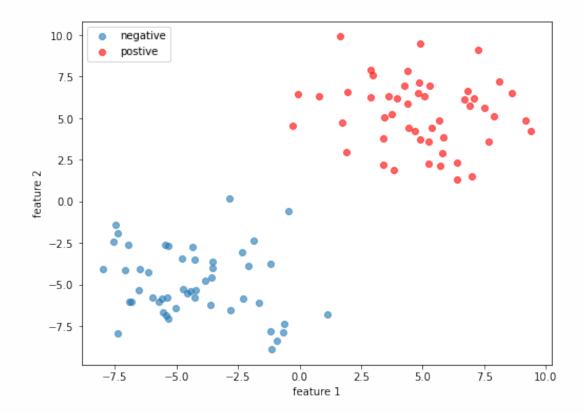
HM 2 CODING

4 Coding for SVM

```
import pandas as pd
import numpy as np
import matplotlib.pyplot as plt
import warnings
warnings.filterwarnings('ignore')
```

4.1 Data generating and plotting

```
mu_neg = np_array([-5, -5])
sigma_neg = 5 * np.identity(2)
mu_pos = np_array([5, 5])
sigma_pos = 5 * np.identity(2)
# neg_1, neg_2 = np.random.multivariate_normal(mu_neg, sigma_neg, 50).T
# pos_1, pos_2 = np.random.multivariate_normal(mu_neg, sigma_pos, 50).T
neg = np.random.multivariate_normal(mu_neg, sigma_neg, 50).T
pos = np.random.multivariate_normal(mu_pos, sigma_pos, 50).T
neg_label = -1 * np.ones(50)
pos_label = np.ones(50)
# scatter plot
fig, ax = plt.subplots(figsize=(8,6))
# ax.subplots(figsize=(8,6))
ax.scatter(neg[0], neg[1], alpha=0.6)
ax.scatter(pos[0], pos[1], alpha=0.6, c='r')
ax.set_xlabel('feature 1')
ax.set_ylabel('feature 2')
ax.legend(['negative', 'postive']);
```



4.2 Draw decision boudary

from sklearn.svm import SVC

```
# concate data
feature1 = np.concatenate((neg[0], pos[0]))
feature2 = np.concatenate((neg[1], pos[1]))
label = np.concatenate((neg_label, pos_label))

X = {
    'feature1': feature1,
    'feature2': feature2
}

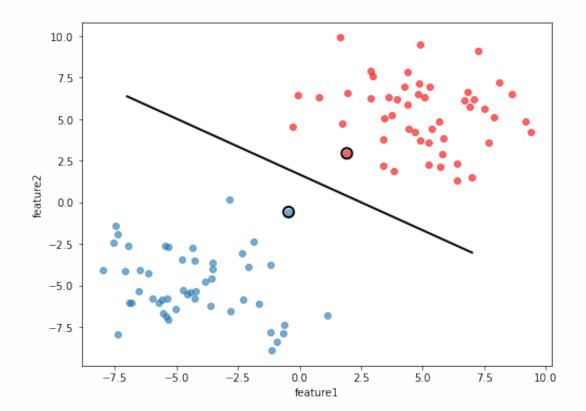
# X_train and y_train
df_X = pd.DataFrame(X)
df_y = pd.DataFrame(label)

df_all = pd.merge(left=df_X, right=df_y, left_index=True, right_index=True)
df_all.columns=['feature1','feature2','label']
```

```
# fit the svc model
svc_clf = SVC(kernel='linear', C=1)
svc_clf.fit(df_X, df_y)
# svc_clf.support_vectors_
```

```
SVC(C=1, kernel='linear')
```

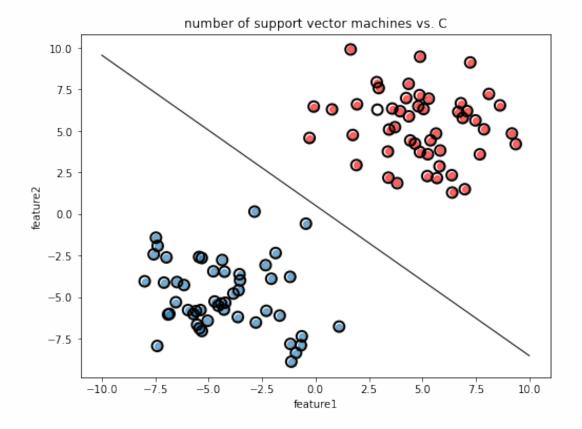
```
# plot the decision function
# get the coefficients and intercept of decision boundary
w = svc_clf.coef_[0]
b = svc_clf.intercept_[0]
x0 = np.linspace(-7, 7) # x
decision_boundary = (-w[0] * x0 - b) / w[1] # line
plt.figure(figsize=(8,6))
plt.plot(x0, decision_boundary, "k-", linewidth=2)
plt.scatter(neg[0], neg[1], alpha=0.6)
plt.scatter(pos[0], pos[1], alpha=0.6, c='r')
# plot support vectors
svs = svc_clf.support_vectors_
plt.scatter(svs[:,0],svs[:,1],s=100,linewidth=2,
                        facecolors='none',edgecolors='k');
plt.xlabel('feature1')
plt.ylabel('feature2');
# plt.grid(True)
```

