000	1.064	-6.4	121	-0.01
29	Dibenzo(a,h)anthracene	1.000	1.070	-7.0	122	0.00
30	Benzo(g,h,i)perylene	1.000	1.057	-5.7	122	0.00

(#) = Out of Range

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