

Data Path : Z:\HPCHEM1\BNA E\Data\BE040216\
 Data File : BE091592.D
 Acq On : 1 Apr 2016 16:05
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Apr 01 23:46:32 2016
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE033016.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 30 18:45:59 2016
 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	103	-0.02
2	1,4-Dioxane	0.400	0.325	18.8	88	-0.02
3	n-Nitrosodimethylamine	0.400	0.352	12.0	90	-0.03
4 S	2-Fluorophenol	0.400	0.368	8.0	95	-0.02
5 S	Phenol-d6	0.400	0.371	7.3	99	-0.02
6 I	Naphthalene-d8	5.000	5.000	0.0	108	0.00
7 S	Nitrobenzene-d5	0.400	0.351	12.3	97	-0.02
8	Naphthalene	0.400	0.389	2.8	102	-0.02
9	2-Methylnaphthalene	0.400	0.383	4.3	102	-0.01
10 I	Acenaphthene-d10	5.000	5.000	0.0	108	-0.03
11 S	2,4,6-Tribromophenol	0.400	0.253	36.8#	58	-0.01
12 S	2-Fluorobiphenyl	0.400	0.384	4.0	102	-0.02
13	Acenaphthylene	0.400	0.386	3.5	117	-0.01
14	Acenaphthene	0.400	0.387	3.3	106	-0.01
15	Fluorene	0.400	0.379	5.3	103	-0.01
16 I	Phenanthrene-d10	5.000	5.000	0.0	110	-0.01
17	Hexachlorobenzene	0.400	0.390	2.5	107	-0.02
18	Pentachlorophenol	0.400	0.273	31.8#	86	-0.02
19	Phenanthrene	0.400	0.378	5.5	103	-0.01
20	Anthracene	0.400	0.360	10.0	101	-0.01
21	Fluoranthene	0.400	0.351	12.3	97	-0.02
22 I	Chrysene-d12	5.000	5.000	0.0	100	-0.02
23	Pyrene	0.400	0.393	1.8	97	0.00
24 S	Terphenyl-d14	0.400	0.380	5.0	93	0.00
25	Benzo(a)anthracene	0.400	0.380	5.0	93	-0.02
26	Chrysene	0.400	0.401	-0.3	98	0.00
27	Indeno(1,2,3-cd)pyrene	0.400	0.403	-0.8	102	-0.05
28 I	Perylene-d12	5.000	5.000	0.0	102	-0.02
29	Benzo(b)fluoranthene	0.400	0.376	6.0	95	-0.02
30	Benzo(k)fluoranthene	0.400	0.389	2.8	102	-0.02
31 C	Benzo(a)pyrene	0.400	0.386	3.5	100	-0.03
32	Dibenzo(a,h)anthracene	0.400	0.388	3.0	100	-0.05
33	Benzo(g,h,i)perylene	0.400	0.392	2.0	100	-0.05

(#) = Out of Range

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