

Data Path : Z:\HPCHEM1\BNA M\DATA\BM013016\
 Data File : BM004147.D
 Acq On : 29 Jan 2016 15:37
 Operator : SJ/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC001

Quant Time: Jan 29 22:50:54 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	88	0.00
2	1,4-Dioxane	1.000	0.977	2.3	90	0.00
3	n-Nitrosodimethylamine	1.000	0.898	10.2	86	0.00
4 S	2-Fluorophenol	1.000	0.864	13.6	84	0.00
5 S	Phenol-d6	1.000	0.875	12.5	86	0.00
6	bis(2-Chloroethyl)ether	1.000	0.915	8.5	89	0.00
7 I	Naphthalene-d8	5.000	5.000	0.0	92	0.00
8 S	Nitrobenzene-d5	1.000	0.890	11.0	91	0.00
9	Nitrobenzene	1.000	0.893	10.7	91	0.00
10	Naphthalene	1.000	0.923	7.7	92	0.00
11	Hexachlorobutadiene	1.000	0.908	9.2	91	0.00
12	2-Methylnaphthalene	1.000	0.932	6.8	94	0.00
13 I	Acenaphthene-d10	5.000	5.000	0.0	97	0.00
14 S	2,4,6-Tribromophenol	1.000	0.888	11.2	95	0.00
15 S	2-Fluorobiphenyl	1.000	0.901	9.9	95	0.00
16	Acenaphthylene	1.000	0.887	11.3	96	0.00
17	Acenaphthene	1.000	0.878	12.2	96	0.00
18	Fluorene	1.000	0.925	7.5	101	0.00
19 I	Phenanthrene-d10	5.000	5.000	0.0	104	0.00
20	4-Bromophenyl-phenylether	1.000	0.894	10.6	95	0.00
21	Hexachlorobenzene	1.000	0.881	11.9	97	0.00
22	Pentachlorophenol	1.000	0.824	17.6	101	0.00
23	Phenanthrene	1.000	0.893	10.7	101	0.00
24	Anthracene	1.000	0.870	13.0	97	0.00
25	Fluoranthene	1.000	0.881	11.9	99	0.00
26 I	Chrysene-d12	5.000	5.000	0.0	85	0.00
27	Benzydine	1.000	1.056	-5.6	92	0.02
28	Pvrene	1.000	1.047	-4.7	97	0.00
29 S	Terphenyl-d14	1.000	1.037	-3.7	98	0.00
30	Benzo(a)anthracene	1.000	0.877	12.3	84	0.00
31	3,3'-Dichlorobenzidine	1.000	1.033	-3.3	96	0.00
32	Chrysene	1.000	1.244	-24.4	116	0.00
33	Bis(2-ethylhexyl)phthalate	1.000	1.105	-10.5	109	0.00
34	Indeno(1,2,3-cd)pyrene	1.000	0.916	8.4	89	0.01
35 I	Perylene-d12	5.000	5.000	0.0	92	0.01
36	Benzo(b)fluoranthene	1.000	0.875	12.5	87	0.00
37	Benzo(k)fluoranthene	1.000	0.971	2.9	101	0.00
38 C	Benzo(a)pyrene	1.000	0.905	9.5	94	0.01
39	Dibenzo(a,h)anthracene	1.000	0.859	14.1	88	0.01
40	Benzo(a,h,i)perylene	1.000	0.859	14.1	89	0.01

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(#) = Out of Range

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