# Assignment 2: Data Preprocessing, Parameter Tuning, Model Evaluation Using SVM and K-NN

# What you will learn

- · Data preprocessing
- · Parameter Tuning
- Model evaluation
- Employing SVM and K-NN methods on the data

# Setup

- Download Anaconda Python 3.6 for consistent environment.
- · Download Pandas library.
- If you use pip environment then make sure your code is compatible with versions of libraries provided within Anaconda's Python 3.6 distribution.

#### **Submission**

- Do not change any variable/function names.
- · Just add your own code and don't change existing code
- Save this file and rename it to be studentid\_lastname.ipynb (student id (underscore) last name.ipynb) where your student id is all numbers
- Export your .ipynb file to PDF (File > Download as > PDF via Latex). Please don't leave this step for final minutes .
- Submit both the notebook and PDF files (NO ZIP, RAR,..).
- If you happen to use any external library not included in Anaconda (mention in Submission Notes section below)

#### In [14]:

```
from sklearn.svm import SVC
from sklearn.pipeline import make pipeline
from sklearn.preprocessing import StandardScaler
from sklearn.model_selection import GridSearchCV
from sklearn.model selection import train test split
from sklearn.model_selection import StratifiedShuffleSplit
from sklearn.cluster import KMeans
from sklearn.metrics import classification report
import itertools
from sklearn.metrics import confusion matrix
import pandas as pd
import numpy as np
import matplotlib.pyplot as plt
%matplotlib inline
# remove the following statements if you like to see warnings
import warnings
warnings.filterwarnings('ignore')
```

### **Submission Notes**

(Please write any notes here that you think I should know during marking)

# [NO MARKS] Warming Up

Various interesting machine learning datasets can be found at:

https://archive.ics.uci.edu/ml/index.php

For this task, we have chosen the Heart Disease dataset, available at:

• https://archive.ics.uci.edu/ml/datasets/heart+Disease

#### **Data Set Information**

- The dataset contains 303 subjects with 76 attributes.
- All the published experiments refer to using a subset of 14 of attributes.
- The goal field in the dataset refers to the presence of heart disease in the patient.
- It is integer valued from 0 (no presence) to 4 (highest presence).

Note: Since the class number 4 is very sparse (just 13 subjects). We have dropped the subjects belonging to class 4 from the data.

Experiments with the Cleveland dataset have concentrated on---attempting to distinguish **the presence** (values 1, 2, 3, 4) from **the absence** (value 0).

#### **Attribute Information**

14 attributes are been used:

- 1. #3 (age)
- 2. #4 (sex)
- 3. #9 (cp)
- 4. #10 (trestbps)
- 5. #12 (chol)
- 6. #16 (fbs)
- 7. #19 (restecg)
- 8. #32 (thalach)
- 9. #38 (exang)
- 10. #40 (oldpeak)
- 11. #41 (slope)
- 12. #44 (ca)
- 13. #51 (thal)
- 14. #58 (num) (the predicted attribute)

# (no marks) Reading the data

```
In [15]:
```

Data-set shape: (303, 14)

# Out[15]:

	age	sex	ср	trestbps	chol	fbs	restecg	thalach	exang	oldpeak	slope	ca	thal	prediction
293	63.0	1.0	4.0	140.0	187.0	0.0	2.0	144.0	1.0	4.0	1.0	2.0	7.0	2
294	63.0	0.0	4.0	124.0	197.0	0.0	0.0	136.0	1.0	0.0	2.0	0.0	3.0	1
295	41.0	1.0	2.0	120.0	157.0	0.0	0.0	182.0	0.0	0.0	1.0	0.0	3.0	0
296	59.0	1.0	4.0	164.0	176.0	1.0	2.0	90.0	0.0	1.0	2.0	2.0	6.0	3
297	57.0	0.0	4.0	140.0	241.0	0.0	0.0	123.0	1.0	0.2	2.0	0.0	7.0	1
298	45.0	1.0	1.0	110.0	264.0	0.0	0.0	132.0	0.0	1.2	2.0	0.0	7.0	1
299	68.0	1.0	4.0	144.0	193.0	1.0	0.0	141.0	0.0	3.4	2.0	2.0	7.0	2
300	57.0	1.0	4.0	130.0	131.0	0.0	0.0	115.0	1.0	1.2	2.0	1.0	7.0	3
301	57.0	0.0	2.0	130.0	236.0	0.0	2.0	174.0	0.0	0.0	2.0	1.0	3.0	1
302	38.0	1.0	3.0	138.0	175.0	0.0	0.0	173.0	0.0	0.0	1.0	?	3.0	0

