

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030516\
 Data File : BM004571.D
 Acq On : 06 Mar 2016 00:07
 Operator : SJ/UM
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC001

Quant Time: Mar 07 08:04:39 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BM030516.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 07 07:54:56 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	110	0.00
2	1,4-Dioxane	0.592	0.484	18.2	109	0.00
3	n-Nitrosodimethylamine	0.385	0.333	13.5	106	0.00
4 S	2-Fluorophenol	1.184	0.982	17.1	101	0.02
5 S	Phenol-d6	1.421	1.194	16.0	102	0.00
6	bis(2-Chloroethyl)ether	1.203	1.033	14.1	102	0.02
7 I	Naphthalene-d8	1.000	1.000	0.0	106	0.00
8 S	Nitrobenzene-d5	0.267	0.228	14.6	101	0.00
9	Nitrobenzene	0.298	0.248	16.8	101	0.00
10	Naphthalene	1.237	1.139	7.9	118	0.01
11	Hexachlorobutadiene	0.216	0.198	8.3	108	0.00
12	2-Methylnaphthalene	0.797	0.705	11.5	105	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	100	-0.02
14 S	2,4,6-Tribromophenol	0.189	0.151	20.1	96	0.00
15 S	2-Fluorobiphenyl	1.639	1.527	6.8	105	0.00
16	Acenaphthylene	2.566	2.289	10.8	101	0.00
17	Acenaphthene	1.547	1.465	5.3	112	-0.02
18	Fluorene	2.075	1.813	12.6	99	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	90	0.00
20	Hexachlorobenzene	0.260	0.225	13.5	98	-0.02
21	Pentachlorophenol	0.027	0.012	55.6#	68	0.00
22	Phenanthrene	1.149	1.205	-4.9	111	-0.02
23	Anthracene	1.134	1.153	-1.7	112	-0.02
24	Fluoranthene	1.460	1.443	1.2	103	0.00
25 I	Chrysene-d12	1.000	1.000	0.0	106	-0.01
26	Benzidine	0.234	0.258	-10.3	161#	0.00
27	Pyrene	1.868	1.538	17.7	103	0.00
28 S	Terphenyl-d14	0.788	0.645	18.1	104	0.00
29	Benzo(a)anthracene	1.620	1.386	14.4	104	0.00
30	3,3'-Dichlorobenzidine	0.500	0.444	11.2	99	0.01
31	Chrysene	1.689	1.390	17.7	103	-0.01
32	Bis(2-ethylhexyl)phthalate	1.215	0.970	20.2	105	-0.01
33	Indeno(1,2,3-cd)pyrene	1.555	1.292	16.9	104	0.00
34 I	Perylene-d12	1.000	1.000	0.0	107	0.00
35 C	Benzo(a)pyrene	1.534	1.387	9.6	106	0.00
36	Dibenzo(a,h)anthracene	1.278	1.096	14.2	105	0.00
37	Benzo(g,h,i)perylene	1.412	1.225	13.2	105	0.00

(#) = Out of Range

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