## Person of Interest Identifier

This project is based on the Enron scandal (https://en.wikipedia.org/wiki/Enron\_scandal) of 2001, the data used here is the email and financial data of enron which was made public and can be found here (https://www.cs.cmu.edu/~./enron/enron\_mail\_20150507.tar.gz). Here we identify people from the numerous enron employees which can be considered as 'person of interest (poi)' i.e. who may have a hand in the scandal. The data has handpicked people classified as poi which were convicted in reality. We use a supervised learning approach to build our poi identifier.

We begin by first importing the required packages. Here we mainly work with the scikit-learn package.

#### In [12]:

```
import os
import sys
import pickle
from feature_format import featureFormat, targetFeatureSplit
from tester import dump_classifier_and_data,test_classifier
from sklearn.cross_validation import train_test_split
from sklearn.decomposition import PCA
from sklearn.feature_selection import SelectKBest
from sklearn.naive_bayes import GaussianNB
from sklearn.svm import SVC
from sklearn.ensemble import RandomForestClassifier
from sklearn.model_selection import GridSearchCV
import matplotlib.pyplot as plt
```

## Next we load our dataset.

## In [3]:

```
with open("final_project_dataset.pkl", "rb") as data_file:
    data_dict = pickle.load(data_file)
my_dataset = data_dict
```

Initially we start with all the features in our data however these may contain noise and give a poor accuracy which will be optimized later. Further we extract the features and labels from this data for analysis. Finally we split these in training and testing data for our model.

#### In [4]:

To get an initial look at our validation figures we use the simple Naive Bayes classifier without any preprocessing. We use a test\_classifier function which evaluates and displays evaluation metrics.

## In [10]:

As we can see we get a fairly low accuracy and recall score owing to the many features contributing as noise. Our first step towards improving these would be using Principal Component Analysis in which the components with maximum variance can be selected while others can be dropped.

## In [35]:

```
num = 12 #This can be tuned to select different number of principal components
print(data.shape)
pca = PCA(n_components = num)
pca.fit(data)
dat = pca.transform(data)
print(dat.shape)

(145, 15)
(145, 12)
```

To further augment our selection we use feature selection. Here we use SelectKBest which selects the 'K' best features which explain the data.

#### In [36]:

```
import warnings
warnings.filterwarnings('ignore') #so that warnings aren't printed

labels, features = targetFeatureSplit(dat)
features_train, features_test, labels_train, labels_test = train_test_split(features, label)
selector = SelectKBest(k=7) #To limit the number of features to seven (i.e. approximately hereatures_train = selector.fit_transform(features_train, labels_train)

mask = selector.get_support()
new_features = []
#Here we printout our K best features
for bool, feature in zip(mask, features_list):
    if bool:
        new_features.append(feature)
print(new_features)
```

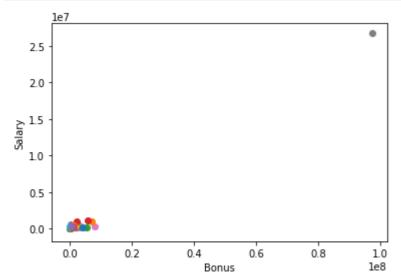
```
['loan_advances', 'bonus', 'restricted_stock_deferred', 'deferred_income', 'total_stock_value', 'expenses', 'exercised_stock_options']
```

Other reasons for poor performance can be Outliers present in the data. This can be investigated by using visualization of our data.

## In [34]:

```
with open("final_project_dataset.pkl", "rb") as data_file:
    data_dict = pickle.load(data_file)
features_list = ['bonus','salary'] #Here we choose bonus & salary for visualization
data = featureFormat(data_dict,features_list)

for point in data:
    bonus = point[0]
    salary = point[1]
    plt.scatter(bonus,salary)
plt.xlabel('Bonus')
plt.ylabel('Salary')
plt.show()
```



We see the presence of a clear outlier in the data which has been hampering our results. We remove this from our original data and proceed with our analysis.

## In [37]:

```
#To detect the outlier:
for i in data_dict:
   if (data_dict[i]['salary'] != "NaN") & (data_dict[i]['bonus'] != "NaN"):
      if (data_dict[i]['salary'] >= 5000000) & (data_dict[i]['bonus'] >= 5000000):
           print("Outlier present is ",i)
```

Outlier present is TOTAL

The outlier was actually the total row with total figures for all the enron employees. We now remove this for further analysis.

#### In [13]:

The next classifier we intend to try is SVM. However, SVMs in general are affected by different ranges of the features and may result in output being dominated by a single feature. To overcome this issue we perform feature scaling, in which we first scale our features to an uniform scale of [0,1].

