03c DEMO SVM

May 15, 2022

1 Machine Learning Foundation

1.1 Course 3, Part c: Support Vector Machines DEMO

1.2 Introduction

We will be using the wine quality data set for these exercises. This data set contains various chemical properties of wine, such as acidity, sugar, pH, and alcohol. It also contains a quality metric (3-9, with highest being better) and a color (red or white). The name of the file is Wine_Quality_Data.csv.

```
[1]: def warn(*args, **kwargs):
    pass
import warnings
warnings.warn = warn
import numpy as np, pandas as pd, matplotlib.pyplot as plt, seaborn as sns
```

1.3 Part 1: Setup

- Import the data.
- $\bullet\,$ Create the target variable y as a 1/0 column where 1 means red.
- Create a pairplot for the dataset.
- Create a bar plot showing the correlations between each column and y
- Pick the most 2 correlated fields (using the absolute value of correlations) and create X
- Use MinMaxScaler to scale X. Note that this will output a np.array. Make it a DataFrame again and rename the columns appropriately.

```
[2]: data = pd.read_csv("https://cf-courses-data.s3.us.cloud-object-storage.

⇔appdomain.cloud/IBM-ML241EN-SkillsNetwork/labs/datasets/Wine_Quality_Data.

⇔csv", sep=',')
```

```
[3]: y = (data['color'] == 'red').astype(int)
fields = list(data.columns[:-1]) # everything except "color"
correlations = data[fields].corrwith(y)
correlations.sort_values(inplace=True)
correlations
```

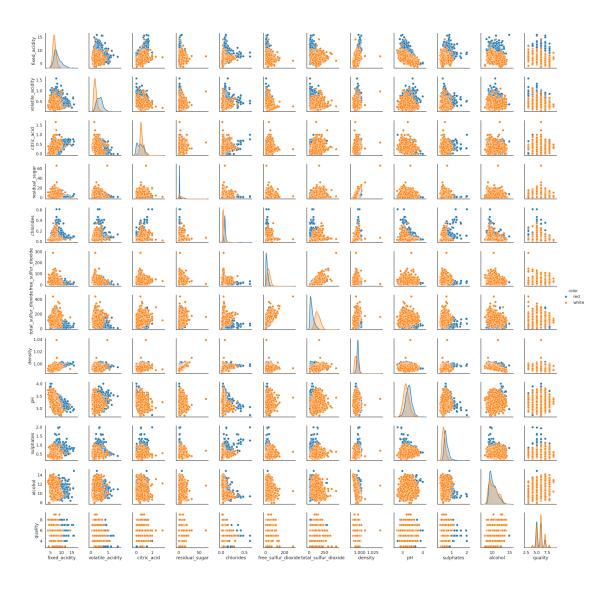
```
[3]: total_sulfur_dioxide -0.700357 free_sulfur_dioxide -0.471644
```

```
residual_sugar
                       -0.348821
citric_acid
                       -0.187397
quality
                       -0.119323
alcohol
                       -0.032970
рΗ
                       0.329129
density
                        0.390645
fixed_acidity
                        0.486740
sulphates
                        0.487218
chlorides
                        0.512678
volatile_acidity
                        0.653036
dtype: float64
```

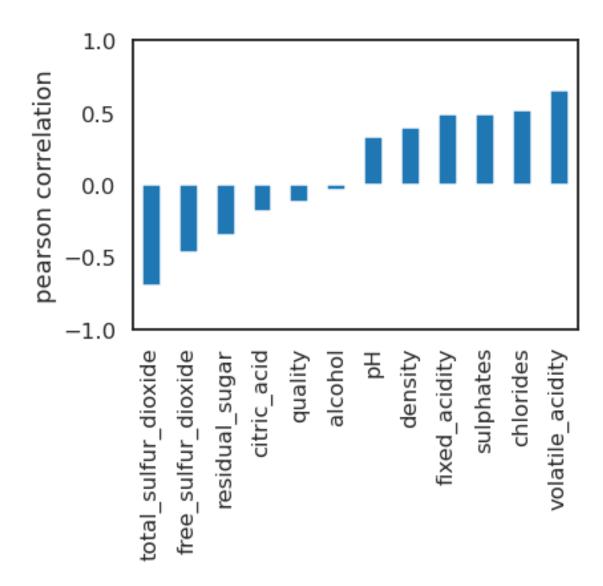
```
[4]: sns.set_context('talk')
#sns.set_palette(palette)
sns.set_style('white')
```

```
[5]: sns.pairplot(data, hue='color')
```

