Machine Learning and Artificial Intelligence 2018/2019 Politecnico di Torino Prof. Barbara Caputo Due Date: 19/01/2019

1 Linear SVM

In the first part of this homework I trained, validated and finally tested a Linear SVM, using the C with the highest accuracy in validation. The essential part of the code is algorithm 1.

Algorithm 1: Searching the best value of C in Linear SVM

```
acc = np.empty(7)
78
        c_i = 1e-3
79
        c_best = 1e-3
80
       a_best = 0
81
       i = 1
82
       xx, yy = make_meshgrid(x[:, 0], x[:, 1])
       while c_i <= 1e3:
84
           clf = svm.LinearSVC(C=c_i, random_state=r_state)
85
           acc[i - 1] = clf.fit(x_train, y_train).score(x_val, y_val) * 100
86
           if acc[i - 1] > a_best:
               c_best = c_i
88
               a_best = acc[i - 1]
89
           ax = fig.add_subplot(4, 2, i)
90
           plot_data(ax, x_train, y_train, xx, yy, clf)
91
           ax.set_xlim(xx.min(), xx.max())
92
           ax.set_ylim(yy.min(), yy.max())
93
           ax.set_xlabel('Sepal length')
94
           ax.set_ylabel('Sepal width')
95
           ax.set_xticks(())
96
           ax.set_yticks(())
97
           ax.legend()
           ax.set_title('C=%2.2E A=%2.1f%%' % (c_i, acc[i - 1]))
99
           c_i = c_i * 10
100
           i = i + 1
```

The plots in figs. 1 and 2 on the next page and on page 3, provide information about one run of the code: the best accuracy found on validation set was 73.33%, with C equal to 1e+1. Clearly, that boundaries become more and more precise on the training set with the increment of C.

C is the hyper parameter of SVM that represents the penalty for misclassifying a data point, so it describes how much SVM has to avoid misclassification during the training. When C is large, the optimization will choose a smaller-margin hyperplane, because the classifier is heavily penalized for misclassified data. The larger is C, the better the classification on training set will be. On the contrary, when C is small, the optimizer will use a larger-margin separating hyperplane, causing the misclassification of more training data points.

After this training, I tested the data on the test set, obtaining a greater accuracy (88.89%), as shown in fig. 3 on page 3. I tried to repeat the code 500 times, producing validation accuracies of $80.03\% \pm 6.62\%$ and testing ones of $73.52\% \pm 7.11\%$ (see section 4 on page 9). The variances are high and this may be due to the initial state of the algorithm of SVM or the splitting of the data in training, validation and test sets: maybe in some cases there is more overfitting on the training data, whereas less in others. To sum up, this type of validation is not stable and may be influenced by overfitting.

