## Homework 3

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## **Question 1**

```
In [2]: import numpy as np
        import pandas as pd
        import matplotlib.pyplot as plt
        from matplotlib.colors import ListedColormap
        from sklearn import neighbors, datasets
        from sklearn import tree
        from sklearn.tree import DecisionTreeRegressor
        from sklearn import svm, datasets, neural network
        from sklearn import preprocessing
        from sklearn.model selection import cross val score, train test split, GridSea
        rchCV, RandomizedSearchCV
        from sklearn.ensemble import RandomForestClassifier, RandomForestRegressor, Ex
        traTreesRegressor, GradientBoostingRegressor
        from sklearn import linear model
        from sklearn.naive bayes import GaussianNB
        from sklearn.metrics import accuracy score, confusion matrix, precision recall
         fscore support, mean absolute error
        from sklearn.metrics import precision recall curve, average precision score, r
        oc curve, auc, mean squared error
        import warnings
        warnings.filterwarnings('ignore')
```

## **Data Loading**

```
In [3]: df = pd.read excel('HW3.xlsx', skiprows=1, header = None)
        df.head()
        features = df.iloc[:,0:22].values
        target = df.iloc[:,24].values
        x_train, x_test, y_train, y_test = train_test_split(features, target, test_siz
        e = 0.2
        features
Out[3]: array([[
                              0, ..., 3662,
                   1,
                        1,
                                               1,
                                                     0],
                              0, ..., 2900,
                                               1,
                                                     1],
                   2,
                         1,
                              0, ..., 3914,
                                                     0],
                   3,
                              0, ..., 3394,
                        1,
               [1998,
                                                     0],
               [1999,
                         1,
                              0, ..., 253,
                                               0,
                                                     1],
                              0, ..., 1844,
               [2000,
                         1,
                                                     0]], dtype=int64)
                                               0,
```

(a)

## **Linear Regression**

```
In [18]: | param_grid ={'fit_intercept':[True,False], 'normalize':[True,False], 'copy_X':
         [True, False]}
         clf_lr = linear_model.LinearRegression()
         grid = GridSearchCV(clf lr, param grid, cv = 10, scoring = 'neg mean absolute
         error')
         grid.fit(x_train,y_train)
         print (grid.best_score_)
         print (grid.best params )
         print (grid.best estimator )
         print("Mean Absolute Error: ",mean_absolute_error(y_test, grid.predict(x_test
         print("Mean Squared Error: ",mean_squared_error(y_test, grid.predict(x_test)))
         -76.84419544679477
         {'copy_X': True, 'fit_intercept': True, 'normalize': True}
         LinearRegression(copy X=True, fit intercept=True, n jobs=None, normalize=Tru
         e)
         Mean Absolute Error: 78.9282734607993
         Mean Squared Error: 17799.383455566836
```

### k-NN

```
In [27]: | clf knn = neighbors.KNeighborsRegressor()
         k range = list(range(1,20))
         weight options = ["uniform", "distance"]
         param_grid = dict(n_neighbors = k_range, weights = weight_options)
         grid = GridSearchCV(clf knn, param grid, cv = 10, scoring = 'neg mean absolute
         error')
         grid.fit(x_train,y_train)
         print (grid.best_score_)
         print (grid.best_params_)
         print (grid.best_estimator_)
         print("Mean Absolute Error: ",mean_absolute_error(y_test, grid.predict(x_test
         print("Mean Squared Error: ",mean_squared_error(y_test, grid.predict(x_test)))
         -94.94527793969209
         {'n_neighbors': 19, 'weights': 'distance'}
         KNeighborsRegressor(algorithm='auto', leaf size=30, metric='minkowski',
                   metric_params=None, n_jobs=None, n_neighbors=19, p=2,
                   weights='distance')
         Mean Absolute Error: 93.93644867800187
         Mean Squared Error: 26909.40331663546
```

