

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081215\
 Data File : BE090480.D
 Acq On : 12 Aug 2015 22:57
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC001

Quant Time: Aug 13 01:22:31 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	114	0.00
2	1,4-Dioxane	1.000	1.030	-3.0	104	0.00
3	n-Nitrosodimethylamine	1.000	1.055	-5.5	121	0.00
4 S	2-Fluorophenol	1.000	1.419	-41.9#	173	0.00
5 S	Phenol-d6	1.000	1.238	-23.8	164	0.00
6 I	Naphthalene-d8	5.000	5.000	0.0	123	0.00
7 S	Nitrobenzene-d5	1.000	1.009	-0.9	134	0.00
8	Nitrobenzene	1.000	1.007	-0.7	142	0.00
9	Naphthalene	1.000	1.016	-1.6	125	0.00
10	Hexachlorobutadiene	1.000	1.078	-7.8	121	0.00
11	2-Methylnaphthalene	1.000	1.078	-7.8	138	0.00
12 I	Acenaphthene-d10	5.000	5.000	0.0	133	0.00
13 S	2,4,6-Tribromophenol	1.000	1.471	-47.1#	236	0.00
14 S	2-Fluorobiphenyl	1.000	1.027	-2.7	135	0.00
15	Acenaphthylene	1.000	1.026	-2.6	138	0.00
16	Acenaphthene	1.000	1.001	-0.1	133	0.00
17	Fluorene	1.000	1.030	-3.0	136	0.00
18 I	Phenanthrene-d10	5.000	5.000	0.0	132	0.00
19	4-Bromophenyl-phenylether	1.000	0.972	2.8	131	0.00
20	Hexachlorobenzene	1.000	0.933	6.7	123	0.00
21	Pentachlorophenol	1.000	1.319	-31.9#	224	0.00
22	Phenanthrene	1.000	0.986	1.4	131	0.00
23	Anthracene	1.000	1.073	-7.3	146	-0.02
24	Fluoranthene	1.000	1.018	-1.8	140	0.00
25 I	Chrysene-d12	5.000	5.000	0.0	125	0.00
26	Pyrene	1.000	1.092	-9.2	142	0.00
27 S	Terphenyl-d14	1.000	1.094	-9.4	138	-0.01
28	Benzo(a)anthracene	1.000	1.058	-5.8	147	0.00
29	Chrysene	1.000	0.972	2.8	119	0.00
30	Bis(2-ethylhexyl)phthalate	1.000	1.156	-15.6	170	0.00
31	Indeno(1,2,3-cd)pyrene	1.000	1.078	-7.8	169	0.00
32 I	Perylene-d12	5.000	5.000	0.0	131	0.00
33	Benzo(b)fluoranthene	1.000	1.082	-8.2	158	0.00
34	Benzo(k)fluoranthene	1.000	0.905	9.5	115	0.00
35 C	Benzo(a)pyrene	1.000	1.060	-6.0	153	0.00
36	Dibenzo(a,h)anthracene	1.000	1.081	-8.1	175	-0.01
37	Benzo(g,h,i)perylene	1.000	1.077	-7.7	146	0.00

(#) = Out of Range

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