

Data Path : Z:\HPCHEM1\BNA M\DATA\BM013016\
 Data File : BM004156.D
 Acq On : 29 Jan 2016 21:01
 Operator : SJ/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC001

Quant Time: Jan 29 22:16:02 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	74	0.00
2	1,4-Dioxane	0.547	0.467	14.6	75	0.00
3	n-Nitrosodimethylamine	0.446	0.419	6.1	76	0.00
4 S	2-Fluorophenol	1.043	1.019	2.3	80	0.00
5 S	Phenol-d6	1.259	1.259	0.0	83	0.00
6	bis(2-Chloroethyl)ether	1.167	1.071	8.2	75	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	77	0.00
8 S	Nitrobenzene-d5	0.219	0.214	2.3	84	0.00
9	Nitrobenzene	0.250	0.240	4.0	83	0.00
10	Naphthalene	1.165	1.060	9.0	76	0.00
11	Hexachlorobutadiene	0.180	0.165	8.3	77	0.00
12	2-Methylnaphthalene	0.762	0.714	6.3	79	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	82	0.00
14 S	2,4,6-Tribromophenol	0.145	0.152	-4.8	98	0.00
15 S	2-Fluorobiphenyl	1.532	1.379	10.0	80	0.00
16	Acenaphthylene	2.158	1.983	8.1	83	0.00
17	Acenaphthene	1.634	1.427	12.7	80	0.00
18	Fluorene	1.987	1.866	6.1	86	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	88	0.00
20	4-Bromophenyl-phenylether	0.163	0.146	10.4	81	0.00
21	Hexachlorobenzene	0.171	0.150	12.3	82	0.00
22	Pentachlorophenol	0.043	0.052	-20.9	143	0.00
23	Phenanthrene	1.074	0.952	11.4	85	0.00
24	Anthracene	1.142	1.053	7.8	87	0.00
25	Fluoranthene	1.355	1.307	3.5	92	0.00
26 I	Chrysene-d12	1.000	1.000	0.0	83	0.00
27	Benzydine	0.241	0.241	0.0	94	0.02
28	Pvrene	1.423	1.399	1.7	90	0.00
29 S	Terphenyl-d14	0.643	0.631	1.9	91	0.00
30	Benzo(a)anthracene	1.518	1.359	10.5	83	0.00
31	3,3'-Dichlorobenzidine	0.338	0.421	-24.6	126	0.00
32	Chrysene	1.372	1.559	-13.6	107	0.00
33	Bis(2-ethylhexyl)phthalate	0.595	0.765	-28.6#	151#	0.00
34	Indeno(1,2,3-cd)pyrene	1.463	1.553	-6.2	99	0.01
35 I	Perylene-d12	1.000	1.000	0.0	96	0.00
36	Benzo(b)fluoranthene	1.551	1.356	12.6	90	0.00
37	Benzo(k)fluoranthene	1.404	1.381	1.6	107	0.00
38 C	Benzo(a)pyrene	1.366	1.296	5.1	102	0.00
39	Dibenzo(a,h)anthracene	1.227	1.141	7.0	99	0.00
40	Benzo(a,h,i)perylene	1.339	1.223	8.7	98	0.00

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