[5]: <seaborn.axisgrid.PairGrid at 0x7f19f12dc2d0>



```
[6]: ax = correlations.plot(kind='bar')
ax.set(ylim=[-1, 1], ylabel='pearson correlation');
```



```
from sklearn.preprocessing import MinMaxScaler

fields = correlations.map(abs).sort_values().iloc[-2:].index
print(fields)
X = data[fields]
scaler = MinMaxScaler()
X = scaler.fit_transform(X)
X = pd.DataFrame(X, columns=['%s_scaled' % fld for fld in fields])
print(X.columns)
```

Index(['volatile_acidity', 'total_sulfur_dioxide'], dtype='object')
Index(['volatile_acidity_scaled', 'total_sulfur_dioxide_scaled'],
dtype='object')

1.4 Part 2: Linear Decision Boundary

Our goal is to look at the decision boundary of a LinearSVC classifier on this dataset. Check out this example in sklearn's documentation.

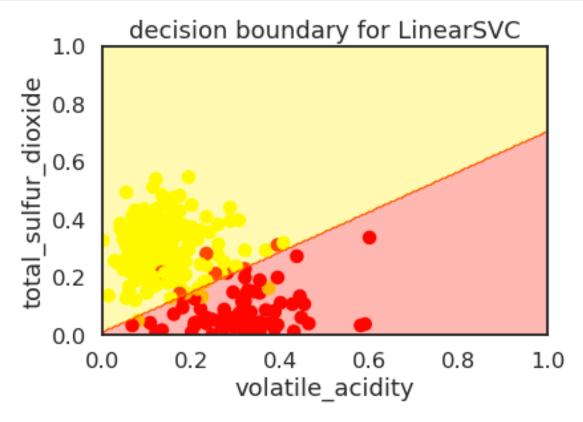
- Fit a Linear Support Vector Machine Classifier to X, y.
- Pick 300 samples from X. Get the corresponding y value. Store them in variables X_color and y_color. This is because original dataset is too large and it produces a crowded plot.
- Modify y_color so that it has the value "red" instead of 1 and 'yellow' instead of 0.
- Scatter plot X_color's columns. Use the keyword argument "color=y_color" to color code samples.
- Use the code snippet below to plot the decision surface in a color coded way.

```
x_axis, y_axis = np.arange(0, 1, .005), np.arange(0, 1, .005)
xx, yy = np.meshgrid(x_axis, y_axis)
xx_ravel = xx.ravel()
yy_ravel = yy.ravel()
X_grid = pd.DataFrame([xx_ravel, yy_ravel]).T
y_grid_predictions = *[YOUR MODEL]*.predict(X_grid)
y_grid_predictions = y_grid_predictions.reshape(xx.shape)
ax.contourf(xx, yy, y_grid_predictions, cmap=colors, alpha=.3)
```

With LinearSVC, it is easy to experiment with different parameter choices and see the decision boundary.

```
[8]: from sklearn.svm import LinearSVC
     LSVC = LinearSVC()
     LSVC.fit(X, y)
     X_color = X.sample(300, random_state=45)
     y_color = y.loc[X_color.index]
     y_color = y_color.map(lambda r: 'red' if r == 1 else 'yellow')
     ax = plt.axes()
     ax.scatter(
         X_color.iloc[:, 0], X_color.iloc[:, 1],
         color=y color, alpha=1)
     x_{axis}, y_{axis} = np.arange(0, 1.005, .005), np.arange(0, 1.005, .005)
     xx, yy = np.meshgrid(x_axis, y_axis)
     xx_ravel = xx.ravel()
     yy_ravel = yy.ravel()
     X_grid = pd.DataFrame([xx_ravel, yy_ravel]).T
     y_grid_predictions = LSVC.predict(X_grid)
     y_grid_predictions = y_grid_predictions.reshape(xx.shape)
     ax.contourf(xx, yy, y_grid_predictions, cmap=plt.cm.autumn_r, alpha=.3)
     # -----
     ax.set(
         xlabel=fields[0],
```

```
ylabel=fields[1],
xlim=[0, 1],
ylim=[0, 1],
title='decision boundary for LinearSVC');
```



1.5 Part 3: Gaussian Kernel

Let's now fit a Gaussian kernel SVC and see how the decision boundary changes.

