

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE070215\
 Data File : BE090051.D
 Acq On : 2 Jul 2015 13:01
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC001

Quant Time: Jul 03 01:26:54 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BE062915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 30 07:01:01 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	82	-0.02
2	1,4-Dioxane	1.000	0.889	11.1	76	-0.02
3	n-Nitrosodimethylamine	1.000	0.833	16.7	71	-0.02
4 S	2-Fluorophenol	1.000	0.916	8.4	75	-0.01
5 S	Phenol-d6	1.000	0.963	3.7	80	-0.01
6 I	Naphthalene-d8	5.000	5.000	0.0	87	-0.01
7 S	Nitrobenzene-d5	1.000	0.845	15.5	75	-0.01
8	Nitrobenzene	1.000	0.843	15.7	76	-0.01
9	Naphthalene	1.000	0.914	8.6	81	-0.01
10	Hexachlorobutadiene	1.000	0.932	6.8	82	-0.01
11	2-Methylnaphthalene	1.000	0.884	11.6	78	-0.01
12 I	Acenaphthene-d10	5.000	5.000	0.0	80	0.00
13 S	2,4,6-Tribromophenol	1.000	0.651	34.9#	53	0.00
14 S	2-Fluorobiphenyl	1.000	0.914	8.6	75	-0.01
15	Acenaphthylene	1.000	0.914	8.6	76	0.00
16	Acenaphthene	1.000	0.931	6.9	76	-0.01
17	Fluorene	1.000	0.889	11.1	73	-0.02
18 I	Phenanthrene-d10	5.000	5.000	0.0	72	-0.02
19	4-Bromophenyl-phenylether	1.000	0.889	11.1	66	-0.02
20	Hexachlorobenzene	1.000	0.908	9.2	68	-0.02
21	Pentachlorophenol	1.000	1.057	-5.7	83	0.00
22	Phenanthrene	1.000	0.898	10.2	67	-0.02
23	Anthracene	1.000	0.890	11.0	66	0.00
24	Fluoranthene	1.000	0.888	11.2	66	-0.01
25 I	Chrysene-d12	5.000	5.000	0.0	69	-0.01
26	Pyrene	1.000	0.943	5.7	66	-0.01
27 S	Terphenyl-d14	1.000	0.898	10.2	62	-0.01
28	Benzo(a)anthracene	1.000	0.909	9.1	64	-0.01
29	Chrysene	1.000	0.900	10.0	63	-0.01
30	Bis(2-ethylhexyl)phthalate	1.000	0.765	23.5	53	-0.01
31	Indeno(1,2,3-cd)pyrene	1.000	0.939	6.1	66	-0.03
32 I	Perylene-d12	5.000	5.000	0.0	66	-0.02
33	Benzo(b)fluoranthene	1.000	1.005	-0.5	66	-0.02
34	Benzo(k)fluoranthene	1.000	1.014	-1.4	67	-0.02
35 C	Benzo(a)pyrene	1.000	1.004	-0.4	66	-0.02
36	Dibenzo(a,h)anthracene	1.000	0.959	4.1	64	-0.03
37	Benzo(g,h,i)perylene	1.000	0.962	3.8	64	-0.04

(#) = Out of Range

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