

Data Path : Z:\HPCHEM1\BNA E\DATA\BE090116\
 Data File : BE092339.D
 Acq On : 1 Sep 2016 21:27
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Sep 02 11:13:11 2016
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE090116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 02 11:02:43 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	95	0.00
2	1,4-Dioxane	0.400	0.454	-13.5	111	0.00
3	n-Nitrosodimethylamine	0.400	0.356	11.0	90	0.00
4 S	2-Fluorophenol	0.400	0.378	5.5	99	0.00
5 S	Phenol-d6	0.400	0.327	18.3	99	0.00
6	bis(2-Chloroethyl)ether	0.400	0.434	-8.5	102	0.02
7 I	Naphthalene-d8	5.000	5.000	0.0	92	0.00
8 S	Nitrobenzene-d5	0.400	0.333	16.8	100	0.00
9	Naphthalene	0.400	0.428	-7.0	106	0.00
10	Hexachlorobutadiene	0.400	0.424	-6.0	105	0.00
11	2-Methylnaphthalene	0.400	0.413	-3.2	104	0.01
12 I	Acenaphthene-d10	5.000	5.000	0.0	92	-0.02
13 S	2,4,6-Tribromophenol	0.400	0.335	16.3	90	0.00
14 S	2-Fluorobiphenyl	0.400	0.401	-0.3	100	0.00
15	Acenaphthylene	0.400	0.409	-2.2	103	0.00
16	Acenaphthene	0.400	0.423	-5.7	104	0.00
17	Fluorene	0.400	0.410	-2.5	101	0.00
18 I	Phenanthrene-d10	5.000	5.000	0.0	92	0.00
19	Hexachlorobenzene	0.400	0.426	-6.5	105	0.00
20	Pentachlorophenol	0.400	0.375	6.3	90	0.00
21	Phenanthrene	0.400	0.435	-8.7	104	0.00
22	Anthracene	0.400	0.402	-0.5	101	0.00
23	Fluoranthene	0.400	0.412	-3.0	99	0.00
24 I	Chrysene-d12	5.000	5.000	0.0	94	0.00
25	Pyrene	0.400	0.393	1.8	98	0.00
26 S	Terphenyl-d14	0.400	0.377	5.8	94	0.00
27	Benzo(a)anthracene	0.400	0.405	-1.3	101	0.00
28	Chrysene	0.400	0.451	-12.7	105	0.00
29	Bis(2-ethylhexyl)phthalate	0.400	0.307	23.3	76	0.00
30	Indeno(1,2,3-cd)pyrene	0.400	0.408	-2.0	100	0.00
31 I	Perylene-d12	5.000	5.000	0.0	90	0.00
32	Benzo(b)fluoranthene	0.400	0.380	5.0	100	0.00
33	Benzo(k)fluoranthene	0.400	0.436	-9.0	99	-0.01
34 C	Benzo(a)pyrene	0.400	0.408	-2.0	100	0.00
35	Dibenzo(a,h)anthracene	0.400	0.403	-0.8	100	-0.02
36	Benzo(g,h,i)perylene	0.400	0.412	-3.0	100	0.00

(#) = Out of Range

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