Linear SVM - C tuning - C_best = 1.00E+01 A_best = 73.3%

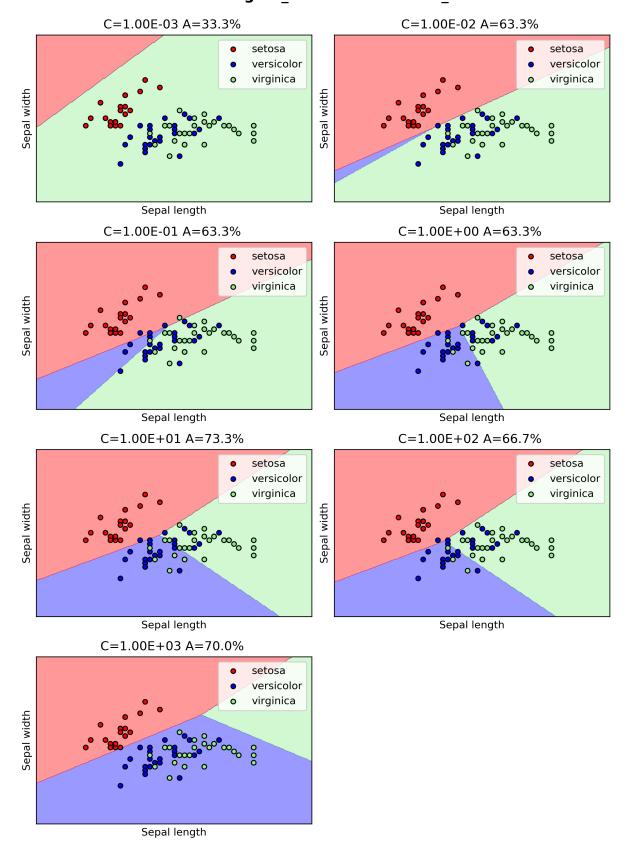


Figure 1: Decision Boundaries changing C in Linear SVM on Training set

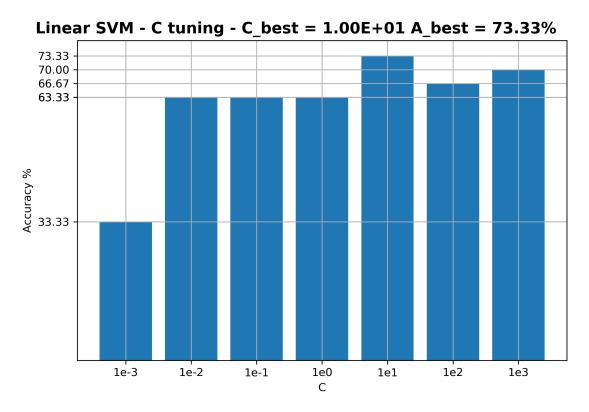


Figure 2: Accuracy changing C in Linear SVM

Linear SVM with validation accuracy of 73.33% $_{\text{C}=1.00\text{E}+01}$ A=88.89%

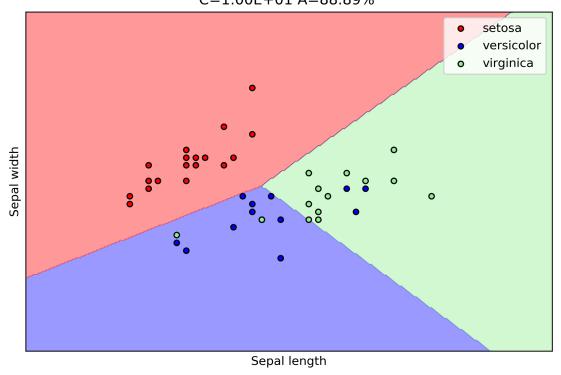


Figure 3: Results on the test set

2 RBF Kernel

In the second part of this homework, I used SVM with Gaussian RBF kernel eq. (1).

$$exp(-\gamma \mid\mid x - x'\mid\mid^2) \tag{1}$$

This time I had to find the best pair of C (explained in section 1 on page 1) and γ (Gamma).

Gamma parameter can be explained as the *spread* of the kernel, therefore of the decision region. When gamma is **low**, the *curve* of the decision boundary is very low, consequently the decision region is very broad. When gamma is **high**, the *curve* of the decision boundary is high and decision-*islands* will arise around data points.

Firstly, I repeated the steps of the previous section in order to find the best accuracy modifying C and using the Gamma calculated by sklearn under the hood (see algorithm 2). This time the boundaries are **not as linear as** the ones of Linear SVM, because of the introduction of the Gaussian RBF kernel. Thanks to this approach, I found out, in the first run, a better training accuracy (76.67%) compared to the Linear SVM one, and a test accuracy near to the validation one (75.56%) with C = 1e-1, as shown in figs. 4 to 6 on pages 5–6.

I used the code of algorithm 1 on page 1, changing the classifier on line 88 with the one in algorithm 2 .

Algorithm 2: RBF Kernel SVM Classifier

```
clf = svm.SVC(kernel='rbf', C=c_i, random_state=r_state)
```

Finally, I performed the grid search of the best pair of C and Gamma with algorithm 3. In the first run I obtained an high accuracy in validation (80.00%) and also on the test set (86.67%) with a C = 1e0 and $\gamma = 1e-1$, as we can see in figs. 7 and 8 on page 7.

Algorithm 3: Searching the best value of C and Gamma in RBF Kernel SVM

```
# Grid Search of C and Gamma
211
212
        c = np.array([1e-3, 1e-2, 1e-1, 1e0, 1e1, 1e2, 1e3])
        gamma = np.array(
213
            [1e-9, 1e-8, 1e-7, 1e-6, 1e-5, 1e-4, 1e-3, 1e-2, 1e-1, 1e0, 1e1, 1e2, 1e3,
                1e4, 1e5, 1e6, 1e7, 1e8, 1e9])
        c_best = c.min()
215
        g_best = gamma.min()
216
        a_best = 0
217
       res = np.zeros([c.shape[0], gamma.shape[0]])
218
        for c_i, i in zip(c, range(0, c.shape[0])):
219
           for gamma_i, j in zip(gamma, range(0, gamma.shape[0])):
               clf = svm.SVC(kernel='rbf', gamma=gamma_i, C=c_i, random_state=r_state)
               res[i, j] = clf.fit(x_train, y_train).score(x_val, y_val) * 100
               if res[i, j] > a_best:
223
                   c_best = c_i
224
                   g_best = gamma_i
225
                   a_best = res[i, j]
```

I repeated this code 500 times and, even then, I found out an high variance in validation $(82.11\% \pm 6.13\%)$ and test $(76.86\% \pm 6.23\%)$ scores (see section 4 on page 9).

