

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE051916\
 Data File : BE091809.D
 Acq On : 19 May 2016 14:26
 Operator : UM/SJ
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE051916

Quant Time: May 19 15:43:39 2016
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE051916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 19 15:36:49 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	99	-0.02
2	1,4-Dioxane	0.649	0.646	0.5	93	0.00
3	n-Nitrosodimethylamine	0.798	0.708	11.3	96	0.00
4 S	2-Fluorophenol	1.307	1.175	10.1	95	0.00
5 S	Phenol-d6	1.840	1.563	15.1	94	0.00
6	bis(2-Chloroethyl)ether	1.477	1.307	11.5	96	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	97	0.00
8 S	Nitrobenzene-d5	0.331	0.286	13.6	94	0.00
9	Nitrobenzene	0.341	0.283	17.0	93	0.00
10	Naphthalene	1.106	1.049	5.2	96	0.00
11	Hexachlorobutadiene	0.164	0.158	3.7	97	-0.02
12	2-Methylnaphthalene	0.744	0.690	7.3	95	-0.01
13 I	Acenaphthene-d10	1.000	1.000	0.0	95	-0.01
14 S	2,4,6-Tribromophenol	0.194	0.160	17.5	91	-0.01
15 S	2-Fluorobiphenyl	1.414	1.336	5.5	95	0.00
16	Acenaphthylene	2.230	2.066	7.4	93	0.00
17	Acenaphthene	1.341	1.229	8.4	93	-0.01
18	Fluorene	1.738	1.631	6.2	96	-0.01
19 I	Phenanthrene-d10	1.000	1.000	0.0	99	-0.01
20	4-Bromophenyl-phenylether	0.187	0.178	4.8	96	0.00
21	Hexachlorobenzene	0.195	0.188	3.6	97	0.00
22	Pentachlorophenol	0.080	0.064	20.0	87	0.00
23	Phenanthrene	1.190	1.176	1.2	97	-0.01
24	Anthracene	1.088	1.010	7.2	95	-0.01
25	Fluoranthene	1.329	1.258	5.3	94	0.00
26 I	Chrysene-d12	1.000	1.000	0.0	101	-0.02
27	Benzydine	0.338	0.243	28.1#	95	0.00
28	Pyrene	1.097	0.984	10.3	95	0.00
29 S	Terphenyl-d14	0.708	0.662	6.5	101	-0.02
30	Benzo(a)anthracene	1.206	1.130	6.3	101	-0.02
31	3,3'-Dichlorobenzidine	0.767	0.606	21.0	91	-0.02
32	Chrysene	1.101	1.033	6.2	102	0.00
33	Bis(2-ethylhexyl)phthalate	0.540	0.372	31.1#	88	-0.02
34	Indeno(1,2,3-cd)pyrene	1.236	1.102	10.8	95	-0.03
35 I	Perylene-d12	1.000	1.000	0.0	99	-0.02
36	Benzo(b)fluoranthene	1.251	1.137	9.1	95	-0.02
37	Benzo(k)fluoranthene	1.281	1.180	7.9	96	-0.02
38 C	Benzo(a)pyrene	1.194	1.080	9.5	95	-0.02
39	Dibenzo(a,h)anthracene	1.166	1.052	9.8	95	-0.05
40	Benzo(g,h,i)perylene	1.203	1.108	7.9	96	-0.05

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(#) = Out of Range

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