

Data Path : Z:\HPCHEM1\BNA M\DATA\BM030116\
 Data File : BM004512.D
 Acq On : 02 Mar 2016 10:10
 Operator : SJ/UM
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC001

Quant Time: Mar 02 15:20:35 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM030116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 02 14:46:57 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	114	0.00
2	1,4-Dioxane	1.000	0.937	6.3	114	0.00
3	n-Nitrosodimethylamine	1.000	0.918	8.2	116	0.00
4 S	2-Fluorophenol	1.000	0.856	14.4	111	0.00
5 S	Phenol-d6	1.000	0.903	9.7	113	0.00
6	bis(2-Chloroethyl)ether	1.000	0.917	8.3	112	0.00
7 I	Naphthalene-d8	5.000	5.000	0.0	113	0.00
8 S	Nitrobenzene-d5	1.000	0.909	9.1	112	0.00
9	Nitrobenzene	1.000	0.898	10.2	113	0.00
10	Naphthalene	1.000	0.913	8.7	113	0.00
11	Hexachlorobutadiene	1.000	0.922	7.8	113	0.00
12	2-Methylnaphthalene	1.000	0.911	8.9	112	0.00
13 I	Acenaphthene-d10	5.000	5.000	0.0	111	0.00
14 S	2,4,6-Tribromophenol	1.000	0.848	15.2	107	0.00
15 S	2-Fluorobiphenyl	1.000	0.940	6.0	111	0.00
16	Acenaphthylene	1.000	0.908	9.2	113	0.00
17	Acenaphthene	1.000	0.950	5.0	112	0.00
18	Fluorene	1.000	0.958	4.2	110	0.00
19 I	Phenanthrene-d10	5.000	5.000	0.0	109	0.00
20	4-Bromophenyl-phenylether	1.000	0.906	9.4	111	0.00
21	Hexachlorobenzene	1.000	1.004	-0.4	112	0.00
22	Pentachlorophenol	1.000	0.794	20.6	105	0.00
23	Phenanthrene	1.000	0.972	2.8	108	0.00
24	Anthracene	1.000	0.873	12.7	108	0.00
25	Fluoranthene	1.000	0.887	11.3	106	0.00
26 I	Chrysene-d12	5.000	5.000	0.0	110	0.00
27	Benzydine	1.000	0.908	9.2	112	0.00
28	Pvrene	1.000	0.938	6.2	107	0.00
29 S	Terphenyl-d14	1.000	0.934	6.6	107	0.00
30	Benzo(a)anthracene	1.000	0.893	10.7	105	0.02
31	3,3'-Dichlorobenzidine	1.000	0.933	6.7	109	0.00
32	Chrysene	1.000	0.932	6.8	108	0.00
33	Bis(2-ethylhexyl)phthalate	1.000	0.790	21.0	95	0.00
34	Indeno(1,2,3-cd)pyrene	1.000	0.957	4.3	112	0.00
35 I	Perylene-d12	5.000	5.000	0.0	111	0.00
36	Benzo(b)fluoranthene	1.000	1.027	-2.7	123	0.00
37	Benzo(k)fluoranthene	1.000	0.903	9.7	113	0.00
38 C	Benzo(a)pyrene	1.000	0.898	10.2	111	0.00
39	Dibenzo(a,h)anthracene	1.000	0.906	9.4	112	0.01
40	Benzo(a,h,i)perylene	1.000	0.909	9.1	112	0.00

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(#) = Out of Range

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