# **Regression Tree**

```
In [25]: complexity values = range(1,25)
         max_nodes = [5, 15, None]
         min purity = [0, 0.5, .1]
         param grid = dict(max depth = complexity values,
                           max leaf nodes = max nodes, min impurity decrease = min puri
         ty)
         clf dt = tree.DecisionTreeRegressor()
         grid = GridSearchCV(clf dt, param grid, cv = 10, scoring = 'neg mean absolute
         error')
         grid.fit(x_train,y_train)
         print (grid.best score )
         print (grid.best_params_)
         print (grid.best_estimator_)
         print("Mean Absolute Error: ",mean absolute error(y test, grid.predict(x test
         )))
         print("Mean Squared Error: ",mean_squared_error(y_test, grid.predict(x_test)))
         -71.66875285786273
         {'max_depth': 6, 'max_leaf_nodes': None, 'min_impurity_decrease': 0.1}
         DecisionTreeRegressor(criterion='mse', max depth=6, max features=None,
                    max_leaf_nodes=None, min_impurity_decrease=0.1,
                    min impurity split=None, min samples leaf=1,
                    min_samples_split=2, min_weight_fraction_leaf=0.0,
                    presort=False, random state=None, splitter='best')
         Mean Absolute Error: 71.5365658351666
         Mean Squared Error: 16169.146597326464
```

# SVM Regression

```
In [ ]: Cs = [0.1, 1.0, 10] # SVM regularization parameter

Gammas = [0.01, 0.1, 1.0]
    param_grid = {'C': [0.1, 1.0, 10], 'gamma' : [0.01, 0.1, 1.0], 'kernel':['line ar', 'rbf']}
    grid = GridSearchCV(svm.SVR(), param_grid, cv=5, scoring = 'neg_mean_absolute_error')
    grid.fit(x_train, y_train)

print (grid.best_score_)
    print (grid.best_params_)
    print (grid.best_estimator_)
    print("Mean Absolute Error: ",mean_absolute_error(y_test, grid.predict(x_test)))
    print("Mean Squared Error: ",mean_squared_error(y_test, grid.predict(x_test)))
```

#### **Neural Network**

```
In [15]:
         mlpr = neural network.MLPRegressor()
         param_list = {"hidden_layer_sizes": [1,50], "activation": ["identity", "logist
         ic", "tanh", "relu"]}
         grid = GridSearchCV(estimator=mlpr, param grid=param list, cv = 5, scoring =
          'neg mean absolute error')
         grid.fit(x train, y train)
         print (grid.best score )
         print (grid.best params )
         print (grid.best_estimator_)
         print("Mean Absolute Error: ",mean_absolute_error(y_test, grid.predict(x_test
         )))
         print("Mean Squared Error: ",mean_squared_error(y_test, grid.predict(x_test)))
         -92.0630236947674
         {'activation': 'identity', 'hidden_layer_sizes': 50}
         MLPRegressor(activation='identity', alpha=0.0001, batch_size='auto',
                beta_1=0.9, beta_2=0.999, early_stopping=False, epsilon=1e-08,
                hidden layer sizes=50, learning_rate='constant',
                learning rate init=0.001, max iter=200, momentum=0.9,
                n_iter_no_change=10, nesterovs_momentum=True, power_t=0.5,
                random state=None, shuffle=True, solver='adam', tol=0.0001,
                validation fraction=0.1, verbose=False, warm start=False)
         Mean Absolute Error: 107.13040362760985
         Mean Squared Error: 29162.450981650185
```

### **Ensemble Methods - Random Forest**

```
In [20]: n estimators = [int(x) \text{ for } x \text{ in np.linspace(start} = 200, stop = 2000, num = 10)
          ) ]
          max features = ['auto', 'sqrt']
          max depth = [int(x) for x in np.linspace(10, 110, num = 11)]
          max depth.append(None)
          min samples split = [2, 5, 10]
          min samples leaf = [1, 2, 4]
          bootstrap = [True, False]
          random_grid = {'n_estimators': n_estimators,
                         'max_features': max_features,
                         'max depth': max depth,
                         'min_samples_split': min_samples_split,
                         'min_samples_leaf': min_samples_leaf,
                         'bootstrap': bootstrap}
          rf = RandomForestRegressor()
          rf random = RandomizedSearchCV(estimator = rf, param distributions = random gr
          id, n_iter = 100, cv = 3, verbose=2, random_state=42, n_jobs = -1)
          rf_random.fit(x_train, y_train)
          print (grid.best score )
          print (grid.best_params_)
          print (grid.best estimator )
          print("Mean Absolute Error: ",mean_absolute_error(y_test, grid.predict(x_test
          print("Mean Squared Error: ",mean squared error(y test, grid.predict(x test)))
```

