

Data Path : Z:\HPCHEM1\BNA M\DATA\BM122415\
 Data File : BM003756.D
 Acq On : 24 Dec 2015 00:24
 Operator : UM/SJ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC001

Quant Time: Dec 24 10:44:16 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM122415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 24 10:08:49 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	100	0.00
2	1,4-Dioxane	1.000	0.905	9.5	100	0.00
3	n-Nitrosodimethylamine	1.000	0.857	14.3	99	0.00
4 S	2-Fluorophenol	1.000	0.839	16.1	98	0.00
5 S	Phenol-d6	1.000	0.819	18.1	98	0.00
6	bis(2-Chloroethyl)ether	1.000	0.863	13.7	99	0.00
7 I	Naphthalene-d8	5.000	5.000	0.0	102	0.00
8 S	Nitrobenzene-d5	1.000	0.825	17.5	100	0.00
9	Nitrobenzene	1.000	0.829	17.1	100	0.00
10	Naphthalene	1.000	0.884	11.6	102	0.00
11	Hexachlorobutadiene	1.000	0.893	10.7	102	0.00
12	2-Methylnaphthalene	1.000	0.879	12.1	102	0.00
13 I	Acenaphthene-d10	5.000	5.000	0.0	105	0.00
14 S	2,4,6-Tribromophenol	1.000	0.817	18.3	101	0.00
15 S	2-Fluorobiphenyl	1.000	0.865	13.5	103	0.00
16	Acenaphthylene	1.000	0.826	17.4	105	0.00
17	Acenaphthene	1.000	0.926	7.4	105	0.00
18	Fluorene	1.000	0.862	13.8	104	0.00
19 I	Phenanthrene-d10	5.000	5.000	0.0	107	0.00
20	4-Bromophenyl-phenylether	1.000	1.303	-30.3#	107	0.00
21	Hexachlorobenzene	1.000	1.245	-24.5	106	0.00
22	Pentachlorophenol	1.000	0.839	16.1	109	0.00
23	Phenanthrene	1.000	1.189	-18.9	108	0.00
24	Anthracene	1.000	0.912	8.8	105	0.00
25	Fluoranthene	1.000	0.950	5.0	107	0.00
26 I	Chrysene-d12	5.000	5.000	0.0	114	0.00
27	Benzydine	1.000	1.148	-14.8	139	0.00
28	Pvrene	1.000	0.864	13.6	108	0.00
29 S	Terphenyl-d14	1.000	0.883	11.7	109	0.00
30	Benzo(a)anthracene	1.000	0.822	17.8	113	0.00
31	3,3'-Dichlorobenzidine	1.000	0.925	7.5	108	0.00
32	Chrysene	1.000	0.913	8.7	109	0.00
33	Bis(2-ethylhexyl)phthalate	1.000	0.842	15.8	106	0.00
34	Indeno(1,2,3-cd)pyrene	1.000	0.879	12.1	108	0.01
35 I	Perylene-d12	5.000	5.000	0.0	112	0.00
36	Benzo(b)fluoranthene	1.000	0.877	12.3	109	0.00
37	Benzo(k)fluoranthene	1.000	0.859	14.1	116	0.00
38 C	Benzo(a)pyrene	1.000	0.847	15.3	112	0.00
39	Dibenzo(a,h)anthracene	1.000	0.844	15.6	108	0.00
40	Benzo(a,h,i)perylene	1.000	0.840	16.0	106	0.00

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

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