### (no marks) Removing the missing data

such data-points.

```
In [16]:
```

```
# Removing all subjects from class 4
df = df.loc[df.prediction != 4]

# Replacing the missing data '?' with NAN values
df.replace('?', np.nan, inplace=True)
df = df.dropna()
df = df.astype(float)
```

### (no marks) Splitting the data

```
In [17]:
```

```
# Separating the data and the labels
X = np.asarray(df[df.columns[:-1]]).astype(np.float32)
y = np.asarray(df.prediction)
# Splitting the data into the train and the test sets
sss = StratifiedShuffleSplit(n splits=1, test size=0.2, random state=0)
sss.get_n_splits(X, y)
train_index, test_index = next(sss.split(X, y))
X_train, X_test = X[train_index], X[test_index]
y train, y test = y[train index], y[test index]
print('Training data: \n', X train)
print('\n')
print('Training labels: \n',y train)
Training data:
 [[45. 0. 4. ... 2. 0. 3.]
 [41. 1. 2. ... 2. 0. 6.]
 [42. 1. 4. ... 1. 0. 3.]
 [67. 1. 4. ... 2. 2. 7.]
[65. 1. 4. ... 2. 1. 7.]
 [62. 0. 4. ... 1. 0. 3.]]
Training labels:
 [0. \ 0. \ 0. \ 1. \ 0. \ 3. \ 0. \ 0. \ 1. \ 1. \ 3. \ 0. \ 1. \ 1. \ 0. \ 2. \ 3. \ 0. \ 1. \ 3. \ 3. \ 0. \ 0.
 0. 2. 0. 0. 3. 0. 0. 0. 3. 1. 2. 0. 0. 3. 0. 0. 1. 1. 1. 1. 1. 1. 0. 0.
 1. 3. 1. 2. 0. 2. 2. 0. 0. 0. 1. 3. 1. 0. 2. 0. 0. 0. 0. 1. 0. 0. 0.
 3. 0. 0. 0. 1. 0. 0. 0. 3. 1. 0. 0. 1. 1. 1. 0. 0. 3. 0. 0. 0. 2. 2. 0.
 0. 0. 0. 0. 0. 3. 0. 2. 2. 2. 0. 1. 0. 0. 1. 1. 0. 0. 0. 0. 0. 0. 0. 2.
 2. 0. 0. 1. 0. 0. 0. 1. 3. 0. 0. 0. 0. 0. 0. 2. 2. 1. 2. 0. 1. 0. 2. 0.
 3. 1. 0. 0. 0. 3. 0. 0. 0. 2. 2. 0. 2. 1. 0. 0. 0. 3. 3. 0. 0. 0. 0. 1.
 0. 1. 3. 0. 0. 1. 0. 1. 0. 0. 0. 0. 3. 0. 1. 0. 0. 3. 0. 2. 3. 0. 0.
 0. 0. 2. 0. 1. 1. 0. 0. 3. 0. 2. 1. 0. 2. 2. 0. 0. 0. 1. 0. 0. 3. 1.
 2. 0. 3. 0. 0. 3. 0. 3. 1. 2. 0.]
```

# [10 marks] Data Exploration

a) (3 marks) Use pandas to find the ratio of the presence of disease versus the absence within the different sex.

Note: 0 is female and 1 is male.

```
In [18]:
```

```
# we have defined a new column which is `true` if there is a presence of disease (i.e., prediction
is [1, 2, 3])
df['has_disease'] = df.apply(lambda x: x.prediction in [1, 2, 3], axis=1)

# use groupby and aggregation
aggr = {'ratio': lambda x: sum(x) / (len(x) - sum(x))}
display(df.groupby('sex')['has_disease'].agg(aggr))
```

```
ratio
sex
0.0 0.323944
1.0 1.134831
```

b) (7 marks) Do the same thing for age. Split the age groups as follows (left included, right isn't):

- 1. [29, 49)
- 2. [49, 69)
- 3. [69, inf)

And then find the average ratio of prevalence of the heart disease within the each group.