Fitting 3 folds for each of 100 candidates, totalling 300 fits

```
[Parallel(n jobs=-1)]: Using backend LokyBackend with 4 concurrent workers.
48.6s
[Parallel(n jobs=-1)]: Done 154 tasks
                                      | elapsed: 3.8min
[Parallel(n jobs=-1)]: Done 300 out of 300 | elapsed: 7.2min finished
-76.84419544679477
{'copy_X': True, 'fit_intercept': True, 'normalize': True}
LinearRegression(copy X=True, fit intercept=True, n jobs=None, normalize=Tru
Mean Absolute Error: 78.9282734607993
Mean Squared Error: 17799.383455566836
```

(b)

### **Data Transformation**

```
In [10]: df2 = pd.read_excel('HW3.xlsx', skiprows=1, header = None)
    df2 = df2[df2[23]==1]
    df2.head()
    features = df.iloc[:,0:22].values
    target = df.iloc[:,24].values
    x_train2, x_test2, y_train2, y_test2 = train_test_split(features, target, test
    _size= 0.2)
```

### **Linear Regression**

```
param_grid ={'fit_intercept':[True,False], 'normalize':[True,False], 'copy_X':
In [21]:
         [True, False]}
         clf lr = linear model.LinearRegression()
         grid = GridSearchCV(clf_lr, param_grid, cv = 10, scoring = 'neg_mean_absolute_
         error')
         grid.fit(x_train2,y_train2)
         print (grid.best score )
         print (grid.best_params_)
         print (grid.best_estimator_)
         print("Mean Absolute Error: ",mean_absolute_error(y_test2, grid.predict(x_test
         2)))
         print("Mean Squared Error: ",mean squared error(y test2, grid.predict(x test2
         )))
         -79.04038025213856
         {'copy_X': True, 'fit_intercept': False, 'normalize': True}
         LinearRegression(copy X=True, fit intercept=False, n jobs=None,
                  normalize=True)
         Mean Absolute Error: 74.88739227110489
         Mean Squared Error: 13981.918741906476
```

### K-NN

```
In [22]: | clf knn = neighbors.KNeighborsRegressor()
         k range = list(range(1,20))
         weight options = ["uniform", "distance"]
         param_grid = dict(n_neighbors = k_range, weights = weight_options)
         grid = GridSearchCV(clf knn, param grid, cv = 10, scoring = 'neg mean absolute
         error')
         grid.fit(x_train2,y_train2)
         print (grid.best_score_)
         print (grid.best_params_)
         print (grid.best_estimator_)
         print("Mean Absolute Error: ",mean_absolute_error(y_test2, grid.predict(x_test
         2)))
         print("Mean Squared Error: ",mean_squared_error(y_test2, grid.predict(x_test2
         )))
         -95.6114089211818
         {'n_neighbors': 19, 'weights': 'distance'}
         KNeighborsRegressor(algorithm='auto', leaf_size=30, metric='minkowski',
                   metric params=None, n jobs=None, n neighbors=19, p=2,
                   weights='distance')
         Mean Absolute Error: 93.55243906949673
         Mean Squared Error: 24551.642715824404
```

## **Regression Tree**

```
In [23]: complexity values = range(1,25)
         max_nodes = [5, 15, None]
         min purity = [0, 0.5, .1]
         param grid = dict(max depth = complexity values,
                           max leaf nodes = max nodes, min impurity decrease = min puri
         ty)
         clf dt = tree.DecisionTreeRegressor()
         grid = GridSearchCV(clf dt, param grid, cv = 10, scoring = 'neg mean absolute
         error')
         grid.fit(x_train2,y_train2)
         print (grid.best score )
         print (grid.best_params_)
         print (grid.best_estimator_)
         print("Mean Absolute Error: ",mean absolute error(y test2, grid.predict(x test
         2)))
         print("Mean Squared Error: ",mean_squared_error(y_test2, grid.predict(x_test2)
         )))
         -76.78214699475116
         {'max depth': 5, 'max leaf nodes': None, 'min impurity decrease': 0.5}
         DecisionTreeRegressor(criterion='mse', max_depth=5, max_features=None,
                    max leaf nodes=None, min impurity decrease=0.5,
                    min impurity split=None, min samples leaf=1,
                    min_samples_split=2, min_weight_fraction_leaf=0.0,
                    presort=False, random state=None, splitter='best')
         Mean Absolute Error: 76.77525493962408
         Mean Squared Error: 20168.69343613505
```

