

Data Path : Z:\HPCHEM1\BNA E\Data\BE080715\
 Data File : BE090394.D
 Acq On : 7 Aug 2015 13:00
 Operator : TP/UM
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC001

Quant Time: Aug 07 23:08:32 2015
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE080515.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 06 12:28:08 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	1.000	0.0	93	0.00
3	n-Nitrosodimethylamine	1.000	0.985	1.5	103	0.00
4 S	2-Fluorophenol	1.000	1.021	-2.1	114	0.00
5 S	Phenol-d6	1.000	0.980	2.0	111	0.00
6 I	Naphthalene-d8	5.000	5.000	0.0	105	0.00
7 S	Nitrobenzene-d5	1.000	0.944	5.6	106	0.00
8	Nitrobenzene	1.000	0.909	9.1	107	0.00
9	Naphthalene	1.000	1.001	-0.1	106	0.00
10	Hexachlorobutadiene	1.000	1.123	-12.3	108	0.00
11	2-Methylnaphthalene	1.000	0.982	1.8	108	0.00
12 I	Acenaphthene-d10	5.000	5.000	0.0	110	0.00
13 S	2,4,6-Tribromophenol	1.000	0.938	6.2	113	0.00
14 S	2-Fluorobiphenyl	1.000	1.006	-0.6	109	0.00
15	Acenaphthylene	1.000	0.969	3.1	108	0.00
16	Acenaphthene	1.000	0.988	1.2	109	0.00
17	Fluorene	1.000	1.033	-3.3	112	0.00
18 I	Phenanthrene-d10	5.000	5.000	0.0	111	0.00
19	4-Bromophenyl-phenylether	1.000	0.945	5.5	107	0.00
20	Hexachlorobenzene	1.000	0.952	4.8	105	0.00
21	Pentachlorophenol	1.000	1.009	-0.9	133	0.00
22	Phenanthrene	1.000	0.975	2.5	109	0.00
23	Anthracene	1.000	0.973	2.7	111	-0.02
24	Fluoranthene	1.000	0.953	4.7	109	0.00
25 I	Chrysene-d12	5.000	5.000	0.0	113	0.00
26	Pyrene	1.000	0.960	4.0	113	0.00
27 S	Terphenyl-d14	1.000	0.967	3.3	111	0.00
28	Benzo(a)anthracene	1.000	0.942	5.8	119	0.00
29	Chrysene	1.000	0.975	2.5	108	0.00
30	Bis(2-ethylhexyl)phthalate	1.000	0.814	18.6	100	0.00
31	Indeno(1,2,3-cd)pyrene	1.000	0.877	12.3	116	0.00
32 I	Perylene-d12	5.000	5.000	0.0	111	0.00
33	Benzo(b)fluoranthene	1.000	1.001	-0.1	121	0.00
34	Benzo(k)fluoranthene	1.000	0.936	6.4	101	0.00
35 C	Benzo(a)pyrene	1.000	0.934	6.6	110	0.00
36	Dibenzo(a,h)anthracene	1.000	0.915	8.5	117	0.00
37	Benzo(g,h,i)perylene	1.000	0.924	7.6	106	0.00

(#) = Out of Range

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