#### In [19]:

```
# write your code here
ranges = [28,48,68,np.inf]
display(df.groupby(pd.cut(df.age, ranges))['has disease'].agg(aggr))
```

#### ratio

```
age
   (28.0,
           0.385965
   (48.0,
           1.042105
    68.01
(68.0, inf] 0.375000
```

# (no marks) Utility function

# In [20]:

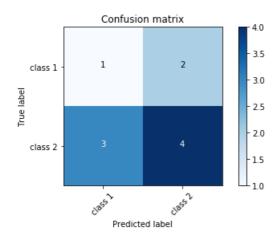
```
# Do not change the function
# This function is adapted from the sklearn website
# This function let you draw a confusion matrix for your problem
def plot_confusion_matrix(cm, classes,
                          normalize=False,
                          title='Confusion matrix',
                          cmap=plt.cm.Blues):
    This function prints and plots the confusion matrix.
    Normalization can be applied by setting `normalize=True`.
    if normalize:
       cm = cm.astype('float') / cm.sum(axis=1)[:, np.newaxis]
       print("Normalized confusion matrix")
    else:
       print('Confusion matrix, without normalization')
    print (cm)
    plt.imshow(cm, interpolation='nearest', cmap=cmap)
    plt.title(title)
    plt.colorbar()
    tick_marks = np.arange(len(classes))
    plt.xticks(tick_marks, classes, rotation=45)
    plt.yticks(tick marks, classes)
    fmt = '.2f' if normalize else 'd'
    thresh = cm.max() / 2.
    for i, j in itertools.product(range(cm.shape[0]), range(cm.shape[1])):
       plt.text(j, i, format(cm[i, j], fmt),
                 horizontalalignment="center",
                 color="white" if cm[i, j] > thresh else "black")
    plt.ylabel('True label')
```

```
plt.xlabel('Predicted label')
plt.tight layout()
```

#### In [21]:

```
# usage
plot confusion matrix(np.array([[1, 2], [3, 4]]), ['class 1', 'class 2'])
```

```
Confusion matrix, without normalization
[[1 2]
 [3 4]]
```



# Task 1 [20 marks]

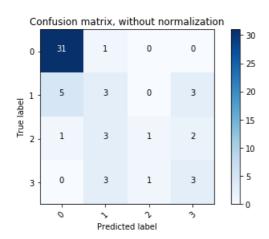
# a) [10 marks] Applying KNN to the data

#### In [22]:

```
from sklearn.neighbors import KNeighborsClassifier
from sklearn.preprocessing import MinMaxScaler
# Task 2
# Add your code in the following part:
# We use min max scaler to normalize the features between [0, 1]
scaler = MinMaxScaler()
# Add your code here instead of ...
scaler.fit(X train)
# Create a knn classifier instance here (If you don't add anything here, your code won't execute!)
knn clf = KNeighborsClassifier()
# Fit the classifier using the train data (If you don't add anything here, your code won't execute
knn_clf.fit(scaler.transform(X_train), y_train)
# Predict the test class labels using the trained classifier (If you don't add anything here, your
code won't execute!)
y pred = knn clf.predict(scaler.transform(X test))
# (If you don't add anything here, your code won't execute!)
print(classification report(y test, y pred))
cnf matrix = confusion_matrix(y_test, y_pred)
plot_confusion_matrix(cnf_matrix, classes=[0,1,2,3],
                      title='Confusion matrix, without normalization')
```

	precision	recall	f1-score	support
0.0	0.84	0.97	0.90	32
1.0	0.30	0.27	0.29	11
2.0	0.50	0.14	0.22	7
3.0	0.38	0.43	0.40	7

```
0.67
                                   0.67
  micro avg
                          0.67
                 0.50
                          0.45
                                  0.45
                                              57
  macro avq
weighted avg
                 0.64
                         0.67
                                  0.64
                                              57
Confusion matrix, without normalization
[[31 1 0 0]
[5 3 0 3]
[1 3 1 2]
 [ 0 3 1 3]]
```



# b) [5 marks] Between \$K=3\$ and \$k=5\$ which one gives more accuracy?

In [23]:

```
from sklearn.metrics import accuracy_score

# write your code here and populate `y_pred_k5` and `y_pred_k3`
# knn_clf defaults to 5 neighbors
knn_clf.n_neighbors = 5
y_pred_k5 = knn_clf.predict(scaler.transform(X_test))

knn_clf.n_neighbors = 3
y_pred_k3 = knn_clf.predict(scaler.transform(X_test))

print(accuracy_score(y_test, y_pred_k5), accuracy_score(y_test, y_pred_k3))
```

# c) [5 marks] Between \$\ell\_1\$, \$\ell\_2\$, and \$cosine\$ similarity which one is better in term of accuracy?

In [24]:

```
# write your code here to experiment with different distance metrics
# use argument `metric` to change to a different distance by default it is euclidean distance

metrics = {
    'll': 'manhattan',
    'l2': 'minkowski',
    'cos': 'cosine'
}

for k, v in metrics.items():
    clf = KNeighborsClassifier(metric=v)
    clf.fit(scaler.transform(X_train), y_train)
    y_pred = clf.predict(scaler.transform(X_test))

    print(k, ': ', accuracy_score(y_test, y_pred))
```

# Task 2 [10 marks]

# Understanding the pipelining architecture of Sklearn

In the code above, you had to call scaler for every prediction by a model. This can be avoided by using a pipeline mechanism within sklearn. Look at the code below:

- 1. We create a data scaler (can be any sclaer with fit and transform functions).
- 2. We create SVC object (again with fit anf transform functions).
- 3. Then we create a pipeline: data  $\operatorname{--}>$  scaler  $\operatorname{--}>$  svc  $\operatorname{--}>$  fit.
- 4. The same transformation is also applied during the prediction phase.