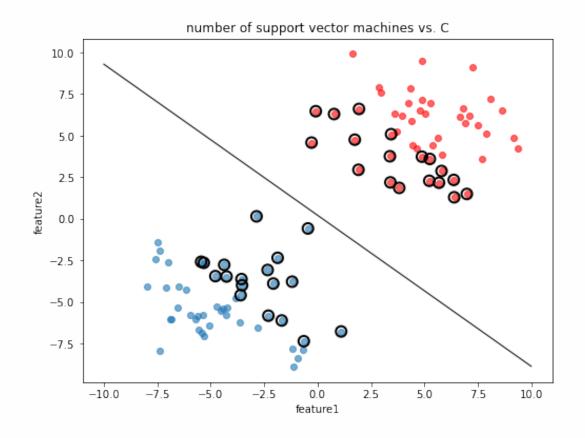


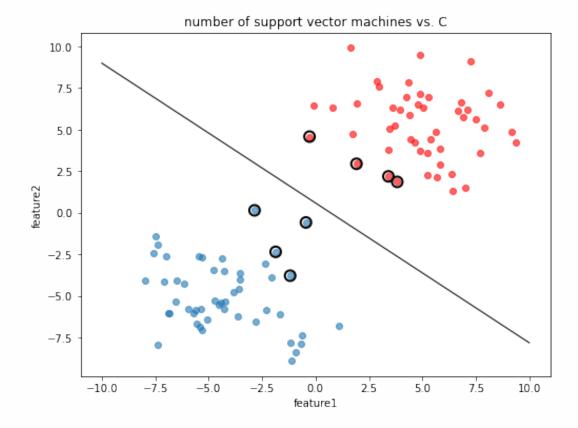
4.3

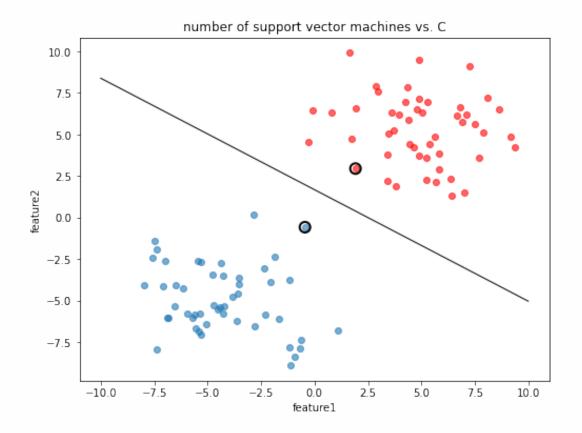
```
def TrainPlotDecisionBoundary(Cs, x0, X_train, y_train,
                             draw_support_vectors, draw_C_svs):
    .....
    A function to plot decision boudaries of SVM.
    ws = [] # store coefs
    bs = [] #store intercepts
    svs_list = []
    n_of_svs = []
    # for each C, train model
    for c in Cs:
        svc_clf = SVC(kernel='linear', C=c)
       svc_clf.fit(X_train, y_train)
       # get the coefficients and intercept of decision boundary
       w = svc_clf.coef_[0]
#
         ws.append(w)
       b = svc_clf.intercept_[0]
         bs.append(b)
#
```

