

Data Path : Z:\HPCHEM1\BNA M\DATA\BM090116\
 Data File : BM007363.D
 Acq On : 01 Sep 2016 17:41
 Operator : UM/SJ
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICV0.4

Quant Time: Sep 02 12:53:36 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM090116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 02 12:51:17 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	91	0.00
2	1,4-Dioxane	0.443	0.404	8.8	81	0.00
3	n-Nitrosodimethylamine	0.476	0.496	-4.2	121	0.00
4 S	2-Fluorophenol	1.032	1.015	1.6	98	0.00
5 S	Phenol-d6	1.400	1.251	10.6	98	0.00
6	bis(2-Chloroethyl)ether	1.086	1.026	5.5	97	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	91	0.00
8 S	Nitrobenzene-d5	0.281	0.237	15.7	99	0.00
9	Nitrobenzene	0.269	0.252	6.3	101	0.00
10	Naphthalene	1.144	1.132	1.0	95	0.00
11	Hexachlorobutadiene	0.209	0.213	-1.9	98	0.00
12	2-Methylnaphthalene	0.817	0.800	2.1	95	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	89	0.00
14 S	2,4,6-Tribromophenol	0.194	0.172	11.3	84	0.00
15 S	2-Fluorobiphenyl	1.553	1.511	2.7	96	0.00
16	Acenaphthylene	2.164	2.016	6.8	96	0.00
17	Acenaphthene	1.495	1.436	3.9	95	0.00
18	Fluorene	1.757	1.749	0.5	98	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	89	0.00
20	4-Bromophenyl-phenylether	0.202	0.186	7.9	96	0.00
21	Hexachlorobenzene	0.256	0.244	4.7	92	0.00
22	Pentachlorophenol	0.079	0.077	2.5	100	0.00
23	Phenanthrene	1.276	1.254	1.7	96	0.00
24	Anthracene	1.283	1.189	7.3	93	0.00
25	Fluoranthene	1.502	1.410	6.1	97	0.00
26 I	Chrysene-d12	1.000	1.000	0.0	95	0.00
27	Benzydine	0.324	0.299	7.7	102	0.02
28	Pvrene	1.425	1.341	5.9	94	0.00
29 S	Terphenyl-d14	0.671	0.623	7.2	95	0.00
30	Benzo(a)anthracene	1.446	1.367	5.5	101	0.00
31	3,3'-Dichlorobenzidine	0.400	0.361	9.8	98	0.00
32	Chrysene	1.408	1.365	3.1	94	0.00
33	Bis(2-ethylhexyl)phthalate	0.753	0.699	7.2	93	0.00
34	Indeno(1,2,3-cd)pyrene	1.237	1.119	9.5	98	0.00
35 I	Perylene-d12	1.000	1.000	0.0	88	0.01
36	Benzo(b)fluoranthene	1.559	1.516	2.8	92	0.00
37	Benzo(k)fluoranthene	1.438	1.465	-1.9	102	0.00
38 C	Benzo(a)pyrene	1.385	1.354	2.2	97	0.01
39	Dibenzo(a,h)anthracene	1.123	1.104	1.7	100	0.01
40	Benzo(a,h,i)perylene	1.165	1.153	1.0	98	0.00

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