

Data Path : S:\HPCHEM1\BNA_M\DATA\BM022916\
 Data File : BM004490.D
 Acq On : 01 Mar 2016 03:47
 Operator : SJ/UM
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC001

Quant Time: Mar 01 07:29:49 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	84	0.00
2	1,4-Dioxane	0.580	0.467	19.5	80	0.00
3	n-Nitrosodimethylamine	0.508	0.443	12.8	78	0.00
4 S	2-Fluorophenol	1.162	1.019	12.3	79	0.00
5 S	Phenol-d6	1.313	1.237	5.8	83	0.02
6	bis(2-Chloroethyl)ether	1.389	1.299	6.5	81	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	88	0.00
8 S	Nitrobenzene-d5	0.254	0.232	8.7	85	0.00
9	Nitrobenzene	0.344	0.311	9.6	83	0.00
10	Naphthalene	1.572	1.340	14.8	86	0.00
11	Hexachlorobutadiene	0.230	0.222	3.5	87	-0.01
12	2-Methylnaphthalene	0.869	0.864	0.6	92	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	92	0.00
14 S	2,4,6-Tribromophenol	0.192	0.166	13.5	89	0.00
15 S	2-Fluorobiphenyl	1.531	1.475	3.7	93	0.00
16	Acenaphthylene	2.861	2.699	5.7	91	0.00
17	Acenaphthene	1.872	1.819	2.8	100	0.00
18	Fluorene	2.285	2.297	-0.5	98	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	87	0.00
20	4-Bromophenyl-phenylether	0.274	0.264	3.6	109	0.00
21	Hexachlorobenzene	0.308	0.316	-2.6	114	0.00
22	Pentachlorophenol	0.074	0.059	20.3	122	0.00
23	Phenanthrene	1.762	1.688	4.2	106	0.00
24	Anthracene	1.485	1.317	11.3	88	0.02
25	Fluoranthene	1.891	1.810	4.3	98	0.00
26 I	Chrysene-d12	1.000	1.000	0.0	79	-0.01
27	Benزيدine	0.535	0.383	28.4#	72	0.00
28	Pyrene	1.977	2.130	-7.7	98	0.00
29 S	Terphenyl-d14	0.643	0.670	-4.2	95	0.00
30	Benzo(a)anthracene	1.845	1.779	3.6	86	-0.01
31	3,3'-Dichlorobenzidine	1.118	1.103	1.3	99	0.00
32	Chrysene	3.757	2.040	45.7#	75	0.00
33	Bis(2-ethylhexyl)phthalate	1.432	1.302	9.1	69	-0.01
34	Indeno(1,2,3-cd)pyrene	1.712	1.783	-4.1	92	-0.01
35 I	Perylene-d12	1.000	1.000	0.0	93	-0.01
36	Benzo(b)fluoranthene	1.949	2.216	-13.7	107	-0.01
37	Benzo(k)fluoranthene	1.920	1.940	-1.0	99	0.00
38 C	Benzo(a)pyrene	1.761	1.743	1.0	95	0.00
39	Dibenzo(a,h)anthracene	1.533	1.465	4.4	93	-0.01
40	Benzo(g,h,i)perylene	1.683	1.615	4.0	93	-0.01

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