

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081715\
 Data File : BE090551.D
 Acq On : 18 Aug 2015 2:11
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC001

Quant Time: Aug 18 06:12:31 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE081415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 14 16:52:42 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	115	0.00
2	1,4-Dioxane	1.000	1.061	-6.1	105	0.00
3	n-Nitrosodimethylamine	1.000	0.966	3.4	112	0.00
4 S	2-Fluorophenol	1.000	1.055	-5.5	120	0.00
5 S	Phenol-d6	1.000	1.045	-4.5	125	0.00
6 I	Naphthalene-d8	5.000	5.000	0.0	117	0.00
7 S	Nitrobenzene-d5	1.000	0.991	0.9	122	0.00
8	Nitrobenzene	1.000	0.986	1.4	124	0.00
9	Naphthalene	1.000	0.977	2.3	115	0.00
10	Hexachlorobutadiene	1.000	0.984	1.6	116	0.00
11	2-Methylnaphthalene	1.000	1.008	-0.8	121	0.00
12 I	Acenaphthene-d10	5.000	5.000	0.0	120	0.00
13 S	2,4,6-Tribromophenol	1.000	1.048	-4.8	138	0.00
14 S	2-Fluorobiphenyl	1.000	1.006	-0.6	121	0.00
15	Acenaphthylene	1.000	1.005	-0.5	122	0.00
16	Acenaphthene	1.000	0.995	0.5	121	0.00
17	Fluorene	1.000	0.967	3.3	116	0.02
18 I	Phenanthrene-d10	5.000	5.000	0.0	114	0.00
19	4-Bromophenyl-phenylether	1.000	1.051	-5.1	123	0.00
20	Hexachlorobenzene	1.000	1.036	-3.6	119	0.00
21	Pentachlorophenol	1.000	0.960	4.0	146	0.00
22	Phenanthrene	1.000	0.995	0.5	116	0.00
23	Anthracene	1.000	1.022	-2.2	121	0.00
24	Fluoranthene	1.000	0.999	0.1	117	0.00
25 I	Chrysene-d12	5.000	5.000	0.0	115	0.00
26	Pyrene	1.000	1.007	-0.7	119	0.00
27 S	Terphenyl-d14	1.000	1.012	-1.2	118	0.00
28	Benzo(a)anthracene	1.000	0.977	2.3	118	0.00
29	Chrysene	1.000	0.993	0.7	116	0.00
30	Bis(2-ethylhexyl)phthalate	1.000	1.122	-12.2	153	0.00
31	Indeno(1,2,3-cd)pyrene	1.000	0.990	1.0	124	0.00
32 I	Perylene-d12	5.000	5.000	0.0	118	0.00
33	Benzo(b)fluoranthene	1.000	1.003	-0.3	122	0.00
34	Benzo(k)fluoranthene	1.000	1.020	-2.0	120	0.00
35 C	Benzo(a)pyrene	1.000	1.018	-1.8	124	0.00
36	Dibenzo(a,h)anthracene	1.000	0.994	0.6	124	0.00
37	Benzo(g,h,i)perylene	1.000	1.018	-1.8	123	0.00

(#) = Out of Range

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