

Data Path : Z:\HPCHEM1\BNA E\DATA\BE060716\
 Data File : BE091898.D
 Acq On : 8 Jun 2016 5:05
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jun 08 07:58:06 2016
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE060716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 07 17:14:38 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	101	0.00
2	1,4-Dioxane	0.400	0.447	-11.7	103	0.00
3	n-Nitrosodimethylamine	0.400	0.388	3.0	95	0.00
4 S	2-Fluorophenol	0.400	0.435	-8.7	111	0.00
5 S	Phenol-d6	0.400	0.430	-7.5	114	0.00
6	bis(2-Chloroethyl)ether	0.400	0.390	2.5	99	0.00
7	n-Nitroso-di-n-propylamine	0.400	0.393	1.8	102	0.00
8 I	Naphthalene-d8	5.000	5.000	0.0	103	0.00
9 S	Nitrobenzene-d5	0.400	0.381	4.8	96	-0.02
10	Nitrobenzene	0.400	0.409	-2.2	108	0.00
11	bis(2-Chloroethoxy)methane	0.400	0.398	0.5	109	0.00
12	Naphthalene	0.400	0.429	-7.2	105	0.00
13	Hexachlorobutadiene	0.400	0.428	-7.0	103	0.00
14	2-Methylnaphthalene	0.400	0.428	-7.0	108	0.00
15 I	Acenaphthene-d10	5.000	5.000	0.0	111	0.00
16 S	2,4,6-Tribromophenol	0.400	0.434	-8.5	129	-0.02
17 S	2-Fluorobiphenyl	0.400	0.411	-2.7	104	0.00
18	Acenaphthylene	0.400	0.385	3.8	105	0.00
19	2,6-Dinitrotoluene	0.400	0.399	0.3	113	-0.02
20	Acenaphthene	0.400	0.415	-3.7	109	0.00
21	2,4-Dinitrotoluene	0.400	0.412	-3.0	107	-0.02
22	Fluorene	0.400	0.410	-2.5	110	-0.02
23	4-Chlorophenyl-phenylether	0.400	0.416	-4.0	108	0.00
24 I	Phenanthrene-d10	5.000	5.000	0.0	96	0.00
25	4-Bromophenyl-phenylether	0.400	0.422	-5.5	108	0.00
26	Hexachlorobenzene	0.400	0.434	-8.5	106	0.00
27	Pentachlorophenol	0.400	0.501	-25.2#	134	0.00
28	Phenanthrene	0.400	0.443	-10.7	109	-0.01
29	Anthracene	0.400	0.428	-7.0	113	-0.01
30	Fluoranthene	0.400	0.444	-11.0	114	-0.02
31 I	Chrysene-d12	5.000	5.000	0.0	115	-0.03
32	Benzidine	0.400	0.408	-2.0	121	0.00
33	Pyrene	0.400	0.512	-28.0#	123	0.00
34 S	Terphenyl-d14	0.400	0.406	-1.5	107	0.00
35	Benzo(a)anthracene	0.400	0.351	12.3	108	0.00
36	3,3'-Dichlorobenzidine	0.400	0.466	-16.5	137	0.00
37	Chrysene	0.400	0.221	44.8#	63	-0.03
38	Bis(2-ethylhexyl)phthalate	0.400	0.350	12.5	104	0.00
39	Indeno(1,2,3-cd)pyrene	0.400	0.392	2.0	104	0.00
40 I	Perylene-d12	5.000	5.000	0.0	102	0.00
41	Benzo(b)fluoranthene	0.400	0.416	-4.0	105	-0.01

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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	0.400	0.427	-6.7	109	0.00
43 C	Benzo(a)pyrene	0.400	0.424	-6.0	111	0.00
44	Dibenzo(a,h)anthracene	0.400	0.415	-3.7	108	-0.02
45	Benzo(a,h,i)perylene	0.400	0.421	-5.2	105	-0.02

(#) = Out of Range

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