

Data Path : Z:\HPCHEM1\BNA M\DATA\BM101515\
 Data File : BM002919.D
 Acq On : 15 Oct 2015 18:08
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC001

Quant Time: Oct 16 06:59:44 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM101515.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 15 16:21:18 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	82	0.00
2	1,4-Dioxane	1.000	0.961	3.9	81	0.00
3	n-Nitrosodimethylamine	1.000	0.984	1.6	79	-0.02
4 S	2-Fluorophenol	1.000	0.954	4.6	80	-0.02
5 S	Phenol-d6	1.000	0.979	2.1	79	0.00
6	bis(2-Chloroethyl)ether	1.000	0.956	4.4	77	0.00
7 I	Naphthalene-d8	5.000	5.000	0.0	83	0.00
8 S	Nitrobenzene-d5	1.000	0.929	7.1	77	-0.02
9	Nitrobenzene	1.000	0.926	7.4	77	-0.01
10	Naphthalene	1.000	0.971	2.9	82	0.00
11	Hexachlorobutadiene	1.000	0.961	3.9	82	0.00
12	2-Methylnaphthalene	1.000	0.976	2.4	83	0.00
13 I	Acenaphthene-d10	5.000	5.000	0.0	87	0.00
14 S	2,4,6-Tribromophenol	1.000	0.836	16.4	77	0.00
15 S	2-Fluorobiphenyl	1.000	0.937	6.3	84	0.00
16	Acenaphthylene	1.000	0.930	7.0	83	0.00
17	Acenaphthene	1.000	0.896	10.4	82	0.00
18	Fluorene	1.000	0.937	6.3	83	0.00
19 I	Phenanthrene-d10	5.000	5.000	0.0	78	0.00
20	4-Bromophenyl-phenylether	1.000	1.007	-0.7	80	0.00
21	Hexachlorobenzene	1.000	1.558	-55.8#	92	0.00
22	Pentachlorophenol	1.000	1.105	-10.5	90	-0.02
23	Phenanthrene	1.000	1.156	-15.6	91	0.00
24	Anthracene	1.000	1.166	-16.6	90	0.00
25	Fluoranthene	1.000	1.097	-9.7	88	-0.01
26 I	Chrysene-d12	5.000	5.000	0.0	92	0.00
27	Benzydine	1.000	0.804	19.6	83	0.00
28	Pvrene	1.000	0.937	6.3	87	-0.01
29 S	Terphenyl-d14	1.000	0.935	6.5	87	0.00
30	Benzo(a)anthracene	1.000	0.893	10.7	80	0.00
31	3,3'-Dichlorobenzidine	1.000	0.867	13.3	84	0.00
32	Chrysene	1.000	0.954	4.6	96	0.00
33	Bis(2-ethylhexyl)phthalate	1.000	0.780	22.0	60	0.00
34	Indeno(1,2,3-cd)pyrene	1.000	0.950	5.0	95	0.00
35 I	Perylene-d12	5.000	5.000	0.0	95	0.00
36	Benzo(b)fluoranthene	1.000	0.949	5.1	90	0.00
37	Benzo(k)fluoranthene	1.000	0.987	1.3	97	0.00
38 C	Benzo(a)pyrene	1.000	0.938	6.2	91	0.00
39	Dibenzo(a,h)anthracene	1.000	0.966	3.4	96	0.00
40	Benzo(a,h,i)perylene	1.000	0.977	2.3	97	-0.01

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

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