

Data Path : Z:\HPCHEM1\BNA M\DATA\BM110515\
 Data File : BM003068.D
 Acq On : 05 Nov 2015 17:34
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC001

Quant Time: Nov 17 13:18:23 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM110415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 17 12:50:51 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	133	0.00
2	1,4-Dioxane	1.000	0.968	3.2	126	0.00
3	n-Nitrosodimethylamine	1.000	0.958	4.2	132	0.00
4 S	2-Fluorophenol	1.000	0.983	1.7	137	0.00
5 S	Phenol-d6	1.000	0.995	0.5	141	0.00
6	bis(2-Chloroethyl)ether	1.000	0.967	3.3	133	-0.02
7 I	Naphthalene-d8	5.000	5.000	0.0	136	0.00
8 S	Nitrobenzene-d5	1.000	1.018	-1.8	146	0.00
9	Nitrobenzene	1.000	1.010	-1.0	145	0.00
10	Naphthalene	1.000	0.971	2.9	134	0.00
11	Hexachlorobutadiene	1.000	0.991	0.9	136	0.00
12	2-Methylnaphthalene	1.000	0.982	1.8	135	0.00
13 I	Acenaphthene-d10	5.000	5.000	0.0	138	0.00
14 S	2,4,6-Tribromophenol	1.000	1.006	-0.6	141	0.00
15 S	2-Fluorobiphenyl	1.000	0.961	3.9	134	0.00
16	Acenaphthylene	1.000	0.998	0.2	144	0.00
17	Acenaphthene	1.000	0.975	2.5	146	0.00
18	Fluorene	1.000	0.950	5.0	128	0.02
19 I	Phenanthrene-d10	5.000	5.000	0.0	169	0.00
20	4-Bromophenyl-phenylether	1.000	0.933	6.7	109	0.00
21	Hexachlorobenzene	1.000	0.958	4.2	119	0.00
22	Pentachlorophenol	1.000	1.041	-4.1	141	0.00
23	Phenanthrene	1.000	0.898	10.2	106	0.00
24	Anthracene	1.000	0.984	1.6	166	0.00
25	Fluoranthene	1.000	0.944	5.6	137	0.01
26 I	Chrysene-d12	5.000	5.000	0.0	151	0.00
27	Benzydine	1.000	0.973	2.7	130	0.00
28	Pvrene	1.000	0.906	9.4	140	0.00
29 S	Terphenyl-d14	1.000	0.917	8.3	134	0.00
30	Benzo(a)anthracene	1.000	0.970	3.0	148	0.00
31	3,3'-Dichlorobenzidine	1.000	0.983	1.7	130	0.00
32	Chrysene	1.000	0.906	9.4	120	0.01
33	Bis(2-ethylhexyl)phthalate	1.000	1.006	-0.6	167	0.01
34	Indeno(1,2,3-cd)pyrene	1.000	1.082	-8.2	145	0.00
35 I	Perylene-d12	5.000	5.000	0.0	141	0.00
36	Benzo(b)fluoranthene	1.000	0.952	4.8	137	0.00
37	Benzo(k)fluoranthene	1.000	0.925	7.5	132	0.00
38 C	Benzo(a)pyrene	1.000	0.927	7.3	144	0.00
39	Dibenzo(a,h)anthracene	1.000	0.989	1.1	147	0.00
40	Benzo(a,h,i)perylene	1.000	0.950	5.0	143	0.00

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

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