

Data Path : Z:\HPCHEM1\BNA M\DATA\BM012216\
 Data File : BM004107.D
 Acq On : 21 Jan 2016 19:52
 Operator : SJ/UM
 Sample : SSTDC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDICC002

Quant Time: Jan 22 01:22:07 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	54	0.00
2	1,4-Dioxane	0.481	0.445	7.5	54	0.00
3	n-Nitrosodimethylamine	0.517	0.457	11.6	48#	-0.02
4 S	2-Fluorophenol	1.023	0.973	4.9	55	-0.02
5 S	Phenol-d6	1.253	1.171	6.5	52	0.00
6	bis(2-Chloroethyl)ether	1.145	1.090	4.8	50	-0.02
7 I	Naphthalene-d8	1.000	1.000	0.0	54	-0.02
8 S	Nitrobenzene-d5	0.246	0.226	8.1	50	-0.01
9	Nitrobenzene	0.266	0.242	9.0	50	-0.01
10	Naphthalene	1.094	1.078	1.5	54	-0.02
11	Hexachlorobutadiene	0.167	0.166	0.6	55	0.00
12	2-Methylnaphthalene	0.739	0.750	-1.5	55	-0.01
13 I	Acenaphthene-d10	1.000	1.000	0.0	60	-0.02
14 S	2,4,6-Tribromophenol	0.116	0.133	-14.7	73	0.00
15 S	2-Fluorobiphenyl	1.612	1.521	5.6	57	-0.01
16	Acenaphthylene	2.086	2.091	-0.2	59	-0.02
17	Acenaphthene	1.457	1.482	-1.7	65	0.00
18	Fluorene	1.734	1.814	-4.6	62	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	63	0.00
20	4-Bromophenyl-phenylether	0.155	0.157	-1.3	68	-0.03
21	Hexachlorobenzene	0.164	0.171	-4.3	73	-0.03
22	Pentachlorophenol	0.043	0.045	-4.7	84	0.00
23	Phenanthrene	1.037	1.017	1.9	65	-0.02
24	Anthracene	1.006	1.116	-10.9	76	0.00
25	Fluoranthene	1.113	1.265	-13.7	75	0.00
26 I	Chrysene-d12	1.000	1.000	0.0	103	0.00
27	Benzydine	0.418	0.195	53.3#	52	0.00
28	Pvrene	1.796	1.412	21.4	81	0.00
29 S	Terphenyl-d14	0.809	0.627	22.5	81	-0.02
30	Benzo(a)anthracene	1.421	1.451	-2.1	108	0.00
31	3,3'-Dichlorobenzidine	0.361	0.340	5.8	105	0.00
32	Chrysene	1.475	1.310	11.2	90	0.00
33	Bis(2-ethylhexyl)phthalate	0.629	0.535	14.9	108	0.00
34	Indeno(1,2,3-cd)pyrene	1.057	1.487	-40.7#	145	-0.01
35 I	Perylene-d12	1.000	1.000	0.0	132	-0.01
36	Benzo(b)fluoranthene	1.528	1.514	0.9	140	0.00
37	Benzo(k)fluoranthene	1.480	1.287	13.0	117	0.00
38 C	Benzo(a)pyrene	1.301	1.293	0.6	139	-0.01
39	Dibenzo(a,h)anthracene	1.110	1.206	-8.6	144	-0.01
40	Benzo(a,h,i)perylene	1.176	1.260	-7.1	143	0.00

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

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