We further use a <code>GridSearchCV</code> for the SVC's parameters tuning.

```
In [25]:
```

```
# Task 2
# Creating a SVM classifier instance
svc = SVC()
# Add a scaler here (If you don't add anything here, your code won't execute!)
data scaler = StandardScaler()
# Update the pipeline by adding the scaler from the previous line
model = make_pipeline(data_scaler, svc)
grid = GridSearchCV(model, param grid)
%time grid.fit(X train, y train)
print( grid.best_estimator_)
Wall time: 189 ms
Pipeline (memory=None,
   steps=[('standardscaler', StandardScaler(copy=True, with mean=True, with std=True)), ('svc',
SVC(C=50, cache size=200, class weight=None, coef0=0.0,
 decision_function_shape='ovr', degree=3, gamma=0.001, kernel='rbf',
 max iter=-1, probability=False, random state=None, shrinking=True,
 tol=0.001, verbose=False))])
In [26]:
# Selecting the best estimator after the parameter search
model = grid.best_estimator_
In [27]:
```

# In [28]:

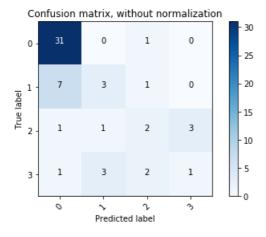
# Predicting the test labels
y pred = model.predict(X test)

```
# Printing the classification report
print(classification_report(y_pred=y_pred,y_true=y_test))
```

		precision	recall	f1-score	support
	0.0	0.78 0.43	0.97 0.27	0.86 0.33	32 11
	2.0	0.33 0.25	0.29 0.14	0.31 0.18	7 7
micro macro weighted	avg	0.65 0.45 0.59	0.65 0.42 0.65	0.65 0.42 0.61	57 57 57

```
In [29]:
```

```
Confusion matrix, without normalization
[[31 0 1 0]
[ 7 3 1 0]
[ 1 1 2 3]
[ 1 3 2 1]]
```



# Task 3 [40 marks]

### How to handle the missing data

More information can be found here: (https://pandas.pydata.org/pandas-docs/stable/user\_guide/missing\_data.html)

a) [6 scores] Name two numeric methods for dealing with the missing data (except dropping):

```
Write the answer here:

1- Interpolation

2- Filling missing values with mean: fillna
```

b) [12 scores] Apply the methods that you mentioned in part (a) to the df with missing data dataframe:

#### In [30]:

c) [22 scores] Apply the steps described in  $Task\ 2$  on df\_1 and df\_2 and show the results using classification\_report and plot confusion matrix.

```
In [31]:
# Task 3 part (c)
# Add your code here
def init grid():
    # Creating a SVM classifier instance
    svc = SVC()
    # Add a scaler here (If you don't add anything here, your code won't execute!)
    data scaler = StandardScaler()
    # Update the pipeline by adding the scaler from the previous line
    model = make pipeline(data scaler, svc)
    param grid = {'svc C': [1, 5, 10, 50],
                  'svc_gamma': [0.0001, 0.0005, 0.001, 0.005]}
    return GridSearchCV(model, param grid)
def split data(df):
    # Separating the data and the labels
    X = np.asarray(df[df.columns[:-1]]).astype(np.float64)
    y = np.asarray(df.prediction)
    # Splitting the data into the train and the test sets
    sss = StratifiedShuffleSplit(n_splits=1, test_size=0.2, random_state=0)
    sss.get_n_splits(X, y)
    train_index, test_index = next(sss.split(X, y))
    X train, X test = X[train index], X[test index]
    y train, y test = y[train_index], y[test_index]
    return X train, X test, y train, y test
def get prediction(data):
    X_train, X_test, y_train, y_test = data
   %time grid.fit(X train, y train)
   grid.best estimator
   model = grid.best_estimator_
    y pred = model.predict(X test)
    return y_pred, y_test
grid = init grid()
method = ('Interpolation', 'Mean')
# Process the 2 data set and display results as instructed
for i, df in enumerate((df 1, df 2)):
    print('\n')
   print('Classification Report For Data Filled By ' + method[i])
   y_p, y_t = get_prediction(split_data(df))
   print(classification_report(y_pred=y_p,y_true=y_t))
    # Computing the confusion matrix for the test data
    cnf_matrix = confusion_matrix(y_t, y_p)
    # Plotting the confusion matrix using the previous function
    plot confusion matrix(cnf matrix, classes=[0,1,2,3],
                          title= method[i] + ' Filled Data Confusion matrix, without normalization'
    # Ensures the plotted image displays during runtime
    plt.show()
Classification Report For Data Filled By Interpolation
Wall time: 166 ms
             precision recall f1-score support
```