Consolidate the code snippets in Question 2 into one function which takes in an estimator, X and y, and produces the final plot with decision boundary. The steps are:
 Fit model

Get sample 300 records from X and the corresponding y's

Create grid, predict, plot using ax.contourf

Add on the scatter plot

- After copying and pasting code, the finished function uses the input estimator and not the LinearSVC model.
- For the following values of gamma, create a Gaussian Kernel SVC and plot the decision boundary.

gammas = [.5, 1, 2, 10]

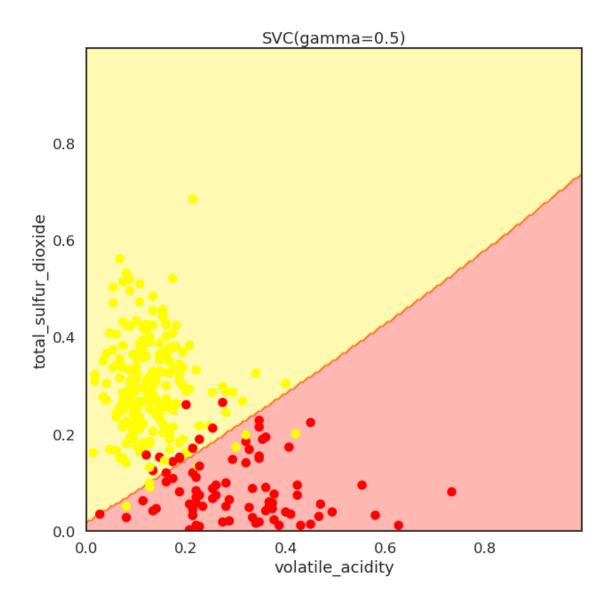
• Holding gamma constant, we plot the decision boundary for various values of C:

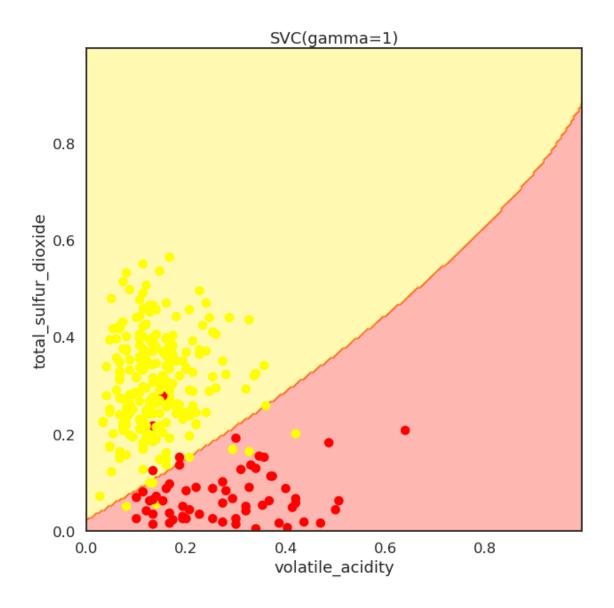
[.1, 1, 10]

