

Data Path : Z:\HPCHEM1\BNA M\DATA\BM012216\
 Data File : BM004115.D
 Acq On : 22 Jan 2016 17:39
 Operator : SJ/UM
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC001

Quant Time: Jan 22 18:17:55 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	96	0.00
2	1,4-Dioxane	0.547	0.465	15.0	97	0.00
3	n-Nitrosodimethylamine	0.446	0.426	4.5	100	-0.02
4 S	2-Fluorophenol	1.043	0.959	8.1	98	-0.02
5 S	Phenol-d6	1.259	1.147	8.9	98	0.00
6	bis(2-Chloroethyl)ether	1.167	1.070	8.3	97	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	98	0.00
8 S	Nitrobenzene-d5	0.219	0.199	9.1	99	0.00
9	Nitrobenzene	0.250	0.226	9.6	100	0.00
10	Naphthalene	1.165	1.074	7.8	98	0.00
11	Hexachlorobutadiene	0.180	0.167	7.2	99	0.00
12	2-Methylnaphthalene	0.762	0.705	7.5	99	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	92	0.00
14 S	2,4,6-Tribromophenol	0.145	0.125	13.8	92	0.00
15 S	2-Fluorobiphenyl	1.532	1.499	2.2	98	0.00
16	Acenaphthylene	2.158	2.019	6.4	96	0.02
17	Acenaphthene	1.634	1.505	7.9	96	0.00
18	Fluorene	1.987	1.911	3.8	99	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	104	0.00
20	4-Bromophenyl-phenylether	0.163	0.131	19.6	86	0.02
21	Hexachlorobenzene	0.171	0.141	17.5	90	0.02
22	Pentachlorophenol	0.043	0.029	32.6#	95	0.00
23	Phenanthrene	1.074	0.918	14.5	97	0.02
24	Anthracene	1.142	0.937	18.0	92	0.00
25	Fluoranthene	1.355	1.156	14.7	96	0.00
26 I	Chrysene-d12	1.000	1.000	0.0	101	0.00
27	Benzydine	0.241	0.122	49.4#	58	0.00
28	Pvrene	1.423	1.213	14.8	95	0.00
29 S	Terphenyl-d14	0.643	0.568	11.7	100	0.00
30	Benzo(a)anthracene	1.518	1.248	17.8	94	0.00
31	3,3'-Dichlorobenzidine	0.338	0.263	22.2	96	-0.01
32	Chrysene	1.372	1.160	15.5	97	-0.01
33	Bis(2-ethylhexyl)phthalate	0.595	0.529	11.1	127	-0.01
34	Indeno(1,2,3-cd)pyrene	1.463	1.263	13.7	99	0.00
35 I	Perylene-d12	1.000	1.000	0.0	97	0.00
36	Benzo(b)fluoranthene	1.551	1.402	9.6	94	0.00
37	Benzo(k)fluoranthene	1.404	1.311	6.6	102	0.00
38 C	Benzo(a)pyrene	1.366	1.258	7.9	100	0.00
39	Dibenzo(a,h)anthracene	1.227	1.125	8.3	99	0.00
40	Benzo(a,h,i)perylene	1.339	1.206	9.9	98	0.00

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(#) = Out of Range

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