

Data Path : Z:\HPCHEM1\BNA E\DATA\BE111616\
 Data File : BE092538.D
 Acq On : 16 Nov 2016 17:51
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Nov 17 14:09:20 2016
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE111616.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 16 15:43:18 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	0.416	-4.0	120	0.00
3	n-Nitrosodimethylamine	0.400	0.468	-17.0	116	0.00
4 S	2-Fluorophenol	0.400	0.398	0.5	102	0.00
5 S	Phenol-d6	0.400	0.403	-0.8	104	0.00
6	bis(2-Chloroethyl)ether	0.400	0.339	15.3	99	0.00
7 I	Naphthalene-d8	5.000	5.000	0.0	102	0.00
8 S	Nitrobenzene-d5	0.400	0.399	0.3	102	0.00
9	Naphthalene	0.400	0.404	-1.0	103	0.00
10	Hexachlorobutadiene	0.400	0.417	-4.2	105	0.00
11	2-Methylnaphthalene	0.400	0.390	2.5	105	0.00
12 I	Acenaphthene-d10	5.000	5.000	0.0	105	0.00
13 S	2,4,6-Tribromophenol	0.400	0.348	13.0	97	0.00
14 S	2-Fluorobiphenyl	0.400	0.397	0.8	104	0.00
15	Acenaphthylene	0.400	0.381	4.8	104	0.00
16	Acenaphthene	0.400	0.393	1.8	103	0.00
17	Fluorene	0.400	0.394	1.5	105	0.00
18 I	Phenanthrene-d10	5.000	5.000	0.0	98	0.00
19	Hexachlorobenzene	0.400	0.400	0.0	110	0.00
20	Pentachlorophenol	0.400	0.448	-12.0	119	0.00
21	Phenanthrene	0.400	0.408	-2.0	104	0.00
22	Anthracene	0.400	0.399	0.3	105	0.00
23	Fluoranthene	0.400	0.402	-0.5	105	0.00
24 I	Chrysene-d12	5.000	5.000	0.0	104	0.00
25	Pyrene	0.400	0.372	7.0	103	0.00
26 S	Terphenyl-d14	0.400	0.373	6.8	105	0.00
27	Benzo(a)anthracene	0.400	0.448	-12.0	110	0.00
28	Chrysene	0.400	0.406	-1.5	102	-0.02
29	Bis(2-ethylhexyl)phthalate	0.400	0.445	-11.2	109	-0.02
30	Indeno(1,2,3-cd)pyrene	0.400	0.421	-5.2	101	0.00
31 I	Perylene-d12	5.000	5.000	0.0	101	-0.01
32	Benzo(b)fluoranthene	0.400	0.403	-0.8	106	0.00
33	Benzo(k)fluoranthene	0.400	0.368	8.0	99	0.00
34 C	Benzo(a)pyrene	0.400	0.371	7.3	102	0.00
35	Dibenzo(a,h)anthracene	0.400	0.369	7.8	100	-0.02
36	Benzo(g,h,i)perylene	0.400	0.380	5.0	101	0.00

(#) = Out of Range

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