# **SVM Regression**

```
In []: Cs = [ 0.1, 1.0, 10] # SVM regularization parameter

Gammas = [0.01, 0.1, 1.0]
    param_grid = {'C': [ 0.1, 1.0, 10], 'gamma' : [0.01, 0.1, 1.0], 'kernel':['lin ear', 'rbf']}
    grid = GridSearchCV(svm.SVR(), param_grid, cv=5, scoring = 'neg_mean_absolute_error')
    grid.fit(x_train2, y_train2)

print (grid.best_score_)
    print (grid.best_params_)
    print (grid.best_estimator_)
    print("Mean Absolute Error: ",mean_absolute_error(y_test2, grid.predict(x_test 2)))
    print("Mean Squared Error: ",mean_squared_error(y_test2, grid.predict(x_test2)))
```

#### **Neural Network**

```
In [24]:
         mlpr = neural network.MLPRegressor()
         param_list = {"hidden_layer_sizes": [1,50], "activation": ["identity", "logist
         ic", "tanh", "relu"]} #, "solver": ["lbfqs", "sqd", "adam"], "alpha": [0.0000
         5,0.0005]}
         grid = GridSearchCV(estimator=mlpr, param grid=param list, cv = 5,scoring = 'n
         eg mean absolute error')
         grid.fit(x_train2, y_train2)
         print (grid.best score )
         print (grid.best_params_)
         print (grid.best estimator )
         print("Mean Absolute Error: ",mean_absolute_error(y_test, grid.predict(x_test
         )))
         print("Mean Squared Error: ",mean squared error(y test, grid.predict(x test)))
         -92.54430459667809
         {'activation': 'identity', 'hidden_layer_sizes': 50}
         MLPRegressor(activation='identity', alpha=0.0001, batch size='auto',
                beta 1=0.9, beta 2=0.999, early stopping=False, epsilon=1e-08,
                hidden_layer_sizes=50, learning_rate='constant',
                learning_rate_init=0.001, max_iter=200, momentum=0.9,
                n iter no change=10, nesterovs momentum=True, power t=0.5,
                random state=None, shuffle=True, solver='adam', tol=0.0001,
                validation fraction=0.1, verbose=False, warm start=False)
         Mean Absolute Error: 85.57882245920304
         Mean Squared Error: 27831.225290051825
```

### **Ensemble Methods - Random Forest**

```
In [25]: n estimators = [int(x) for x in np.linspace(start = 200, stop = 2000, num = 10)]
         max features = ['auto', 'sqrt']
         max depth = [int(x) for x in np.linspace(10, 110, num = 11)]
         max depth.append(None)
         min samples split = [2, 5, 10]
         min samples leaf = [1, 2, 4]
         bootstrap = [True, False]
         random_grid = {'n_estimators': n_estimators,
                         'max_features': max_features,
                         'max depth': max depth,
                         'min_samples_split': min_samples_split,
                         'min_samples_leaf': min_samples_leaf,
                         'bootstrap': bootstrap}
         rf = RandomForestRegressor()
         rf random = RandomizedSearchCV(estimator = rf, param distributions = random gr
         id, n_iter = 100, cv = 3, verbose=2, random_state=42, n_jobs = -1)
         rf_random.fit(x_train2, y_train2)
         print (grid.best score )
         print (grid.best_params_)
         print (grid.best estimator )
         print("Mean Absolute Error: ",mean_absolute_error(y_test2, grid.predict(x_test
         print("Mean Squared Error: ",mean squared error(y test2, grid.predict(x test2
         )))
         Fitting 3 folds for each of 100 candidates, totalling 300 fits
         [Parallel(n jobs=-1)]: Using backend LokyBackend with 4 concurrent workers.
         [Parallel(n jobs=-1)]: Done 33 tasks
                                                     | elapsed:
                                                                  41.6s
         [Parallel(n jobs=-1)]: Done 154 tasks
                                                     | elapsed: 3.5min
         [Parallel(n jobs=-1)]: Done 300 out of 300 | elapsed: 6.6min finished
         -92.54430459667809
         {'activation': 'identity', 'hidden_layer_sizes': 50}
         MLPRegressor(activation='identity', alpha=0.0001, batch_size='auto',
```

```
beta_1=0.9, beta_2=0.999, early_stopping=False, epsilon=1e-08,
       hidden_layer_sizes=50, learning_rate='constant',
       learning rate init=0.001, max iter=200, momentum=0.9,
       n iter no change=10, nesterovs momentum=True, power t=0.5,
       random state=None, shuffle=True, solver='adam', tol=0.0001,
       validation fraction=0.1, verbose=False, warm start=False)
Mean Absolute Error: 73.66728669313828
Mean Squared Error: 20557.77195385305
```

