

Data Path : Z:\HPCHEM1\BNA E\DATA\BE100715\
 Data File : BE090987.D
 Acq On : 7 Oct 2015 14:04
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC001

Quant Time: Oct 08 01:44:51 2015
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE092815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 07 17:11:43 2015
 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	176	0.00
2	1,4-Dioxane	1.000	2.338	-133.8#	367	0.00
3	n-Nitrosodimethylamine	1.000	0.679	32.1#	124	0.00
4 S	2-Fluorophenol	1.000	1.092	-9.2	201	0.00
5 S	Phenol-d6	1.000	0.957	4.3	189	0.00
6	bis(2-Chloroethyl)ether	1.000	2.213	-121.3#	357	0.00
7 I	Naphthalene-d8	5.000	5.000	0.0	174	0.00
8 S	Nitrobenzene-d5	1.000	1.237	-23.7	247	0.00
9	Nitrobenzene	1.000	1.275	-27.5#	266	0.00
10	Naphthalene	1.000	0.993	0.7	173	0.00
11	Hexachlorobutadiene	1.000	0.850	15.0	147	0.00
12	2-Methylnaphthalene	1.000	1.053	-5.3	195	0.00
13 I	Acenaphthene-d10	5.000	5.000	0.0	169	0.00
14 S	2,4,6-Tribromophenol	1.000	0.934	6.6	192	0.00
15 S	2-Fluorobiphenyl	1.000	1.004	-0.4	182	0.00
16	Acenaphthylene	1.000	0.967	3.3	178	0.00
17	Acenaphthene	1.000	0.957	4.3	172	0.00
18	Fluorene	1.000	0.989	1.1	179	0.00
19 I	Phenanthrene-d10	5.000	5.000	0.0	177	0.00
20	4-Bromophenyl-phenylether	1.000	0.978	2.2	188	0.00
21	Hexachlorobenzene	1.000	0.123	87.7#	39	0.00
22	Pentachlorophenol	1.000	0.828	17.2	177	0.00
23	Phenanthrene	1.000	0.967	3.3	179	0.00
24	Anthracene	1.000	1.017	-1.7	194	0.00
25	Fluoranthene	1.000	1.042	-4.2	201	0.00
26 I	Chrysene-d12	5.000	5.000	0.0	195	0.00
27	Benzydine	1.000	1.845	-84.5#	415	0.00
28	Pvrene	1.000	0.937	6.3	199	0.00
29 S	Terphenyl-d14	1.000	1.032	-3.2	221	0.00
30	Benzo(a)anthracene	1.000	1.077	-7.7	221	0.00
31	3,3'-Dichlorobenzidine	1.000	0.022	97.8#	0	-20.84#
32	Chrysene	1.000	0.922	7.8	187	0.00
33	Bis(2-ethylhexyl)phthalate	1.000	1.456	-45.6#	367	0.00
34	Indeno(1,2,3-cd)pyrene	1.000	1.224	-22.4	266	0.00
35 I	Perylene-d12	5.000	5.000	0.0	239	0.00
36	Benzo(b)fluoranthene	1.000	1.056	-5.6	281	0.00
37	Benzo(k)fluoranthene	1.000	0.872	12.8	222	0.00
38 C	Benzo(a)pyrene	1.000	1.097	-9.7	304	0.00
39	Dibenzo(a,h)anthracene	1.000	1.270	-27.0#	361	0.00
40	Benzo(a,h,i)perylene	1.000	1.207	-20.7	332	0.00

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Compound	Amount	Calc.	%Dev Area	% Dev(min)

(#) = Out of Range				
SPCC's out = 0 CCC's out = 0				