

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE091015\
 Data File : BE090781.D
 Acq On : 11 Sep 2015 3:13
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC001

Quant Time: Sep 11 03:55:46 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE090115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 02 11:40:50 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	110	0.00
2	1,4-Dioxane	1.000	0.788	21.2	85	0.00
3	n-Nitrosodimethylamine	1.000	0.779	22.1	85	0.00
4 S	2-Fluorophenol	1.000	0.868	13.2	94	0.00
5 S	Phenol-d6	1.000	0.897	10.3	104	0.00
6 I	Naphthalene-d8	5.000	5.000	0.0	121	0.00
7 S	Nitrobenzene-d5	1.000	0.830	17.0	101	0.00
8	Nitrobenzene	1.000	0.845	15.5	106	0.00
9	Naphthalene	1.000	0.860	14.0	105	0.00
10	Hexachlorobutadiene	1.000	0.842	15.8	99	0.00
11	2-Methylnaphthalene	1.000	0.905	9.5	113	0.00
12 I	Acenaphthene-d10	5.000	5.000	0.0	132	0.00
13 S	2,4,6-Tribromophenol	1.000	0.864	13.6	116	0.00
14 S	2-Fluorobiphenyl	1.000	0.823	17.7	111	0.00
15	Acenaphthylene	1.000	0.850	15.0	117	0.00
16	Acenaphthene	1.000	0.847	15.3	111	0.00
17	Fluorene	1.000	0.855	14.5	114	0.00
18 I	Phenanthrene-d10	5.000	5.000	0.0	127	0.00
19	4-Bromophenyl-phenylether	1.000	0.837	16.3	111	0.00
20	Hexachlorobenzene	1.000	0.834	16.6	106	0.00
21	Pentachlorophenol	1.000	0.956	4.4	99	0.00
22	Phenanthrene	1.000	0.854	14.6	109	-0.02
23	Anthracene	1.000	0.877	12.3	116	0.00
24	Fluoranthene	1.000	0.825	17.5	114	0.00
25 I	Chrysene-d12	5.000	5.000	0.0	121	0.00
26	Pyrene	1.000	0.917	8.3	117	0.00
27 S	Terphenyl-d14	1.000	0.891	10.9	113	0.00
28	Benzo(a)anthracene	1.000	0.887	11.3	109	0.00
29	Chrysene	1.000	0.903	9.7	107	0.00
30	Bis(2-ethylhexyl)phthalate	1.000	0.814	18.6	108	0.00
31	Indeno(1,2,3-cd)pyrene	1.000	0.873	12.7	115	0.00
32 I	Perylene-d12	5.000	5.000	0.0	125	0.00
33	Benzo(b)fluoranthene	1.000	0.846	15.4	109	0.00
34	Benzo(k)fluoranthene	1.000	0.884	11.6	112	0.00
35 C	Benzo(a)pyrene	1.000	0.873	12.7	117	0.00
36	Dibenzo(a,h)anthracene	1.000	0.871	12.9	114	0.00
37	Benzo(g,h,i)perylene	1.000	0.829	17.1	110	-0.01

(#) = Out of Range

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