

Data Path : Z:\HPCHEM1\BNA_E\Data\BE081815\
 Data File : BE090553.D
 Acq On : 18 Aug 2015 13:50
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC001

Quant Time: Aug 19 02:29:12 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	116	0.00
2	1,4-Dioxane	1.000	1.085	-8.5	108	0.00
3	n-Nitrosodimethylamine	1.000	0.991	0.9	116	0.00
4 S	2-Fluorophenol	1.000	1.040	-4.0	120	0.00
5 S	Phenol-d6	1.000	1.030	-3.0	124	0.00
6 I	Naphthalene-d8	5.000	5.000	0.0	118	0.00
7 S	Nitrobenzene-d5	1.000	0.980	2.0	122	0.00
8	Nitrobenzene	1.000	0.980	2.0	123	0.00
9	Naphthalene	1.000	0.975	2.5	116	0.00
10	Hexachlorobutadiene	1.000	0.978	2.2	115	0.00
11	2-Methylnaphthalene	1.000	1.003	-0.3	121	0.00
12 I	Acenaphthene-d10	5.000	5.000	0.0	120	0.00
13 S	2,4,6-Tribromophenol	1.000	1.005	-0.5	133	0.00
14 S	2-Fluorobiphenyl	1.000	1.009	-0.9	121	0.00
15	Acenaphthylene	1.000	1.003	-0.3	123	0.00
16	Acenaphthene	1.000	0.987	1.3	120	0.00
17	Fluorene	1.000	1.017	-1.7	123	0.00
18 I	Phenanthrene-d10	5.000	5.000	0.0	117	0.00
19	4-Bromophenyl-phenylether	1.000	1.027	-2.7	123	0.00
20	Hexachlorobenzene	1.000	1.002	-0.2	118	0.00
21	Pentachlorophenol	1.000	0.961	3.9	150	0.00
22	Phenanthrene	1.000	0.951	4.9	113	0.00
23	Anthracene	1.000	0.989	1.1	120	0.00
24	Fluoranthene	1.000	0.979	2.1	117	0.00
25 I	Chrysene-d12	5.000	5.000	0.0	115	0.00
26	Pyrene	1.000	1.014	-1.4	119	0.00
27 S	Terphenyl-d14	1.000	1.028	-2.8	120	0.00
28	Benzo(a)anthracene	1.000	0.983	1.7	118	0.00
29	Chrysene	1.000	0.996	0.4	115	0.00
30	Bis(2-ethylhexyl)phthalate	1.000	1.080	-8.0	146	0.00
31	Indeno(1,2,3-cd)pyrene	1.000	0.985	1.5	122	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	121	0.00
34	Benzo(k)fluoranthene	1.000	1.018	-1.8	119	0.00
35 C	Benzo(a)pyrene	1.000	1.008	-0.8	123	0.00
36	Dibenzo(a,h)anthracene	1.000	0.988	1.2	123	0.00
37	Benzo(g,h,i)perylene	1.000	1.009	-0.9	121	0.00

(#) = Out of Range

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