

Data Path : Z:\HPCHEM1\BNA M\DATA\BM010216\
 Data File : BM003922.D
 Acq On : 31 Dec 2015 19:52
 Operator : SJ/UM
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC001

Quant Time: Jan 04 13:59:14 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM010216.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 04 13:33:21 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	93	0.00
2	1,4-Dioxane	0.481	0.450	6.4	95	0.00
3	n-Nitrosodimethylamine	0.517	0.503	2.7	91	0.00
4 S	2-Fluorophenol	1.023	0.911	10.9	88	0.00
5 S	Phenol-d6	1.253	1.162	7.3	88	0.00
6	bis(2-Chloroethyl)ether	1.145	1.136	0.8	91	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	92	0.00
8 S	Nitrobenzene-d5	0.246	0.229	6.9	87	0.00
9	Nitrobenzene	0.266	0.246	7.5	87	0.00
10	Naphthalene	1.094	1.082	1.1	92	-0.02
11	Hexachlorobutadiene	0.167	0.161	3.6	91	0.00
12	2-Methylnaphthalene	0.739	0.716	3.1	90	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	90	0.00
14 S	2,4,6-Tribromophenol	0.116	0.100	13.8	82	0.00
15 S	2-Fluorobiphenyl	1.612	1.618	-0.4	91	0.00
16	Acenaphthylene	2.086	2.035	2.4	87	0.00
17	Acenaphthene	1.457	1.423	2.3	93	0.00
18	Fluorene	1.734	1.763	-1.7	90	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	90	0.00
20	4-Bromophenyl-phenylether	0.155	0.138	11.0	86	0.00
21	Hexachlorobenzene	0.164	0.148	9.8	90	0.00
22	Pentachlorophenol	0.043	0.027	37.2#	72	0.00
23	Phenanthrene	1.037	0.928	10.5	85	0.00
24	Anthracene	1.006	0.910	9.5	88	0.00
25	Fluoranthene	1.113	0.991	11.0	84	0.00
26 I	Chrysene-d12	1.000	1.000	0.0	90	0.00
27	Benzydine	0.418	0.256	38.8#	60	0.00
28	Pvrene	1.796	1.677	6.6	83	0.00
29 S	Terphenyl-d14	0.809	0.742	8.3	83	-0.02
30	Benzo(a)anthracene	1.421	1.411	0.7	92	0.00
31	3,3'-Dichlorobenzidine	0.361	0.258	28.5#	69	0.00
32	Chrysene	1.475	1.176	20.3	71	0.00
33	Bis(2-ethylhexyl)phthalate	0.629	0.311	50.6#	55	0.00
34	Indeno(1,2,3-cd)pyrene	1.057	1.032	2.4	88	0.00
35 I	Perylene-d12	1.000	1.000	0.0	85	0.00
36	Benzo(b)fluoranthene	1.528	1.464	4.2	88	0.00
37	Benzo(k)fluoranthene	1.480	1.329	10.2	79	0.00
38 C	Benzo(a)pyrene	1.301	1.192	8.4	83	0.00
39	Dibenzo(a,h)anthracene	1.110	1.135	-2.3	88	0.00
40	Benzo(a,h,i)perylene	1.176	1.203	-2.3	88	0.00

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(#) = Out of Range

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