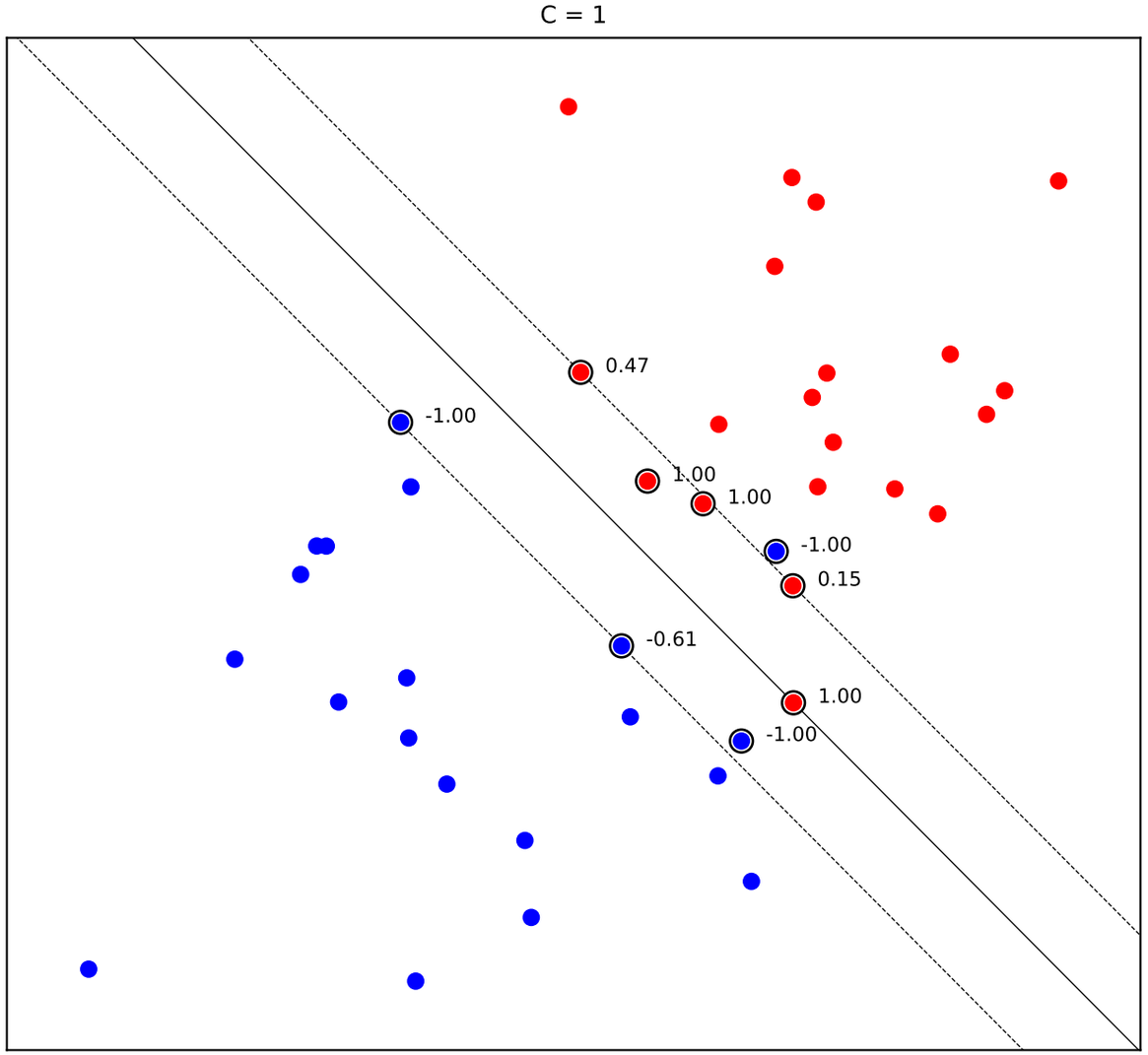
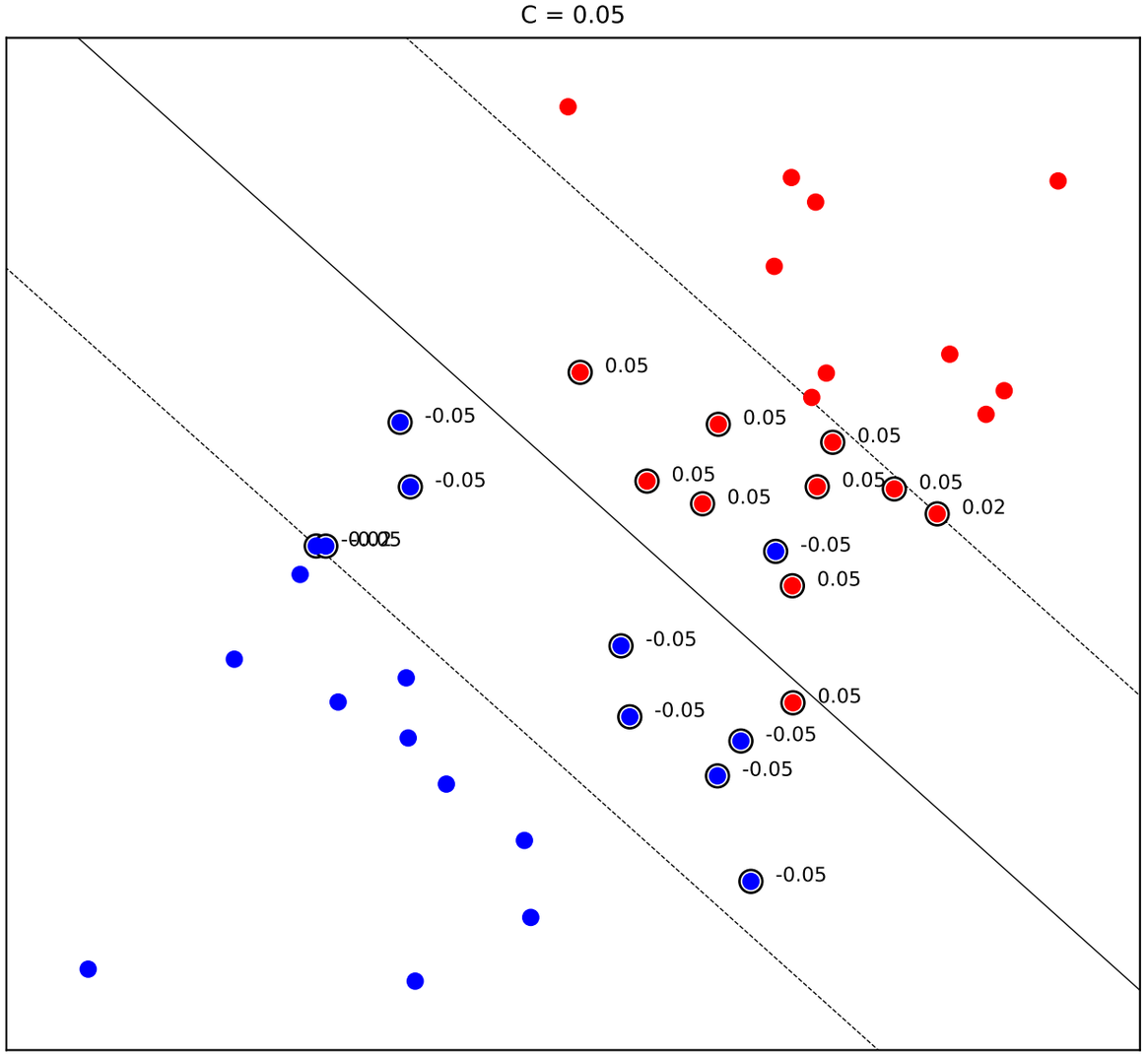

C = 1
0.47
-1.00
1.00
1.00
-1.00
0.15
-0.61
1.00
-1.00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0.05
-0.05
0.05
0.05
-0.05
0.05
0.05
0.05
0.05
0.02
-0.05
0.05
-0.05
0.05
-0.05
-0.05
-0.05
-0.05