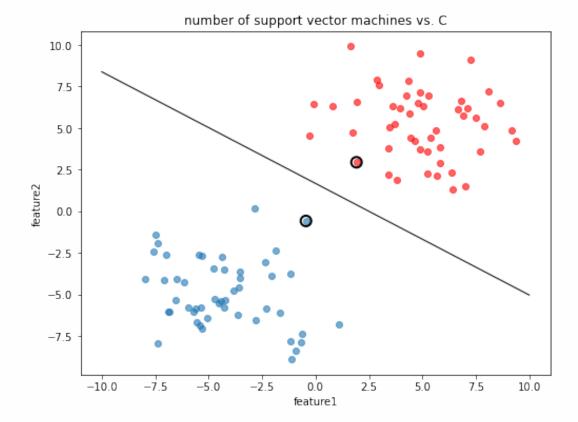
```
svs = svc clf.support vectors
#
          svs_list.append(svs)
        n_svs = len(svs)
        n_of_svs.append(n_svs)
        fig1, ax1 = plt.subplots(figsize=(8,6))
        decision_boundary = (-w[0] * x0 - b) / w[1] # line
        ax1.plot(x0, decision_boundary, "k-", linewidth=1)
        plt.scatter(X_train.iloc[:50,0], X_train.iloc[:50,1], alpha=0.6)
        plt.scatter(X_train.iloc[51:,0], X_train.iloc[51:,1], alpha=0.6, c='r')
        if draw_support_vectors == True:
      plt.figure(figsize=(8,6))
#
          for w,b in zip(ws, bs):
#
            # draw vector machines
            plt.scatter(svs[:,0],svs[:,1],s=100,linewidth=2,
                        facecolors='none',edgecolors='k');
            plt.xlabel('feature1')
            plt.ylabel('feature2');
            plt.title('number of support vector machines vs. C')
    if draw_C_svs == True:
        fig2, ax2 = plt.subplots(figsize=(8,6))
        ax2.plot(np.log(Cs), n_of_svs)
        plt.xlabel('C')
        plt.ylabel('number of support vectors')
```

