

Data Path : Z:\HPCHEM1\BNA M\DATA\BM012216\
 Data File : BM004110.D
 Acq On : 22 Jan 2016 14:39
 Operator : SJ/UM
 Sample : SSTDICV001
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ICVBM012216

Quant Time: Jan 22 15:17:53 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	107	0.00
2	1,4-Dioxane	0.547	0.441	19.4	102	0.00
3	n-Nitrosodimethylamine	0.446	0.400	10.3	105	-0.02
4 S	2-Fluorophenol	1.043	0.915	12.3	104	0.00
5 S	Phenol-d6	1.259	1.096	12.9	104	0.00
6	bis(2-Chloroethyl)ether	1.167	1.023	12.3	104	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	106	0.00
8 S	Nitrobenzene-d5	0.219	0.198	9.6	106	0.00
9	Nitrobenzene	0.250	0.222	11.2	106	0.00
10	Naphthalene	1.165	1.029	11.7	102	0.00
11	Hexachlorobutadiene	0.180	0.161	10.6	104	0.00
12	2-Methylnaphthalene	0.762	0.670	12.1	102	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	107	0.00
14 S	2,4,6-Tribromophenol	0.145	0.120	17.2	101	0.00
15 S	2-Fluorobiphenyl	1.532	1.343	12.3	102	0.00
16	Acenaphthylene	2.158	1.872	13.3	102	0.00
17	Acenaphthene	1.634	1.374	15.9	101	0.00
18	Fluorene	1.987	1.705	14.2	102	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	106	0.00
20	4-Bromophenyl-phenylether	0.163	0.150	8.0	100	0.00
21	Hexachlorobenzene	0.171	0.158	7.6	103	0.00
22	Pentachlorophenol	0.043	0.033	23.3	109	0.00
23	Phenanthrene	1.074	0.951	11.5	102	0.00
24	Anthracene	1.142	1.027	10.1	102	0.00
25	Fluoranthene	1.355	1.212	10.6	102	0.00
26 I	Chrysene-d12	1.000	1.000	0.0	100	0.00
27	Benzydine	0.241	0.244	-1.2	115	0.00
28	Pvrene	1.423	1.311	7.9	102	0.00
29 S	Terphenyl-d14	0.643	0.596	7.3	104	0.00
30	Benzo(a)anthracene	1.518	1.316	13.3	98	0.00
31	3,3'-Dichlorobenzidine	0.338	0.314	7.1	113	0.00
32	Chrysene	1.372	1.379	-0.5	114	0.00
33	Bis(2-ethylhexyl)phthalate	0.595	0.497	16.5	119	0.00
34	Indeno(1,2,3-cd)pyrene	1.463	1.360	7.0	105	0.00
35 I	Perylene-d12	1.000	1.000	0.0	109	0.00
36	Benzo(b)fluoranthene	1.551	1.334	14.0	101	0.00
37	Benzo(k)fluoranthene	1.404	1.245	11.3	109	0.00
38 C	Benzo(a)pyrene	1.366	1.195	12.5	107	0.00
39	Dibenzo(a,h)anthracene	1.227	1.065	13.2	105	0.00
40	Benzo(a,h,i)perylene	1.339	1.140	14.9	104	0.00

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(#) = Out of Range

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