

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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	102	0.00
2	1,4-Dioxane	0.781	0.774	0.9	96	0.00
3	n-Nitrosodimethylamine	0.836	0.795	4.9	94	0.00
4 S	2-Fluorophenol	1.066	1.098	-3.0	106	0.00
5 S	Phenol-d6	1.320	1.400	-6.1	113	0.00
6	bis(2-Chloroethyl)ether	1.431	1.386	3.1	100	0.00
7	n-Nitroso-di-n-propylamine	1.130	1.061	6.1	99	0.00
8 I	Naphthalene-d8	1.000	1.000	0.0	104	0.00
9 S	Nitrobenzene-d5	0.336	0.347	-3.3	104	0.00
10	Nitrobenzene	0.304	0.300	1.3	105	0.00
11	bis(2-Chloroethoxy)methane	0.507	0.476	6.1	104	0.00
12	Naphthalene	1.079	1.146	-6.2	105	0.00
13	Hexachlorobutadiene	0.161	0.173	-7.5	105	0.00
14	2-Methylnaphthalene	0.696	0.707	-1.6	104	0.00
15 I	Acenaphthene-d10	1.000	1.000	0.0	110	0.00
16 S	2,4,6-Tribromophenol	0.142	0.154	-8.5	129	-0.02
17 S	2-Fluorobiphenyl	1.517	1.652	-8.9	110	0.00
18	Acenaphthylene	2.263	2.173	4.0	105	0.00
19	2,6-Dinitrotoluene	0.240	0.210	12.5	105	-0.02
20	Acenaphthene	1.383	1.410	-2.0	107	0.00
21	2,4-Dinitrotoluene	0.285	0.227	20.4	106	-0.02
22	Fluorene	1.691	1.674	1.0	106	-0.02
23	4-Chlorophenyl-phenylether	0.829	0.847	-2.2	105	0.00
24 I	Phenanthrene-d10	1.000	1.000	0.0	109	0.00
25	4-Bromophenyl-phenylether	0.180	0.162	10.0	105	0.00
26	Hexachlorobenzene	0.184	0.178	3.3	108	0.00
27	Pentachlorophenol	0.070	0.052	25.7#	114	0.00
28	Phenanthrene	1.158	1.098	5.2	106	-0.01
29	Anthracene	1.032	0.934	9.5	109	0.00
30	Fluoranthene	1.250	1.135	9.2	106	-0.02
31 I	Chrysene-d12	1.000	1.000	0.0	100	0.00
32	Benzidine	0.333	0.306	8.1	102	0.00
33	Pyrene	1.262	1.636	-29.6#	108	0.00
34 S	Terphenyl-d14	0.891	1.243	-39.5#	114	0.00
35	Benzo(a)anthracene	7.222	8.749	-21.1	102	0.00
36	3,3'-Dichlorobenzidine	0.349	0.318	8.9	101	0.00
37	Chrysene	5.098	4.170	18.2	81	0.00
38	Bis(2-ethylhexyl)phthalate	0.716	0.538	24.9	90	0.00
39	Indeno(1,2,3-cd)pyrene	1.524	1.797	-17.9	97	0.00
40 I	Perylene-d12	1.000	1.000	0.0	99	0.00
41	Benzo(b)fluoranthene	1.222	1.190	2.6	95	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	1.268	1.272	-0.3	100	0.00
43 C	Benzo(a)pyrene	1.146	1.123	2.0	100	0.00
44	Dibenzo(a,h)anthracene	1.068	1.029	3.7	97	-0.02
45	Benzo(a,h,i)perylene	1.120	1.099	1.9	96	-0.02

(#) = Out of Range

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