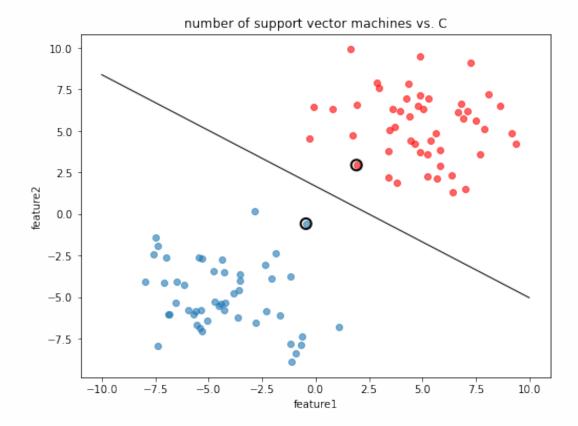


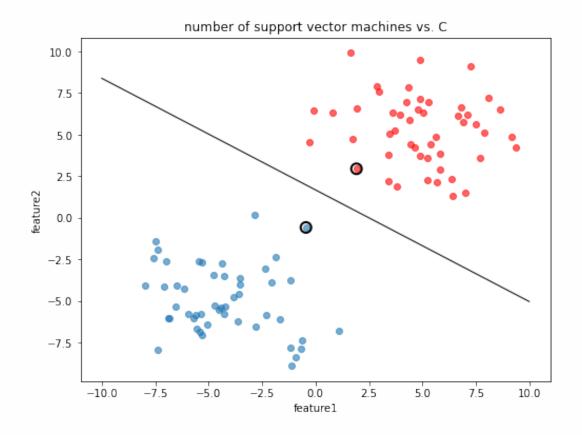


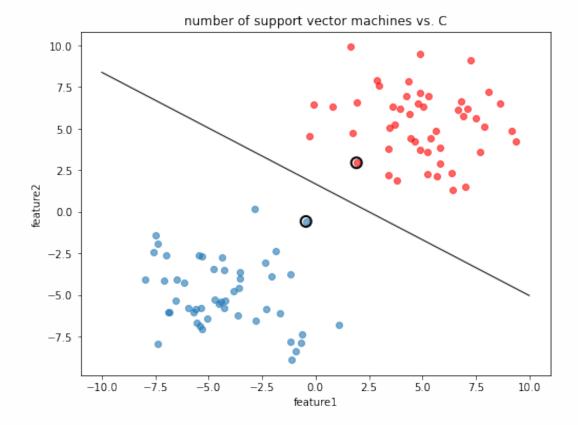


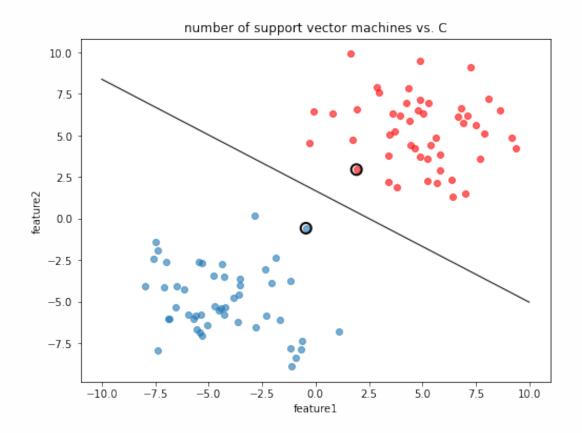


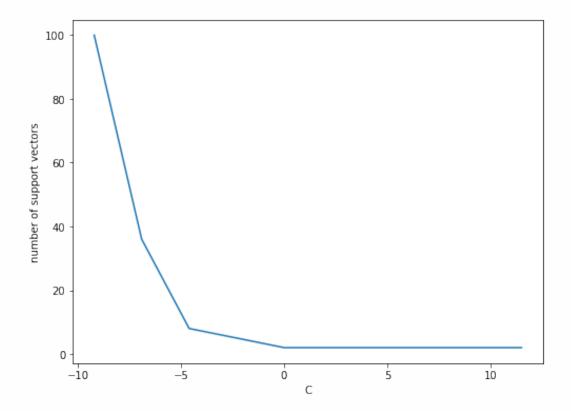












The number of support vectors decreases as C increases.

Because C identifies the penalty of misclassification. The higher C is, the larger the penalty will be. Thus, when C is relatively higher, the model will reduce the number of the violations so the margin is narrower, so there are fewer support vectors. When C is smaller, the penalty reduces, and margin is wider which allows more support vectors.

4.4 Rescale data

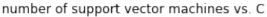
```
# from sklearn.preprocessing import MinMaxScaler

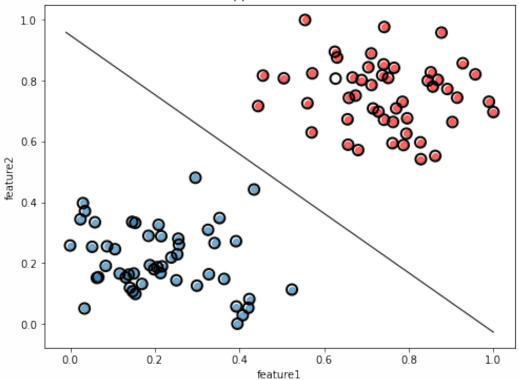
X_train = df_X
y_train = df_y
f1 = X_train.iloc[:, 0]
f2 = X_train.iloc[:, 1]
# min_max_scaler = MinMaxScaler()
min_f1 = min(f1)
max_f1 = max(f1)
min_f2 = min(f2)
max_f2 = max(f2)

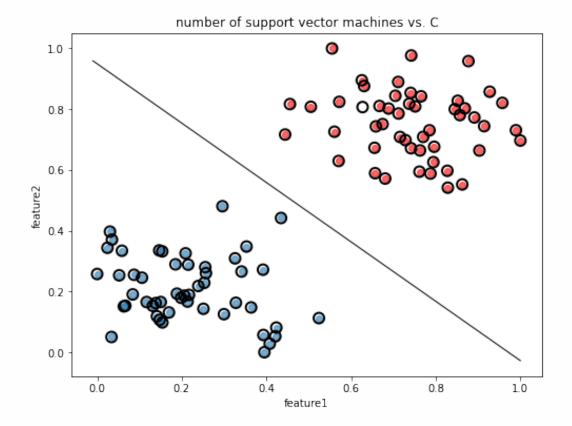
f1_scaled = (f1 - min_f1) / (max_f1 - min_f1)
f2_scaled = (f2 - min_f2) / (max_f2 - min_f2)
```

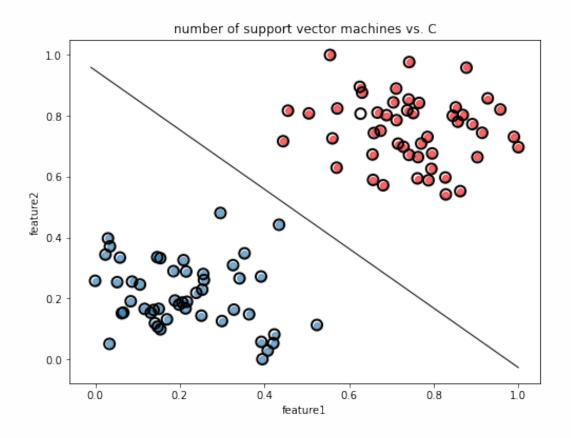
```
scaled = {}
scaled['f1_scaled'] = f1_scaled
scaled['f2_scaled'] = f2_scaled
X_train_scaled = pd.DataFrame(scaled)

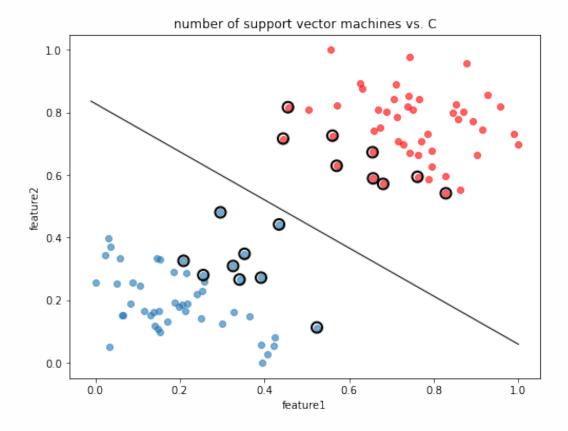
# X_train_scaled
```