(c)

As of 4:45PM, my SVM models have not finished running, despite doing so for the last 2.5 hours. In the interest of time, I'll be elminiating these models on the sheer basis of being computationally inefficient. I believe the code for SVM Regression runs properly in both scenarios, but I do not have the resources to generate results for it.

That being said, I managed to successfully generate 5 of the models.

In part A, the model that generated the best results was the Regression Tree, with a mean absolute error of 71. This was followed by both the Linear Regression and Random Forest models which for some reason generated identical results.

In part B, model with the best results was the Random Forest Model with a Mean Absolute Error of 73. However, the Linear Regression and Regression Tree models both generate a lower Mean Squared Error, which punishes larger errors. The choice for part B depends on what the end user values more.

In general, it appears that Part B results generated less Mean Absolute Error and Mean Squared Error. This may be because removing all the Non-Purchase variables also removed all the 0 values from the target variables. This removed any skew that these number may have been having on the results and allowed the model to fit better.

## Question 2

## **Data Loading and Transformation**

```
sb df = pd.read csv("spambase.data", header = None)
In [38]:
           sb df.head()
Out[38]:
                  0
                        1
                             2
                                  3
                                              5
                                                         7
                                                               8
                                                                              48
                                                                                    49
                                                                                         50
                                                                                                51
                                                                                                       52
                                                    6
                                                                     9
               0.00
                     0.64
                           0.64
                                0.0
                                     0.32
                                           0.00
                                                 0.00
                                                      0.00
                                                            0.00
                                                                  0.00
                                                                            0.00
                                                                                 0.000
                                                                                         0.0
                                                                                             0.778
                                                                                                    0.000
                           0.50
                                           0.28
                                                            0.00
                                                                  0.94
               0.21
                     0.28
                                0.0
                                     0.14
                                                 0.21
                                                      0.07
                                                                            0.00
                                                                                 0.132
                                                                                        0.0
                                                                                             0.372
                                                                                                    0.180
               0.06
                    0.00
                           0.71
                                0.0
                                     1.23
                                          0.19
                                                 0.19
                                                      0.12
                                                            0.64
                                                                  0.25
                                                                            0.01
                                                                                 0.143 0.0
                                                                                             0.276
                                                                                                    0.184
                     0.00
                           0.00
                                0.0
                                     0.63
                                           0.00
                                                 0.31
                                                       0.63
                                                            0.31
                                                                  0.63
                                                                            0.00
                                                                                 0.137
                                                                                             0.137
                                                                                                    0.000
               0.00 \quad 0.00 \quad 0.00 \quad 0.0 \quad 0.63 \quad 0.00 \quad 0.31 \quad 0.63 \quad 0.31 \quad 0.63
                                                                           0.00 0.135 0.0
                                                                                            0.135
                                                                                                    0.000
           5 rows × 58 columns
           features = sb df.iloc[:,0:56].values
In [39]:
           target = sb df.iloc[:,57].values
            x_train, x_test, y_train, y_test = train_test_split(features, target, test_siz
           e= 0.2, stratify = target)
```