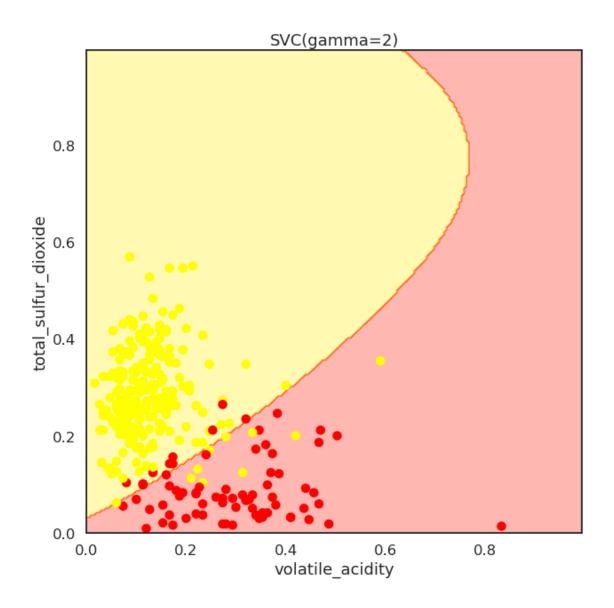
```
[9]: def plot_decision_boundary(estimator, X, y):
         estimator.fit(X, y)
         X_{color} = X.sample(300)
         y_color = y.loc[X_color.index]
         y_color = y_color.map(lambda r: 'red' if r == 1 else 'yellow')
         x_axis, y_axis = np.arange(0, 1, .005), np.arange(0, 1, .005)
         xx, yy = np.meshgrid(x_axis, y_axis)
         xx_ravel = xx.ravel()
         yy_ravel = yy.ravel()
         X_grid = pd.DataFrame([xx_ravel, yy_ravel]).T
         y_grid_predictions = estimator.predict(X_grid)
         y_grid_predictions = y_grid_predictions.reshape(xx.shape)
         fig, ax = plt.subplots(figsize=(10, 10))
         ax.contourf(xx, yy, y_grid_predictions, cmap=plt.cm.autumn_r, alpha=.3)
         ax.scatter(X_color.iloc[:, 0], X_color.iloc[:, 1], color=y_color, alpha=1)
         ax.set(
             xlabel=fields[0],
             ylabel=fields[1],
             title=str(estimator))
```

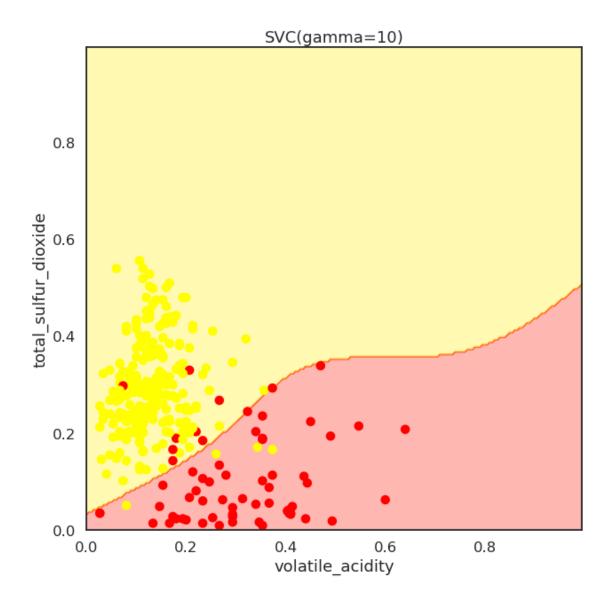
```
[10]: from sklearn.svm import SVC

gammas = [.5, 1, 2, 10]
for gamma in gammas:
    SVC_Gaussian = SVC(kernel='rbf', gamma=gamma)
    plot_decision_boundary(SVC_Gaussian, X, y)
```