RBF Kernel - C/G tuning - C_best=1.00E-01 A_best=76.67%

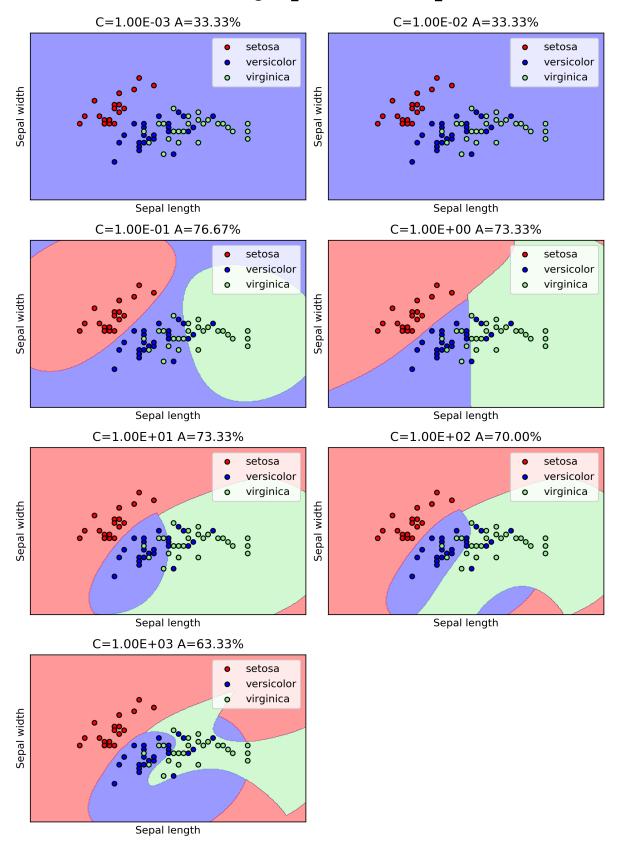


Figure 4: Decision Boundaries changing C in RBF Kernel SVM on Training set

RBF Kernel - C/G tuning - C_best=1.00E-01 A_best=76.67%

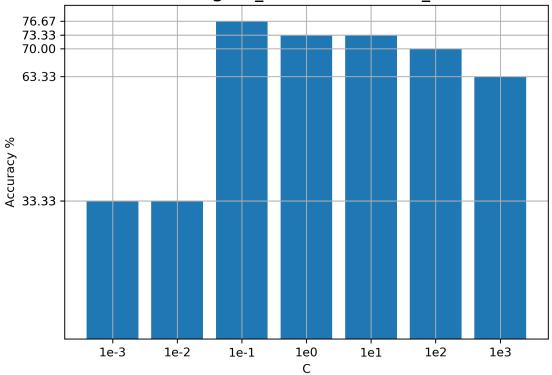


Figure 5: Accuracy changing C in RBF Kernel SVM

RBF Kernel with validation accuracy of 76.67% $_{\text{C}=1.00\text{E-}01}$ $_{\text{A}=75.56\%}$