### **Decision Tree**

```
In [40]:
         complexity values = range(1,25)
         weight_options = ["gini", "entropy"]
         max nodes = [5, 15, None]
         min purity = [0, 0.5, .1]
         param_grid = dict(max_depth = complexity_values, criterion = weight_options,
                           max leaf nodes = max nodes, min impurity decrease = min puri
         ty)
         clf_dt = tree.DecisionTreeClassifier()
         grid = GridSearchCV(clf_dt, param_grid, cv = 10, scoring = 'accuracy')
         grid.fit(x_train,y_train)
         print (grid.best_score_)
         print (grid.best params )
         print (grid.best_estimator_)
         print("Prediction Accuracy: ",accuracy_score(y_test, grid.predict(x_test)))
         0.9192934782608696
         {'criterion': 'gini', 'max depth': 11, 'max leaf nodes': None, 'min impurity
         decrease': 0}
         DecisionTreeClassifier(class_weight=None, criterion='gini', max_depth=11,
                     max features=None, max leaf nodes=None,
                     min impurity decrease=0, min impurity split=None,
                     min samples leaf=1, min samples split=2,
                     min weight fraction leaf=0.0, presort=False, random state=None,
                     splitter='best')
         Prediction Accuracy: 0.9239956568946797
In [41]: | clf dt = grid.fit(x train,y train)
         y predict = clf dt.predict(x test)
         confusion matrix(y test,y predict)
Out[41]: array([[526, 32],
                [ 35, 328]], dtype=int64)
In [42]: precision recall fscore support(y test, y predict)
Out[42]: (array([0.93761141, 0.91111111]),
          array([0.94265233, 0.90358127]),
          array([0.94012511, 0.90733057]),
          array([558, 363], dtype=int64))
```

## Logistic Regression

```
In [43]: | param_grid ={'C': [0.001, 0.01, 0.1, 1, 10, 100, 1000]}
         clf lr = linear model.LogisticRegression()
         grid = GridSearchCV(clf lr, param grid, cv = 10, scoring = 'accuracy')
         grid.fit(x_train,y_train)
         print (grid.best score )
         print (grid.best_params_)
         print (grid.best estimator )
         print("Prediction Accuracy: ",accuracy_score(y_test, grid.predict(x_test)))
         0.925
         {'C': 1}
         LogisticRegression(C=1, class weight=None, dual=False, fit intercept=True,
                   intercept_scaling=1, max_iter=100, multi_class='warn',
                   n_jobs=None, penalty='12', random_state=None, solver='warn',
                   tol=0.0001, verbose=0, warm start=False)
         Prediction Accuracy: 0.9305103148751357
         clf lr = grid.fit(x train,y train)
In [44]:
         y_predict = clf_lr.predict(x_test)
         confusion_matrix(y_test,y_predict)
Out[44]: array([[537, 21],
                [ 43, 320]], dtype=int64)
In [45]: precision recall fscore support(y test, y predict)
Out[45]: (array([0.92586207, 0.93841642]),
          array([0.96236559, 0.8815427]),
          array([0.94376098, 0.90909091]),
          array([558, 363], dtype=int64))
```

## K- Nearest Neighbor

```
In [46]: | clf knn = neighbors.KNeighborsClassifier()
         k range = list(range(1,20))
         weight options = ["uniform", "distance"]
         param_grid = dict(n_neighbors = k_range, weights = weight_options)
         grid = GridSearchCV(clf knn, param grid, cv = 10, scoring = 'accuracy')
         grid.fit(x train,y train)
         print (grid.best score )
         print (grid.best_params_)
         print (grid.best_estimator_)
         print("Prediction Accuracy: ",accuracy_score(y_test, grid.predict(x_test)))
         0.9160326086956522
         {'n_neighbors': 6, 'weights': 'distance'}
         KNeighborsClassifier(algorithm='auto', leaf_size=30, metric='minkowski',
                    metric params=None, n jobs=None, n neighbors=6, p=2,
                    weights='distance')
         Prediction Accuracy: 0.9098805646036916
In [47]: | clf_knn = grid.fit(x_train,y_train)
         y_predict = clf_knn.predict(x_test)
         confusion_matrix(y_test,y_predict)
Out[47]: array([[525, 33],
                [ 50, 313]], dtype=int64)
In [48]: | precision_recall_fscore_support(y_test, y_predict)
Out[48]: (array([0.91304348, 0.90462428]),
          array([0.94086022, 0.86225895]),
          array([0.92674316, 0.88293371]),
          array([558, 363], dtype=int64))
```