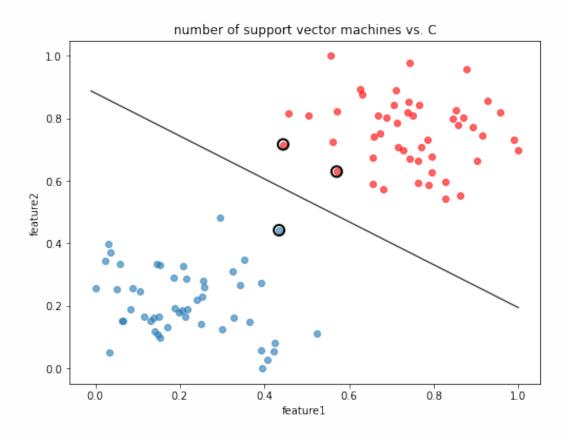


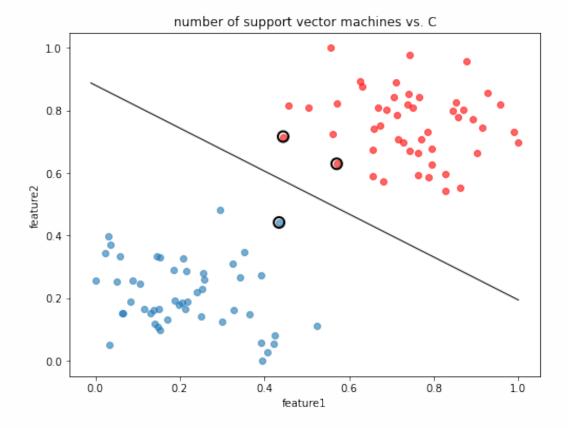


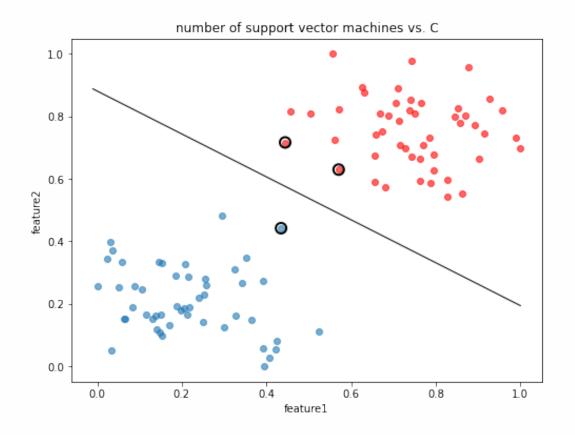


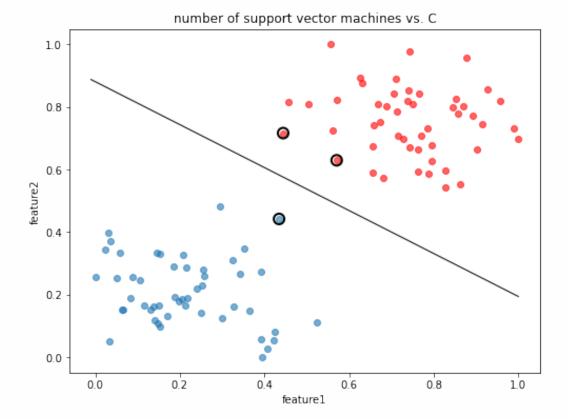


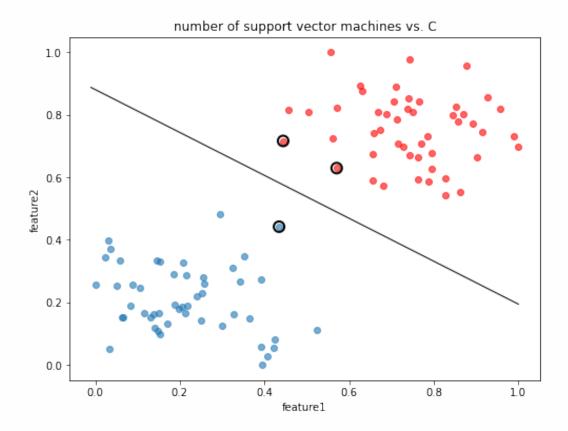


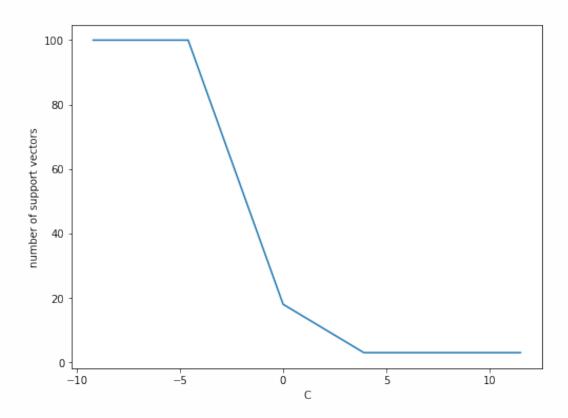












Yes, they are different from the previous question.

The geometric margin decreases while the effect of each feature on the predictions increases.

4.5 boosted decsion trees

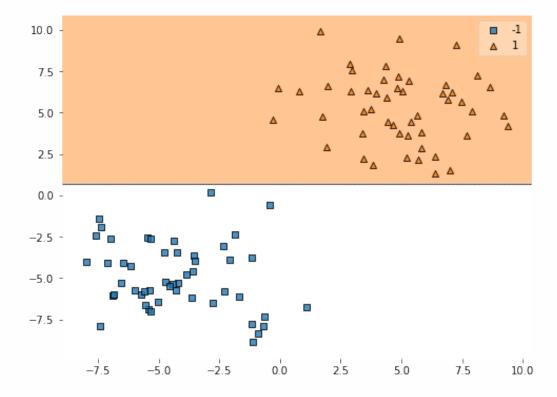
```
from sklearn.ensemble import GradientBoostingClassifier
# from sklearn.inspection import DecisionBoundaryDisplay

gb_clf = GradientBoostingClassifier().fit(X_train, y_train)
gb_clf_rescaled = GradientBoostingClassifier().fit(X_train_scaled, y_train)
# disp = DecisionBoundaryDisplay.from_estimator(
# gb_clf, X_train, response_method="predict",
# xlabel=X_train.columns[:-1], ylabel=y_train.columns[-1],
# alpha=0.5)
```

```
import matplotlib.gridspec as gridspec
from mlxtend.plotting import plot_decision_regions

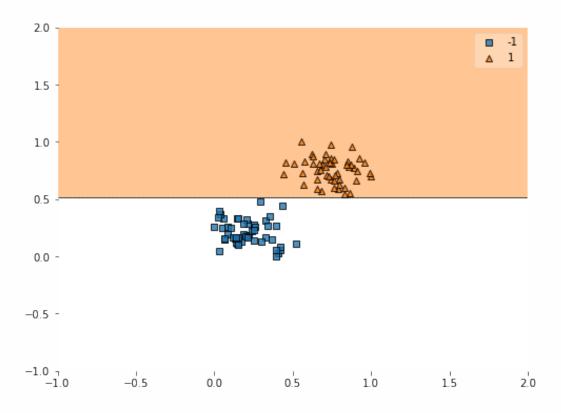
fig = plt.subplots(figsize=(8,6))
plot_decision_regions(np.array(X_train), label.astype(int), gb_clf)
# plot_decision_regions(np.array(X_train_scaled), label.astype(int), gb_clf_rescaled)
```

<AxesSubplot:>



```
fig = plt.subplots(figsize=(8,6))
plot_decision_regions(np.array(X_train_scaled), label.astype(int), gb_clf_rescaled)
```

<AxesSubplot:>



The plots and decision boundaries do not change after rescaling the data when using tree model. That is consistent with my expectation.

5 Coding for Logistic Regression

```
import csv
import math
from sklearn.model_selection import train_test_split
```

5.2 Preprocessing