0.0

1.0

2.0

3.0

0.82

0.17

0.25

0.25

0 00

0.94

0.18

0.14

0.14

0 00

0.87

0.17

0.18

0.18

0 00

33

11

7

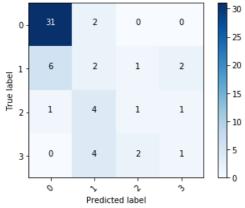
7

	micro	avg	U.6U	U.6U	U.6U	つと
	macro	avg	0.37	0.35	0.35	58
W	eighted	avq	0.56	0.60	0.57	58

Confusion matrix, without normalization

[[31 2 0 0] [6 2 1 2] [1 4 1 1] [0 4 2 1]]

Interpolation Filled Data Confusion matrix, without normalization



Classification Report For Data Filled By Mean

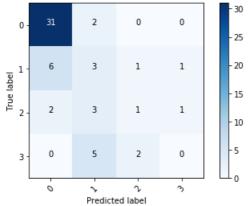
Wall time: 160 ms

		precision	recall	f1-score	support
	0.0	0.79	0.94	0.86	33
	1.0	0.23	0.27	0.25	11
	2.0	0.25	0.14	0.18	7
	3.0	0.00	0.00	0.00	7
micro	avg	0.60	0.60	0.60	58
macro	avg	0.32	0.34	0.32	58
weighted	avg	0.53	0.60	0.56	58

Confusion matrix, without normalization

[[31 2 0 0] [6 3 1 1] [2 3 1 1] [0 5 2 0]]

Mean Filled Data Confusion matrix, without normalization



# Task 4 [20 marks]

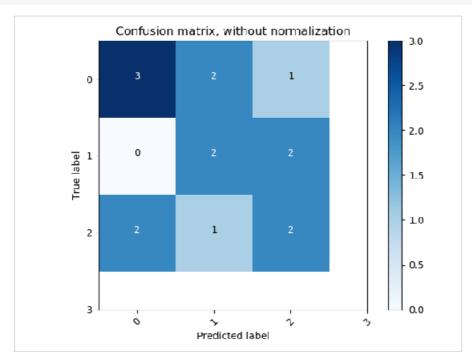
# **Model Evaluation**

For the given confusion matrix, answer the following questions.

In [32]:

```
I = plt.imread('foo.png')
```

```
fig = plt.figure(figsize= (10,10))
plt.imshow(I)
plt.axis('off')
plt.show()
```



#### Calculate the following parameters (Use macro-average definition)

You can find all these definitions on <a href="https://scikit-learn.org/stable/auto\_examples/model\_selection/plot\_precision\_recall.html">https://scikit-learn.org/stable/auto\_examples/model\_selection/plot\_precision\_recall.html</a> and <a href="https://en.wikipedia.org/wiki/Confusion\_matrix">https://en.wikipedia.org/wiki/Confusion\_matrix</a>

Write your answer in front of each parameter:

- 1- [2 scores] total number of instances = 3+2+1+0+2+2+2+1+2 = 15
- 2- [2 scores] number of classes = 3
- 3- [2 scores] True positive (TP) =  $(3+2+2)/3=\sim 2.33$
- 4- [2 scores] True negative (TN) =  $(7+8+7)/3 = \sim 7.33$
- 5- [2 scores] False positive (FP) = (2+3+3)/3=~ 2.67
- 6- [2 scores] False negative (FN) = (3+2+3)/3=~ 2.67
- 7- [2 scores] Sensitivity, recall, hit rate, or true positive rate (TPR) =  $(0.5+0.5+0.4)/3=\sim 0.47$
- 8- [2 scores] Specificity, selectivity or true negative rate (TNR) = (0.78+0.73+0.70)/3=~ 0.74
- 9- [4 scores] F1-Score = (0.55+0.44+0.40)/3=~ 0.46

### In [33]:

# Assignment end