#### **SVM**

```
In [54]: Cs = [0.001, 0.01, 0.1, 1.0, 10] # SVM regularization parameter
         Gammas = [0.001, 0.01, 0.1, 1.0]
         param grid = \{'C': [0.001, 0.01, 0.1, 1.0, 10], 'gamma': [0.001, 0.01, 0.1,
         1.0], 'kernel':['linear', 'rbf']}
         grid = GridSearchCV(svm.SVC(), param_grid, cv=10, scoring = 'accuracy')
         grid.fit(x train, y train)
         print (grid.best score )
         print (grid.best_params_)
         print (grid.best estimator )
         print("Prediction Accuracy: ",accuracy_score(y_test, grid.predict(x_test)))
         0.9385869565217392
         {'C': 10, 'gamma': 0.01, 'kernel': 'rbf'}
         SVC(C=10, cache size=200, class weight=None, coef0=0.0,
           decision_function_shape='ovr', degree=3, gamma=0.01, kernel='rbf',
           max_iter=-1, probability=False, random_state=None, shrinking=True,
           tol=0.001, verbose=False)
         Prediction Accuracy: 0.9283387622149837
         clf svm linear = grid.fit(x train,y train)
In [55]:
         y predict = clf svm linear.predict(x test)
         confusion_matrix(y_test,y_predict)
Out[55]: array([[528, 30],
                [ 36, 327]], dtype=int64)
In [56]: precision recall fscore support(y test, y predict)
Out[56]: (array([0.93617021, 0.91596639]),
          array([0.94623656, 0.90082645]),
          array([0.94117647, 0.90833333]),
          array([558, 363], dtype=int64))
```

#### **Model Selections**

**No Misclassification Costs** 

Out of the four classfication models I generated (Decision Tree, Logistic Regression, K-NN, and SVM), the one that produced the highest accuracy is SVM. The overall accuracy however, is not necessarily the most useful metric when comparin the models. The accuracy of SVM also isn't much higher than the other three. Precision and Recall are also important metrics to look at, and these metrics measure the following.

Precision: Out of how many emails we identified as spam, how many were truly spam.

Recall: Our of all the emails that are truly spam, how many of them did we correctly identify.

In this case, I'll be using recall as the more important metric. Emails that get flagged as spam can be further reviewed by a user checking the spam inbox, while failed identification can lead to annoyances at best and security issues or scams at worst. The Logisitic Regression model in this case has the highest recall at 96% without trading off much of the accuracy from SVM. Both accuracies are in the 94% range, and the precision also is not that much lower. Without misclassification costs, I would choose Logistic Regression.

#### With Misclassification Costs

With Misclassification costs, the game changes a little bit. False positives are the emails that the model identifies as spam, but are not, while false negatives are emails in which the model says are not spam but really are. The more important one depends on what standpoint you're looking from. False positives mean that a user might miss some important emails that get flagged into your spam box and false negatives might pose a security risk with the severity depending on the context.

In this case, I'll treat false negatives as the more important case. A rational user would occassionally check their spam box for any emails that got unfortunately flagged if they were expecting something important, while they would likely trust the spam filter to catch spam emails and could be caught off guard by malicious emails even if they're being careful. With my experience in information security, I see the latter case being more common and costly than the first case, and I prefer erring on the side of caution.

We can apply the cost ratio of 10:1 to each of the confusion matricies generated by the models. From there we can calculate the misclassification costs for each of the models.

Decision Tree: 382

Logistic Regression: 451

KNN: 533

SVM: 390

In this case, the model that generated the lowest cost was The decision tree, with SVM coming in a close second. The logisitic regression model which I chose for the evaluation without the cost produced a much higher misclassification cost. In choosing between the Decision Tree and SVM models however, the SVM model shows a greater accuracy and similar precision, recall, and fscore measure while not trading off as much cost. With this information, I would give the edge to SVM, even if Decision Tree costs slightly lower.