

Data Path : Z:\HPCHEM1\BNA E\DATA\BE052516\
 Data File : BE091874.D
 Acq On : 26 May 2016 00:52
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: May 26 02:04:23 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	126	0.00
2	1,4-Dioxane	0.400	0.369	7.8	105	0.00
3	n-Nitrosodimethylamine	0.400	0.334	16.5	120	0.00
4 S	2-Fluorophenol	0.400	0.346	13.5	119	0.00
5 S	Phenol-d6	0.400	0.389	2.8	140	0.00
6	bis(2-Chloroethyl)ether	0.400	0.373	6.8	129	-0.02
7	n-Nitroso-di-n-propylamine	0.400	0.368	8.0	135	0.00
8 I	Naphthalene-d8	5.000	5.000	0.0	141	0.00
9 S	Nitrobenzene-d5	0.400	0.368	8.0	141	0.00
10	Nitrobenzene	0.400	0.366	8.5	146	0.00
11	bis(2-Chloroethoxy)methane	0.400	0.308	23.0	136	0.00
12	Naphthalene	0.400	0.389	2.8	145	0.00
13	Hexachlorobutadiene	0.400	0.384	4.0	139	0.00
14	2-Methylnaphthalene	0.400	0.386	3.5	146	0.00
15 I	Acenaphthene-d10	5.000	5.000	0.0	142	0.00
16 S	2,4,6-Tribromophenol	0.400	0.316	21.0	126	0.00
17 S	2-Fluorobiphenyl	0.400	0.385	3.8	142	-0.02
18	Acenaphthylene	0.400	0.392	2.0	145	0.00
19	2,6-Dinitrotoluene	0.400	0.307	23.3	130	-0.02
20	Acenaphthene	0.400	0.390	2.5	143	0.00
21	2,4-Dinitrotoluene	0.400	0.433	-8.2	133	0.00
22	Fluorene	0.400	0.368	8.0	139	0.00
23	4-Chlorophenyl-phenylether	0.400	0.331	17.3	133	0.00
24 I	Phenanthrene-d10	5.000	5.000	0.0	124	0.00
25	4-Bromophenyl-phenylether	0.400	0.393	1.8	131	0.00
26	Hexachlorobenzene	0.400	0.410	-2.5	132	0.00
27	Pentachlorophenol	0.400	0.377	5.8	151	-0.02
28	Phenanthrene	0.400	0.365	8.8	126	0.00
29	Anthracene	0.400	0.385	3.8	131	0.00
30	Fluoranthene	0.400	0.327	18.3	117	-0.02
31 I	Chrysene-d12	5.000	5.000	0.0	99	0.00
32	Benzidine	0.400	0.486	-21.5	136	0.00
33	Pyrene	0.400	0.441	-10.2	119	0.00
34 S	Terphenyl-d14	0.400	0.418	-4.5	110	0.00
35	Benzo(a)anthracene	0.400	0.418	-4.5	110	0.00
36	3,3'-Dichlorobenzidine	0.400	0.410	-2.5	112	0.00
37	Chrysene	0.400	0.346	13.5	105	0.00
38	Bis(2-ethylhexyl)phthalate	0.400	0.752	-88.0#	219	0.00
39	Indeno(1,2,3-cd)pyrene	0.400	0.399	0.3	106	0.00
40 I	Perylene-d12	5.000	5.000	0.0	100	0.00
41	Benzo(b)fluoranthene	0.400	0.365	8.8	104	0.00

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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	0.400	0.419	-4.7	102	-0.01
43 C	Benzo(a)pyrene	0.400	0.398	0.5	106	0.00
44	Dibenzo(a,h)anthracene	0.400	0.395	1.3	105	0.00
45	Benzo(a,h,i)perylene	0.400	0.397	0.8	105	0.00

(#) = Out of Range

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