

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM012716\
 Data File : BM004130.D
 Acq On : 27 Jan 2016 11:03
 Operator : SJ/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC001

Quant Time: Jan 28 00:01:47 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BM012216.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 22 14:46:41 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	82	0.00
2	1,4-Dioxane	1.000	0.963	3.7	83	0.00
3	n-Nitrosodimethylamine	1.000	0.891	10.9	80	0.00
4 S	2-Fluorophenol	1.000	0.906	9.4	82	0.00
5 S	Phenol-d6	1.000	0.898	10.2	82	0.00
6	bis(2-Chloroethyl)ether	1.000	0.913	8.7	83	0.00
7 I	Naphthalene-d8	5.000	5.000	0.0	85	0.00
8 S	Nitrobenzene-d5	1.000	0.918	8.2	86	0.00
9	Nitrobenzene	1.000	0.914	8.6	87	0.00
10	Naphthalene	1.000	0.915	8.5	84	0.00
11	Hexachlorobutadiene	1.000	0.908	9.2	84	0.00
12	2-Methylnaphthalene	1.000	0.924	7.6	86	0.00
13 I	Acenaphthene-d10	5.000	5.000	0.0	90	0.00
14 S	2,4,6-Tribromophenol	1.000	0.897	10.3	89	0.00
15 S	2-Fluorobiphenyl	1.000	0.885	11.5	87	0.00
16	Acenaphthylene	1.000	0.894	10.6	89	0.00
17	Acenaphthene	1.000	0.867	13.3	88	0.00
18	Fluorene	1.000	0.875	12.5	88	0.00
19 I	Phenanthrene-d10	5.000	5.000	0.0	97	0.00
20	4-Bromophenyl-phenylether	1.000	0.901	9.9	90	0.00
21	Hexachlorobenzene	1.000	0.909	9.1	93	0.00
22	Pentachlorophenol	1.000	0.901	9.9	107	0.00
23	Phenanthrene	1.000	0.922	7.8	97	0.00
24	Anthracene	1.000	0.908	9.2	95	0.00
25	Fluoranthene	1.000	0.910	9.0	96	0.00
26 I	Chrysene-d12	5.000	5.000	0.0	98	0.00
27	Benzydine	1.000	0.952	4.8	93	0.02
28	Pyrene	1.000	0.878	12.2	94	0.00
29 S	Terphenyl-d14	1.000	0.878	12.2	95	0.00
30	Benzo(a)anthracene	1.000	0.778	22.2	87	0.00
31	3,3'-Dichlorobenzidine	1.000	0.854	14.6	89	0.00
32	Chrysene	1.000	0.969	3.1	106	0.00
33	Bis(2-ethylhexyl)phthalate	1.000	0.915	8.5	101	-0.02
34	Indeno(1,2,3-cd)pyrene	1.000	0.799	20.1	89	0.00
35 I	Perylene-d12	5.000	5.000	0.0	92	0.00
36	Benzo(b)fluoranthene	1.000	0.945	5.5	94	0.00
37	Benzo(k)fluoranthene	1.000	0.882	11.8	92	0.00
38 C	Benzo(a)pyrene	1.000	0.912	8.8	94	0.00
39	Dibenzo(a,h)anthracene	1.000	0.842	15.8	86	0.00
40	Benzo(g,h,i)perylene	1.000	0.891	10.9	92	0.00

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

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