

Data Path : S:\HPCHEM1\BNA_M\DATA\BM022916\
 Data File : BM004492.D
 Acq On : 01 Mar 2016 05:42
 Operator : SJ/UM
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC001

Quant Time: Mar 01 08:01:47 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	94	0.00
2	1,4-Dioxane	0.580	0.469	19.1	90	0.00
3	n-Nitrosodimethylamine	0.508	0.436	14.2	86	0.00
4 S	2-Fluorophenol	1.162	1.020	12.2	88	0.00
5 S	Phenol-d6	1.313	1.200	8.6	91	0.02
6	bis(2-Chloroethyl)ether	1.389	1.296	6.7	91	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	95	0.00
8 S	Nitrobenzene-d5	0.254	0.227	10.6	90	0.00
9	Nitrobenzene	0.344	0.307	10.8	89	0.00
10	Naphthalene	1.572	1.339	14.8	93	0.00
11	Hexachlorobutadiene	0.230	0.224	2.6	94	-0.01
12	2-Methylnaphthalene	0.869	0.839	3.5	96	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	100	0.00
14 S	2,4,6-Tribromophenol	0.192	0.151	21.4	88	0.00
15 S	2-Fluorobiphenyl	1.531	1.419	7.3	97	0.00
16	Acenaphthylene	2.861	2.687	6.1	99	0.00
17	Acenaphthene	1.872	1.664	11.1	99	0.00
18	Fluorene	2.285	2.147	6.0	99	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	104	0.00
20	4-Bromophenyl-phenylether	0.274	0.199	27.4#	99	0.00
21	Hexachlorobenzene	0.308	0.224	27.3#	97	0.00
22	Pentachlorophenol	0.074	0.050	32.4#	123	0.00
23	Phenanthrene	1.762	1.314	25.4#	99	0.00
24	Anthracene	1.485	1.251	15.8	100	0.00
25	Fluoranthene	1.891	1.571	16.9	101	-0.02
26 I	Chrysene-d12	1.000	1.000	0.0	98	0.00
27	Benzydine	0.535	0.383	28.4#	89	0.00
28	Pyrene	1.977	1.782	9.9	101	0.00
29 S	Terphenyl-d14	0.643	0.556	13.5	97	0.00
30	Benzo(a)anthracene	1.845	1.653	10.4	98	0.00
31	3,3'-Dichlorobenzidine	1.118	0.894	20.0	99	0.00
32	Chrysene	3.757	1.889	49.7#	86	0.00
33	Bis(2-ethylhexyl)phthalate	1.432	1.319	7.9	86	0.00
34	Indeno(1,2,3-cd)pyrene	1.712	1.597	6.7	102	-0.01
35 I	Perylene-d12	1.000	1.000	0.0	103	0.00
36	Benzo(b)fluoranthene	1.949	1.771	9.1	94	0.00
37	Benzo(k)fluoranthene	1.920	1.831	4.6	103	0.00
38 C	Benzo(a)pyrene	1.761	1.717	2.5	103	0.00
39	Dibenzo(a,h)anthracene	1.533	1.462	4.6	102	0.00
40	Benzo(g,h,i)perylene	1.683	1.626	3.4	103	-0.01

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