

Data Path : Z:\HPCHEM1\BNA E\DATA\BE060816\
 Data File : BE091907.D
 Acq On : 8 Jun 2016 18:48
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jun 09 05:03:51 2016
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE060716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 08 16:11:34 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	102	0.00
2	1,4-Dioxane	0.400	0.408	-2.0	96	0.00
3	n-Nitrosodimethylamine	0.400	0.380	5.0	94	0.00
4 S	2-Fluorophenol	0.400	0.412	-3.0	106	0.00
5 S	Phenol-d6	0.400	0.424	-6.0	113	0.00
6	bis(2-Chloroethyl)ether	0.400	0.387	3.3	100	0.00
7	n-Nitroso-di-n-propylamine	0.400	0.376	6.0	99	0.00
8 I	Naphthalene-d8	5.000	5.000	0.0	104	0.00
9 S	Nitrobenzene-d5	0.400	0.402	-0.5	104	0.00
10	Nitrobenzene	0.400	0.395	1.3	105	0.00
11	bis(2-Chloroethoxy)methane	0.400	0.375	6.3	104	0.00
12	Naphthalene	0.400	0.425	-6.2	105	0.00
13	Hexachlorobutadiene	0.400	0.431	-7.7	105	0.00
14	2-Methylnaphthalene	0.400	0.406	-1.5	104	0.00
15 I	Acenaphthene-d10	5.000	5.000	0.0	110	0.00
16 S	2,4,6-Tribromophenol	0.400	0.432	-8.0	129	-0.02
17 S	2-Fluorobiphenyl	0.400	0.436	-9.0	110	0.00
18	Acenaphthylene	0.400	0.384	4.0	105	0.00
19	2,6-Dinitrotoluene	0.400	0.381	4.8	105	-0.02
20	Acenaphthene	0.400	0.408	-2.0	107	0.00
21	2,4-Dinitrotoluene	0.400	0.411	-2.7	106	-0.02
22	Fluorene	0.400	0.396	1.0	106	-0.02
23	4-Chlorophenyl-phenylether	0.400	0.408	-2.0	105	0.00
24 I	Phenanthrene-d10	5.000	5.000	0.0	109	0.00
25	4-Bromophenyl-phenylether	0.400	0.360	10.0	105	0.00
26	Hexachlorobenzene	0.400	0.387	3.3	108	0.00
27	Pentachlorophenol	0.400	0.387	3.3	114	0.00
28	Phenanthrene	0.400	0.379	5.3	106	-0.01
29	Anthracene	0.400	0.362	9.5	109	0.00
30	Fluoranthene	0.400	0.363	9.3	106	-0.02
31 I	Chrysene-d12	5.000	5.000	0.0	100	0.00
32	Benzidine	0.400	0.399	0.3	102	0.00
33	Pyrene	0.400	0.519	-29.8#	108	0.00
34 S	Terphenyl-d14	0.400	0.487	-21.7	114	0.00
35	Benzo(a)anthracene	0.400	0.405	-1.3	102	0.00
36	3,3'-Dichlorobenzidine	0.400	0.408	-2.0	101	0.00
37	Chrysene	0.400	0.327	18.3	81	0.00
38	Bis(2-ethylhexyl)phthalate	0.400	0.350	12.5	90	0.00
39	Indeno(1,2,3-cd)pyrene	0.400	0.418	-4.5	97	0.00
40 I	Perylene-d12	5.000	5.000	0.0	99	0.00
41	Benzo(b)fluoranthene	0.400	0.390	2.5	95	0.00

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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	0.400	0.401	-0.3	100	0.00
43 C	Benzo(a)pyrene	0.400	0.392	2.0	100	0.00
44	Dibenzo(a,h)anthracene	0.400	0.385	3.8	97	-0.02
45	Benzo(a,h,i)perylene	0.400	0.393	1.8	96	-0.02

(#) = Out of Range

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