```
# word dictionary
word_dict = {}
with open('dict.txt') as file:

for line in file:
    (word, index) = line.strip().split(' ')
    index_int = int(index)
    word_dict[index_int] = str(word)

n_words = len(word_dict)
```

```
def GetKey(dictionary, target):
    get the key of certain value query
    for key, value in dictionary.items():
        if value == target:
            return key
# review_dict = {}
label dict = {} # store labels of the reviews
review dict = {} # store the reviews like
                   # { 0: {23:1, 24:1, 25:1, ...}
                    # 1: {13:1, 14:1, 15:1, ...}
                    # n: {34:1, 36:1, ....} }
                # these inner dicts might have different lengths
                # because each review has different number of words
# open .tsv file
with open('moviereview.tsv') as file:
    tsv_file = csv.reader(file, delimiter="\t")
    file list = list(tsv file)
    for i in range(len(file_list)): # get the index of each line
        label = file list[i][0]
        review = file_list[i][1]
        label_int = int(label) # convert string to int
        label_dict[i] = label_int # store labels
        sentence dict = {} # for each line of review,
                            # create a dict to store the Booleans
                            # if the word appears in dictionary:
                                # set to 1
                            # else: ignore
        for word in review.split(' '):
```

```
if word in word_dict.values():
          word_idx = GetKey(word_dict, word)
          sentence_dict[word_idx] = 1

review_dict[i] = sentence_dict

len(review_dict[5])
len(review_dict[10])
```