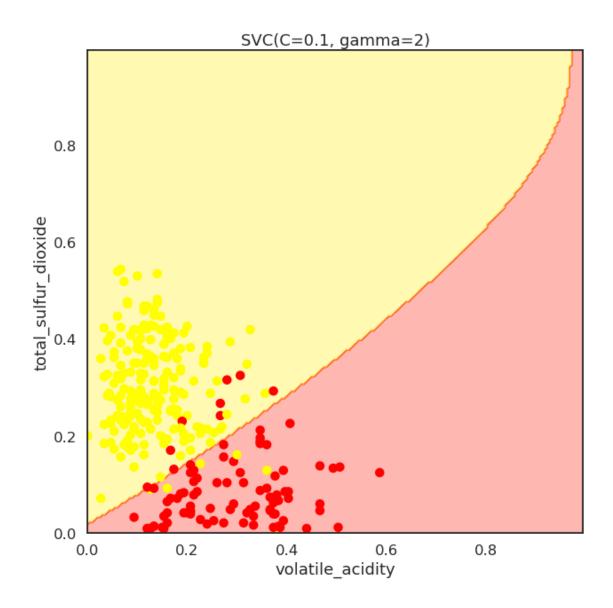


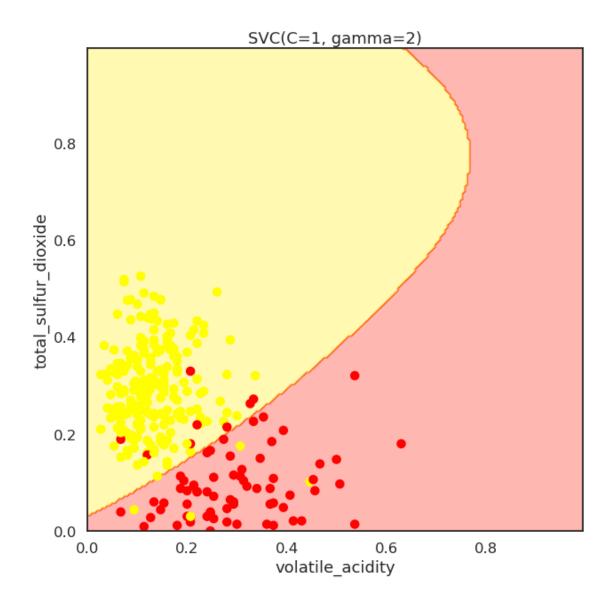


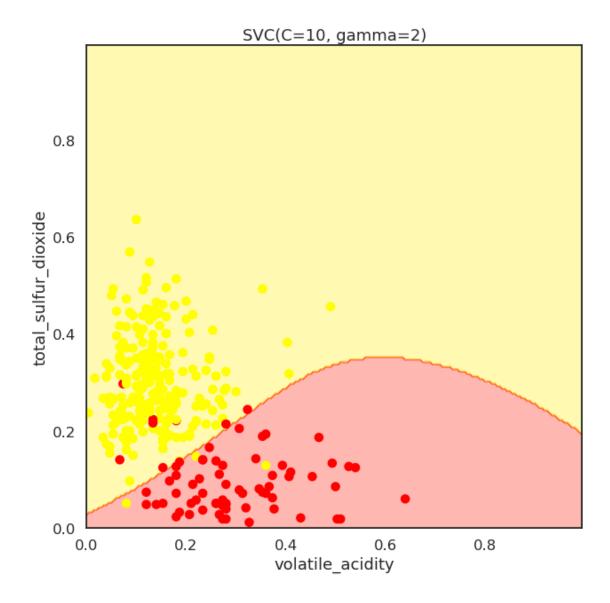




```
[11]: Cs = [.1, 1, 10]
for C in Cs:
    SVC_Gaussian = SVC(kernel='rbf', gamma=2, C=C)
    plot_decision_boundary(SVC_Gaussian, X, y)
```







1.6 Part 4: Comparing Kernel Execution Times

In this exercise, we will compare the fitting times between SVC vs Nystroem with rbf kernel. Jupyter Notebooks provide a useful magic function **%timeit** which executes a line and prints out the time it took to fit. If we type **%%timeit** in the beginning of the cell, it will output the execution time.

We proceed with the following steps:

- Create y from data.color, and X from the rest of the columns.
- Use %%timeit to get the time for fitting an SVC with rbf kernel.
- Use %%timeit to get the time for the following: fit_transform the data with Nystroem and then fit a SGDClassifier.

Nystroem+SGD will take much less to fit. This difference will be more pronounced if the dataset

was bigger.

- Make 5 copies of X and concatenate them
- Make 5 copies of y and concatenate them
- Compare the time it takes to fit the both methods above

```
[12]: from sklearn.kernel_approximation import Nystroem
      from sklearn.svm import SVC
      from sklearn.linear_model import SGDClassifier
      y = data.color == 'red'
      X = data[data.columns[:-1]]
      kwargs = {'kernel': 'rbf'}
      svc = SVC(**kwargs)
      nystroem = Nystroem(**kwargs)
      sgd = SGDClassifier()
[13]: %%timeit
      svc.fit(X, y)
     587 ms \pm 16.7 ms per loop (mean \pm std. dev. of 7 runs, 1 loop each)
[14]: %%timeit
      X_transformed = nystroem.fit_transform(X)
      sgd.fit(X_transformed, y)
     277 ms \pm 44 ms per loop (mean \pm std. dev. of 7 runs, 1 loop each)
[15]: X2 = pd.concat([X]*5)
      y2 = pd.concat([y]*5)
      print(X2.shape)
      print(y2.shape)
     (32485, 12)
     (32485,)
 []: %timeit svc.fit(X2, y2)
 []: | %%timeit
      X2_transformed = nystroem.fit_transform(X2)
      sgd.fit(X2_transformed, y2)
```

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