

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE091015\
 Data File : BE090772.D
 Acq On : 10 Sep 2015 21:45
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC001

Quant Time: Sep 11 03:42:55 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	0.824	17.6	88	0.00
3	n-Nitrosodimethylamine	1.000	0.767	23.3	83	0.00
4 S	2-Fluorophenol	1.000	0.784	21.6	84	0.00
5 S	Phenol-d6	1.000	0.874	12.6	101	0.00
6 I	Naphthalene-d8	5.000	5.000	0.0	123	0.00
7 S	Nitrobenzene-d5	1.000	0.798	20.2	99	0.00
8	Nitrobenzene	1.000	0.818	18.2	104	0.00
9	Naphthalene	1.000	0.855	14.5	106	0.00
10	Hexachlorobutadiene	1.000	0.832	16.8	99	0.00
11	2-Methylnaphthalene	1.000	0.903	9.7	115	0.00
12 I	Acenaphthene-d10	5.000	5.000	0.0	139	0.00
13 S	2,4,6-Tribromophenol	1.000	0.807	19.3	112	0.00
14 S	2-Fluorobiphenyl	1.000	0.816	18.4	115	0.00
15	Acenaphthylene	1.000	0.833	16.7	121	0.00
16	Acenaphthene	1.000	0.857	14.3	118	0.00
17	Fluorene	1.000	0.846	15.4	119	0.00
18 I	Phenanthrene-d10	5.000	5.000	0.0	134	0.00
19	4-Bromophenyl-phenylether	1.000	0.834	16.6	116	0.00
20	Hexachlorobenzene	1.000	0.824	17.6	110	0.00
21	Pentachlorophenol	1.000	0.867	13.3	85	0.00
22	Phenanthrene	1.000	0.856	14.4	115	-0.02
23	Anthracene	1.000	0.864	13.6	120	0.00
24	Fluoranthene	1.000	0.778	22.2	112	0.00
25 I	Chrysene-d12	5.000	5.000	0.0	124	0.00
26	Pyrene	1.000	0.896	10.4	117	0.00
27 S	Terphenyl-d14	1.000	0.865	13.5	112	0.00
28	Benzo(a)anthracene	1.000	0.861	13.9	108	0.00
29	Chrysene	1.000	0.901	9.9	109	0.00
30	Bis(2-ethylhexyl)phthalate	1.000	0.718	28.2#	96	0.00
31	Indeno(1,2,3-cd)pyrene	1.000	0.849	15.1	114	0.00
32 I	Perylene-d12	5.000	5.000	0.0	124	0.00
33	Benzo(b)fluoranthene	1.000	0.852	14.8	109	0.00
34	Benzo(k)fluoranthene	1.000	0.862	13.8	109	0.00
35 C	Benzo(a)pyrene	1.000	0.847	15.3	113	0.00
36	Dibenzo(a,h)anthracene	1.000	0.860	14.0	111	-0.01
37	Benzo(g,h,i)perylene	1.000	0.806	19.4	106	0.00

(#) = Out of Range

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