493

```
# split data
# X_train5, y_train5, X_test5, y_test5 = train_test_split(review_dict, label_dict,
                                                           test_size=0.20,
random_state=10)
def dict_slice(dict_, start, end):
    a function to divide the dictionary
    into expected slices
    .......
    keys = dict_.keys()
    dict_slice = {}
    for k in list(keys)[start:end]:
        dict_slice[k] = dict_[k]
    return dict_slice
# split train set and test set
X_train5 = dict_slice(review_dict, 0, round(0.8*len(review_dict)))
X_test5 = dict_slice(review_dict, round(0.8*len(review_dict)), len(review_dict))
y train5 = dict slice(label dict, 0, round(0.8*len(label dict)))
y_test5 = dict_slice(label_dict, round(0.8*len(label_dict)), len(label_dict))
```

```
# initialization
T = 30
eta = 0.1
theta = [0] * n_words
bias = 0
n_lines = len(X_train5) # total number of reviews in the training set
n_words = len(word_dict) # toal number of words in the dictionary
# 30 epochs
for t in range(T):
```

```
grad vector = [0] * n lines #initialize gradient vector
   for i in range(n_lines): # interate each row of reviews
        sum1 = bias
        for index in X_train5[i]: # for each line, iterate each word index
            sum1 += theta[index] # sum out theta with corresponding index
                                   # this is a faster way to calculate dot product
       y_label = y_train5[i] # get the label for each review
       grad_vector[i] = y_label - 1 / (1 + math.exp(-sum1)) # calculate loss
# print(grad_vector)
   theta_sum = [0] * n_words
   bias sum = 0
   for i in range(n_lines):
       for index in X_train5[i]:
            theta sum[index] += grad vector[i]
        bias_sum += grad_vector[i]
   for index in range(n_words):
        theta[index] += eta * theta_sum[index] / n_lines
   bias += eta * bias_sum / n_lines
```

```
# run test set

best_theta = theta
sup_script = 0
pred_labels = []
for key, reviews in X_test5.items():
    for idx, values in reviews.items():
        sup_script += best_theta[idx]

P_pos = 1 / (1 + math.exp(sup_script))
    if P_pos > 0.5:
        pred_labels.append(1)

else:
    pred_labels.append(0)

def Accuracy(true_label, predict_label):
    """"
    a function to calculate prediction accuracy
    """"
    correct = 0
```

```
incorrect = 0
for i,j in zip(true_label, predict_label):
    if i == j:
        correct += 1

    elif i != j:
        incorrect += 1

accuracy_rate = correct / len(predict_label)

return accuracy_rate

Accuracy(y_test5.values(), pred_labels)
```

0.9208333333333333

6 Boosted Decision Trees and Random Forest

6.1 Preprocessing

```
from sklearn.ensemble import RandomForestClassifier
from sklearn.ensemble import AdaBoostClassifier
from sklearn import metrics
from sklearn.metrics import plot_roc_curve, roc_auc_score
from sklearn.preprocessing import OneHotEncoder
from sklearn.model_selection import KFold
from sklearn.model_selection import train_test_split, GridSearchCV
```

```
titanic = pd.read_csv('Titanic.csv')
titanic.head()
```

```
.dataframe tbody tr th {
    vertical-align: top;
}
.dataframe thead th {
    text-align: right;
}
```

	Survived	Pclass	Sex	Age	SibSp	Parch	Fare	Cabin	Embarked
0	0	3	male	22.0	1	0	7.2500	NaN	S
1	1	1	female	38.0	1	0	71.2833	C85	С
2	1	3	female	26.0	0	0	7.9250	NaN	S
3	1	1	female	35.0	1	0	53.1000	C123	S
4	0	3	male	35.0	0	0	8.0500	NaN	S

```
# drop NaN columns
df_ti = titanic.dropna(axis='columns')
df_ti.shape
df_ti1 = df_ti.copy()
```

```
(891, 6)
```

```
# encoding sex

def SexEncoder(df):

    if df['Sex'] == 'male':
        return 0
    elif df['Sex'] == 'female':
        return 1

df_encoded = pd.DataFrame(df_ti.apply(SexEncoder, axis=1), columns=['Sex_encoding'])
    df_ti_encoded = pd.merge(left=df_ti, right=df_encoded, left_index=True, right_index=True)
    df_ti_encoded1 = df_ti_encoded.drop(columns='Sex')
```

```
# split training set and test set

# split X and y
X = df_ti_encoded1.drop(columns='Survived')
y = df_ti_encoded1['Survived']

X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.20, random_state=10)
```