#### In [14]:

```
print('Before scaling = ',data[0,3])
from sklearn.preprocessing import MinMaxScaler
scaler = MinMaxScaler()
scaler.fit(data)
scaled_data = scaler.transform(data)
print('After scaling = ',scaled_data[0,3]) #Example showing feature scaling effect
labels, features = targetFeatureSplit(scaled_data)

features_train, features_test, labels_train, labels_test = train_test_split(features, label)

Before scaling = -126027.0
After scaling = 0.09634567351381695
```

Now we build our SVM classifier and train it on our data.

```
In [15]:
```

```
clf = SVC()
clf.fit(features_train,labels_train)
test_classifier(clf,data_dict,features_list)

Got a divide by zero when trying out: SVC(C=1.0, cache_size=200, class_weigh
t=None, coef0=0.0,
    decision_function_shape='ovr', degree=3, gamma='auto', kernel='rbf',
    max_iter=-1, probability=False, random_state=None, shrinking=True,
    tol=0.001, verbose=False)
Precision or recall may be undefined due to a lack of true positive predicit
ons.
```

The result from the SVM classifier show that it doesn't perform well here as the number of true positives is very less.

Next we try out the Random Forest algorithm with a different approach. Here we use GridSearch technique to find parameters which yield the best performance for our classifier.

## In [25]:

```
labels, features = targetFeatureSplit(data)
features_train, features_test, labels_train, labels_test = train_test_split(features, label

rfc = RandomForestClassifier()
parameters = {'n_estimators':[10,50,100], 'max_features':('auto','log2')}

clf = GridSearchCV(rfc,parameters)
clf.fit(features_train,labels_train)
print(clf.best_params_)

{'max_features': 'log2', 'n_estimators': 50}
```

The above parameters are tested seperately below as test\_classifier doesn't work well with GridSearchCV.

```
In [27]:
```

```
clf = RandomForestClassifier(n_estimators=50, max_features='log2')
clf.fit(features_train,labels_train)
test classifier(clf,data dict,features list)
RandomForestClassifier(bootstrap=True, class_weight=None, criterion='gini',
            max_depth=None, max_features='log2', max_leaf_nodes=None,
            min_impurity_decrease=0.0, min_impurity_split=None,
            min_samples_leaf=1, min_samples_split=2,
            min weight fraction leaf=0.0, n estimators=50, n jobs=1,
            oob score=False, random state=None, verbose=0,
            warm_start=False)
        Accuracy: 0.86957
                                Precision: 0.59977
                                                         Recall: 0.26150 F1:
0.36421 F2: 0.29475
        Total predictions: 14000
                                        True positives: 523
                                                                 False positi
ves: 349
                False negatives: 1477
                                        True negatives: 11651
```

The result shows that this classifier performs fairly well, which can be clearly seen from the above scores. A huge reason for this is here we perform parameter tunning by testing different parameters

and selecting the best ones. Changing parameters changes the performance of the classifier and can have impactful effects on its performance.

Speaking about the evaluation metrics, creating a classifier has no meaning until its performace is tested on data other than the data on which it was trained. This is called as validation, the train\_test\_split we have been performing, splits the data for training and testing for this purpose.

We obviously cannot rely on accuracy as a sole measure as it can be doctored easily by skewness present in our data. Hence we use a variety of metrics with accuracy to evaluate our classifier. The precision shows how well the classifier predicts a person of interest i.e. how well the classifier identifies a person as a poi and it is correct. On the other hand recall shows the fraction of total true cases, i.e. the probability of the classifier detecting a poi provided it is actually a poi.

Finally, another approach that is often used is creating new features out of the existing features that may help contribute towards the classification. Looking at our features, here we design a new feature called 'savings' which will be the net amout with a person including their salary and bonus and after excluding their expenses, this residual amount with a person can help in identifying their financial status which in this sense could mean their involvement in the fraud.

## In [17]:

```
with open("final project dataset.pkl", "rb") as data file:
    data_dict = pickle.load(data_file)
#New feature 'savings' created as follows
for key,i in data_dict.items():
    if i['salary']=='NaN':
        i['salary']=0
    if i['bonus']=='NaN':
        i['bonus']=0
    if i['expenses']=='NaN':
        i['expenses']=0
    i['savings'] = i['salary']+i['bonus']-i['expenses']
data dict.pop('TOTAL') #Outlier removal
my_dataset = data_dict
#Our feature list this time will contain our new feature which will replace the two feature
features_list = ['poi', 'loan_advances', 'restricted_stock_deferred',
                  'deferred_income', 'total_stock_value', 'savings',
                 'exercised stock options']
data = featureFormat(my_dataset, features_list, sort_keys = True)
labels, features = targetFeatureSplit(data)
features_train, features_test, labels_train, labels_test = train_test_split(features, label
```

We next validate this new feature with our classifier and observe minor improvemets in accuracy.

However when using our new feature list with the earlier Naive Bayes classifier, improvements in classification are clearly visible.

#### In [18]:

```
clf = RandomForestClassifier(n_estimators=50,max_features='log2')
clf.fit(features_train,labels_train)
test_classifier(clf,data_dict,features_list)
```

ves: 441 False negatives: 1465 True negatives: 12559

# In [19]:

```
clf = GaussianNB()
clf.fit(features_train,labels_train)
test_classifier(clf,data_dict,features_list)
```

GaussianNB(priors=None)

Accuracy: 0.77387 Precision: 0.28236 Recall: 0.45150 F1:

0.34744 F2: 0.40320

Total predictions: 15000 True positives: 903 False positi

ves: 2295 False negatives: 1097 True negatives: 10705

Thus we have performed analysis on the Enron dataset classifying the employees as person of interests. Another approach which can be utilized for the same is performing text learning on the text data i.e. Enron email data.