

Data Path : S:\HPCHEM1\BNA_M\DATA\BM022916\
 Data File : BM004500.D
 Acq On : 01 Mar 2016 10:29
 Operator : SJ/UM
 Sample : SSTDC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDC001

Quant Time: Mar 01 11:57:37 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BM022916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 01 00:52:28 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	90	0.00
2	1,4-Dioxane	0.580	0.491	15.3	90	0.00
3	n-Nitrosodimethylamine	0.508	0.462	9.1	88	0.00
4 S	2-Fluorophenol	1.162	1.065	8.3	89	0.02
5 S	Phenol-d6	1.313	1.295	1.4	94	0.02
6	bis(2-Chloroethyl)ether	1.389	1.371	1.3	92	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	94	0.00
8 S	Nitrobenzene-d5	0.254	0.239	5.9	93	0.00
9	Nitrobenzene	0.344	0.319	7.3	91	0.00
10	Naphthalene	1.572	1.385	11.9	95	0.00
11	Hexachlorobutadiene	0.230	0.230	0.0	96	-0.01
12	2-Methylnaphthalene	0.869	0.892	-2.6	101	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	99	0.00
14 S	2,4,6-Tribromophenol	0.192	0.171	10.9	99	0.00
15 S	2-Fluorobiphenyl	1.531	1.520	0.7	104	0.00
16	Acenaphthylene	2.861	2.716	5.1	99	0.00
17	Acenaphthene	1.872	1.894	-1.2	112	0.00
18	Fluorene	2.285	2.354	-3.0	108	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	89	0.00
20	4-Bromophenyl-phenylether	0.274	0.290	-5.8	123	0.00
21	Hexachlorobenzene	0.308	0.344	-11.7	128	0.00
22	Pentachlorophenol	0.074	0.062	16.2	132	0.00
23	Phenanthrene	1.762	1.859	-5.5	120	0.00
24	Anthracene	1.485	1.456	2.0	100	0.02
25	Fluoranthene	1.891	1.964	-3.9	109	0.00
26 I	Chrysene-d12	1.000	1.000	0.0	84	0.01
27	Benzydine	0.535	0.804	-50.3#	161#	0.00
28	Pyrene	1.977	2.209	-11.7	108	0.00
29 S	Terphenyl-d14	0.643	0.714	-11.0	107	0.00
30	Benzo(a)anthracene	1.845	1.771	4.0	91	0.01
31	3,3'-Dichlorobenzidine	1.118	1.686	-50.8#	161#	0.00
32	Chrysene	3.757	2.269	39.6#	89	0.00
33	Bis(2-ethylhexyl)phthalate	1.432	1.190	16.9	67	-0.01
34	Indeno(1,2,3-cd)pyrene	1.712	1.860	-8.6	102	-0.01
35 I	Perylene-d12	1.000	1.000	0.0	100	0.00
36	Benzo(b)fluoranthene	1.949	2.032	-4.3	104	0.00
37	Benzo(k)fluoranthene	1.920	1.900	1.0	104	0.00
38 C	Benzo(a)pyrene	1.761	1.781	-1.1	103	0.00
39	Dibenzo(a,h)anthracene	1.533	1.509	1.6	102	0.00
40	Benzo(g,h,i)perylene	1.683	1.662	1.2	102	-0.01

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