6.2 Fit models and estimate training time

```
RF_clf = RandomForestClassifier()
RF_clf.fit(X_train, y_train)
%time
```

```
CPU times: user 3 μs, sys: 1 μs, total: 4 μs
Wall time: 5.96 μs
```

```
Ada_clf = AdaBoostClassifier()
Ada_clf.fit(X_train, y_train)
%time
```

```
CPU times: user 2 \mus, sys: 1 \mus, total: 3 \mus Wall time: 5.01 \mus
```

Adaboost required less time to train.

Because Adaboost and boosted tree models constraints the dpeth of trees and number of leaves so that they train faster, while random forest has not restrictions on this parameters.

Adaboost default parameters:

- 1. n_estimator (default=50)(The maximum number of estimators at which boosting is terminated. In case of perfect fit, the learning procedure is stopped early.).
- 2. Learning rate (default=1)(Weight applied to each classifier at each boosting iteration.)
- 3. ...

Random Forest default parameters:

- 1. n_estimators (default=100)
- 2. max_depth=None
- 3. ...

6.3 Tuning and plotting

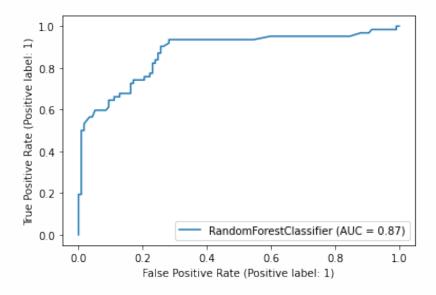
```
# parameters
params_rf = {
    'criterion': ['gini', 'entropy'],
    'max_depth': [5, 10, 15, 20],
    'min_samples_split': [2, 4, 8, 12, 16],
    'min_samples_leaf': [1, 5, 10]
}
params_ab = {
    'n_estimators': [5, 10, 20, 30, 50, 75, 100],
    'learning_rate': [0.001, 0.01, 0.1, 0.2, 0.5, 1]
}
scores={}
# hyperparameter tuning
rf search = GridSearchCV(RF clf, params rf, cv=5, scoring='accuracy')
ab_search = GridSearchCV(Ada_clf, params_ab, cv=5, scoring='accuracy')
rf_search.fit(X_train, y_train)
ab_search.fit(X_train, y_train)
rf_best_param = rf_search.best_params_
ab_best_param = ab_search.best_params_
rf_best_score = rf_search.best_score_
ab_best_score = ab_search.best_score_
print(f'Best parameters for RandomForest: {rf_best_param}')
print(
    f'Mean cross-validated accuracy score of the best_estimator: '+ \
    f'{rf best score:.3f}'
scores['RandomForest'] = rf_best_score
print('----')
print(f'Best parameters for AdaBoost: {ab best param}')
print(
    f'Mean cross-validated accuracy score of the best_estimator: '+ \
    f'{ab_best_score:.3f}'
scores['AdaBoost'] = ab_best_score
```

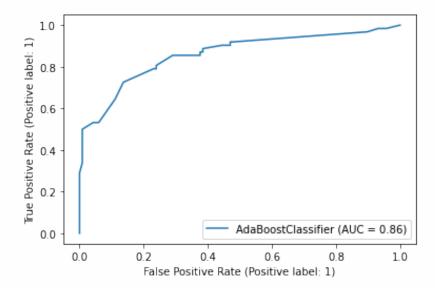
```
Best parameters for RandomForest: {'criterion': 'gini', 'max_depth': 20, 'min_samples_leaf': 5, 'min_samples_split': 2}

Mean cross-validated accuracy score of the best_estimator: 0.813
----

Best parameters for AdaBoost: {'learning_rate': 0.1, 'n_estimators': 75}

Mean cross-validated accuracy score of the best_estimator: 0.806
```





The ROC curves of two models looks very similar. And AUC score is also roughly the same.

Differences:

Random Forest use boostrap method to generate many trees with different splitting criterions, and use voting to finalize the tree model. While Adaboost combines several weak classifiers into one classifier, which would generalize better.

In this problem, the prediction result of two models are similar because the data is balanced and has resonable size. However, if sample size is small and noisy, random forest can leads to overfitting.

7 Generalized Additive Models