

Data Path : Z:\HPCHEM1\BNA E\DATA\BE052516\
 Data File : BE091867.D
 Acq On : 25 May 2016 20:26
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: May 26 01:49:34 2016
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE052516.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 25 19:42:30 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	113	0.00
2	1,4-Dioxane	0.400	0.414	-3.5	106	0.00
3	n-Nitrosodimethylamine	0.400	0.337	15.8	108	0.00
4 S	2-Fluorophenol	0.400	0.370	7.5	114	0.00
5 S	Phenol-d6	0.400	0.383	4.3	123	0.00
6	bis(2-Chloroethyl)ether	0.400	0.375	6.3	116	-0.02
7	n-Nitroso-di-n-propylamine	0.400	0.372	7.0	123	0.00
8 I	Naphthalene-d8	5.000	5.000	0.0	122	0.00
9 S	Nitrobenzene-d5	0.400	0.370	7.5	123	0.00
10	Nitrobenzene	0.400	0.367	8.3	128	0.00
11	bis(2-Chloroethoxy)methane	0.400	0.319	20.3	123	0.00
12	Naphthalene	0.400	0.386	3.5	125	0.00
13	Hexachlorobutadiene	0.400	0.384	4.0	121	0.00
14	2-Methylnaphthalene	0.400	0.391	2.3	129	0.00
15 I	Acenaphthene-d10	5.000	5.000	0.0	128	0.00
16 S	2,4,6-Tribromophenol	0.400	0.331	17.3	120	0.00
17 S	2-Fluorobiphenyl	0.400	0.379	5.3	126	0.00
18	Acenaphthylene	0.400	0.387	3.3	129	0.00
19	2,6-Dinitrotoluene	0.400	0.327	18.3	125	-0.02
20	Acenaphthene	0.400	0.396	1.0	130	0.00
21	2,4-Dinitrotoluene	0.400	0.453	-13.2	130	0.00
22	Fluorene	0.400	0.378	5.5	129	0.00
23	4-Chlorophenyl-phenylether	0.400	0.343	14.2	123	0.00
24 I	Phenanthrene-d10	5.000	5.000	0.0	118	0.00
25	4-Bromophenyl-phenylether	0.400	0.387	3.3	122	0.00
26	Hexachlorobenzene	0.400	0.459	-14.8	140	0.00
27	Pentachlorophenol	0.400	0.395	1.3	150	-0.02
28	Phenanthrene	0.400	0.361	9.8	118	0.00
29	Anthracene	0.400	0.382	4.5	123	0.00
30	Fluoranthene	0.400	0.346	13.5	116	-0.02
31 I	Chrysene-d12	5.000	5.000	0.0	95	0.00
32	Benzidine	0.400	0.445	-11.2	116	0.00
33	Pyrene	0.400	0.444	-11.0	115	0.00
34 S	Terphenyl-d14	0.400	0.431	-7.7	109	0.00
35	Benzo(a)anthracene	0.400	0.433	-8.2	109	0.00
36	3,3'-Dichlorobenzidine	0.400	0.408	-2.0	107	0.00
37	Chrysene	0.400	0.350	12.5	102	0.00
38	Bis(2-ethylhexyl)phthalate	0.400	0.490	-22.5	137	0.00
39	Indeno(1,2,3-cd)pyrene	0.400	0.420	-5.0	107	0.00
40 I	Perylene-d12	5.000	5.000	0.0	101	0.00
41	Benzo(b)fluoranthene	0.400	0.414	-3.5	119	0.00

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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	0.400	0.369	7.8	91	0.00
43 C	Benzo(a)pyrene	0.400	0.397	0.8	107	0.00
44	Dibenzo(a,h)anthracene	0.400	0.396	1.0	106	0.00
45	Benzo(a,h,i)perylene	0.400	0.399	0.3	107	0.00

(#) = Out of Range

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