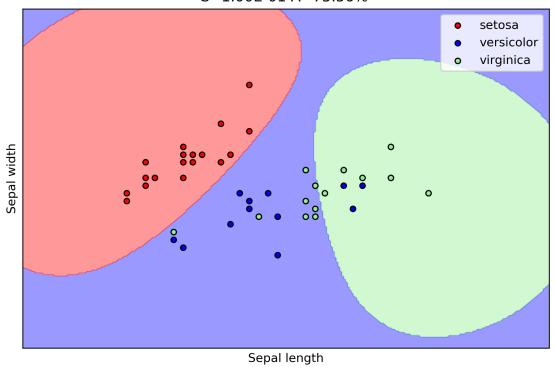


Figure 6: Results on the test set

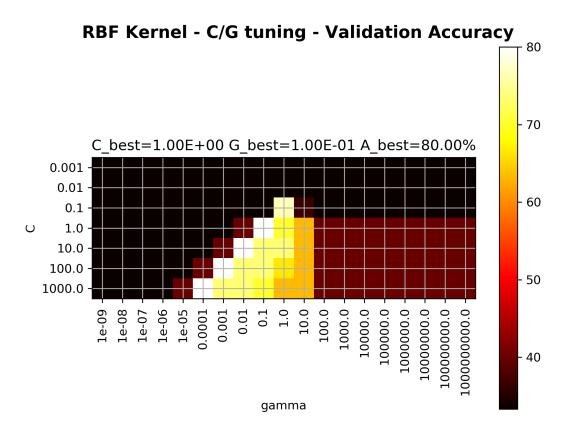


Figure 7: Grid Search of C and Gamma in RBF Kernel SVM on Training set

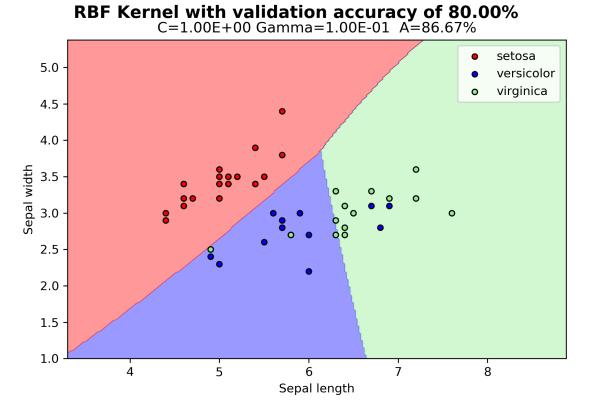


Figure 8: Results on the test set

3 K-Fold

In the last part of the homework, I merged training and validation sets and I performed a **k-fold** crossvalidation with k = 5 using algorithm 4.

Algorithm 4: Validation using k-fold with k = 5

```
c_best = c.min()
265
       g_best = gamma.min()
266
       a_best = 0
267
       k = 5
268
       kf = KFold(n_splits=k, shuffle=True, random_state=r_state)
       res = np.zeros([c.shape[0], gamma.shape[0]])
        for i, c_i in enumerate(c):
           for j, gamma_i in enumerate(gamma):
272
               temp = np.zeros(kf.n_splits)
               for k_i, (train_index, test_index) in enumerate(kf.split(x_train)):
                   clf = svm.SVC(kernel='rbf', gamma=gamma_i, C=c_i)
                   clf.fit(x_train[train_index], y_train[train_index])
                   temp[k_i] = clf.score(x_train[test_index], y_train[test_index]) * 100
               res[i, j] = np.average(temp)
278
               if res[i, j] > a_best:
                   c_best = c_i
280
                   g_best = gamma_i
281
                   a_best = res[i, j]
282
```

In the first run, I obtained a validation accuracy equal to 76.19%, but a test accuracy of 82.22% with a C = 1e0 and $\gamma = 1e-1$ (see figs. 9 and 10 on this page and on the next page).

The final score is satisfying, although it is lower than some of the final scores found in previous sections. I tried to run the code 500 times obtaining a validation accuracy of $81.17\% \pm 2.44\%$ and a testing one of $78.14\% \pm 5.38\%$ (see section 4 on the following page).

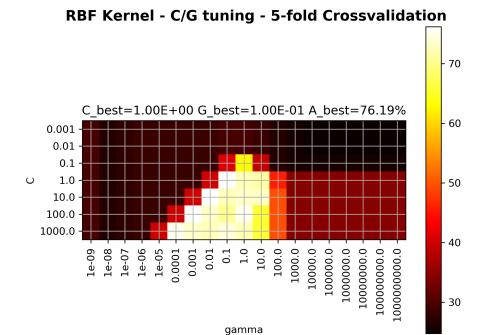


Figure 9: Grid Search of C and Gamma in RBF Kernel SVM with 5-fold crossvalidation



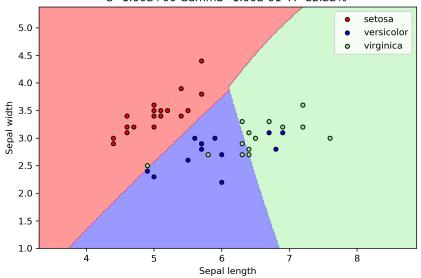


Figure 10: Results on the test set

4 Final Comments

I decided to implement a simple Python script (testing.py) in order to calculate the mean and the variance of the results and to to better evaluate the three types of parameter tuning algorithm explained in this report . The result obtained over 500 repetitions can be found in table 1.

Table	1:	Accuracy	dis	tri	but	ion
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Accuracy (%) (500 rep.)	AVG Validation	STD Validation	AVG Test	STD Test
Linear SVM	80.03	6.62	73.52	7.11
RBF Kernel SVM	82.11	6.13	76.86	6.23
RBF Kernel SVM (k-fold)	81.17	2.44	78.14	5.38

At this point, the data gathered are definitely noteworthy: the first two type of parameter tuning algorithms have a larger variance than the last one. Therefore, the validation averages are almost the same, but not the test ones. Indeed, we can find an higher test accuracy on the third type.

In conclusion, taking all the results into account, the validation done with k-fold crossvalidation is more stable than the other ones and it is useful to avoid overfitting problem.

5 Code Execution

5.1 Requirements

- Python 3
- All dependencies in requirements.txt.
 - \$ pip install -r requirements.txt to install them

5.2 Usage

• \$ python main.py
Execute the code

5.3 Reproducibility

In order to reproduce the same data for this experiment you have to change the global variable r_state (line) from None to 252894 which is my badge number.

Attachments

- source_code folder:
 - main.py
 - requirements.txt
 - testing.py Code used for testing in final comments