Project 5: The Enron Dataset ¶

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The Data Set & Outliers

There are 146 people in the Enron Dataset, 21 features each, 18 POI (persons of interest). A tossed out three data points as being unusual artifacts. I did not throw out outliers because they generally seemed to have useful information to identify POIs.

Feature Selection & Engineering

I selected 7 features from the original dataset, and created 3 new ones. Looking at correlation to the POI label (from the correlation matrix below), it seems that my new features are not particularly useful for classification. And in fact, 3-5 of best (selectKbest) of the original 7 features seem to be good enough to do the job. If I were to spend more time on this project I would engineer additional features and determine what would be the smallest number of features that would yield good results. Since there was not a huge number of features, I did not do PCA.

Training & Testing Classifiers

As a starting point, I tried NB, DTC, KNN classifiers a few times using different test sizes. NB worked OK, DTC did not, KNN worked well. So I fcoused on KNN for in my grid search. I ran gridSearchCSV with 10, 5, 3, and 1 features with test sizes of 0.3 and 0.1. Of these 10 searches, 5 resulted in classifier 1 and 5 resulted in classifier 2. Classifier 1 had a better F1 score (0.45 vs. 0.37) so I called this my best classifier and wrote it to the pkl file.

For reasons discussed here, it seems like it isn't a good idea to use a pipeline to search through a range of selectKbest features: https://discussions.udacity.com/t/selectkbest-pca-and-pipelines/22986/9 (https://discussions.udacity.com/t/selectkbest-pca-and-pipelines/22986/9

I suppose I could have written a loop to try to the entire range, but I just picked 10, 5, 3, 1 features as stated above.

```
gridSearchCSV() param_grid = {"knnn_neighbors": [1, 3, 5, 8, 10, 12, 14], "knnalgorithm": ["auto", "ball_tree", "kd_tree", "brute"], "knnleaf_size": range(3,10,1), "knnp": [1,2] }
```

```
Classifier 1 KNeighborsClassifier(algorithm='auto', leaf_size=3, metric='minkowski', metric_params=None, n_jobs=1, n_neighbors=1, p=1, weights='uniform')
F1 = 0.45
```

```
Classifier 2 KNeighborsClassifier(algorithm='auto', leaf_size=3, metric='minkowski', metric_params=None, n_jobs=1, n_neighbors=1, p=1, weights='uniform')
F1 = 0.37
```

I also ran AdaBoost using NB and DTC as the base classifiers (KNN is incompatible with AdaBoost). The resulting classifiers had lower F1 scores than Classifier 1, so I stuck with that one.

```
DTC = DecisionTreeClassifier(random_state = 42, max_features = "auto", class_weight = "balanced",max_depth = None) ada = AdaBoostClassifier(base_estimator=DTC)

param_grid = {"adabase_estimatorcriterion" : ["gini", "entropy"], "adabase_estimatorsplitter" : ["best", "random"], "ada__n_estimators": [1, 10, 50, 100, 200] }
```

AdaBoostClassifier(algorithm='SAMME.R', base_estimator=DecisionTreeClassifier(class_weight='balanced', criterion='entropy', max_depth=None, max_features='auto', max_leaf_nodes=None, min_impur...dom_state=42, splitter='random'), learning_rate=1.0, n_estimators=10, random_state=None))]) {'adabase_estimator} i'random', 'adan_estimators':10'adabase_estimator__criterion': 'entropy'} F1 = 0.32

```
NB = GaussianNB() adaN = AdaBoostClassifier(base_estimator=NB)

param_grid = { "ada__n_estimators": [1, 10, 50, 100, 200] }

AdaBoostClassifier(algorithm='SAMME.R', base_estimator=GaussianNB(priors=None), learning_rate=1.0, n_estimators=1, random_state=None))]) {'ada__n_estimators': 1}

F1 = 0.25
```

Tuning & Validation

By choosing different classifiers and trying parameters, it is possible to get a classifier that works really well (has very high scores) on the test data. This is called tuning, and is part of the art of machine learning. A classifier out of the box may not do a good job of correctly labeling data, but with parameter turning we can improve this.

But if we try to optimize too much, we can be guilty of overfitting (overtuning). In fact we could manually create a classifier that perfectly fits the training data with an absurd amount of overfit. Because of possible idiosyncracies and outliers in the training data, it is important to consider possible overfitting.

After fitting the classifier on training data, we test (or validate) it with test data before releasing it into the real world. It is a fairly easy thing, with overtuning, to create a classifier that works very well on the training data. But if we don't test the classifier on other data, we may have built a pretty, but useless, classifier. Among other things, outliers in the training data in can skew the classifier fit. Outliers in the testing data can skew the validations scores.

Validation can be a bit tricky if there is a limited data, because we don't have much data to train on, much less allocate as between training and test data. This is definitely true of our data set, with only 143 people and 18 POIs). A large number of features relative to the number of data points can be a problem as well, though we can address this to some degree with PCA.

Cross validation techniques help us avoid this, by radomly distributing and redistributing data between training and test, we can avoid classifier training problems such as overfitting.

I used StratifiedShuffleSplit() with 100 splits. This means that there were 100 randomized train-test splits. The shuffling balances data between training and testing. It also randomzies data that might be ordered. I also could also have used .KFold, but the random shuffling of StratifiedShuffleSplit() is useful because of our small data set.

Stratification is the process of distributing class labels relatively evenly between testing and training to prevent bias in either. StratifiedShuffleSplit() and StratifiedKFold() do this for us.

Precision & Recall

Since the label is POI:

precision = POI true positives / (POI true positives + POI false positives)

recall = POI true positives / (POI true positives + POI false negatives)

Thus if our classifier identified 10 POIs, with 5 true positives and 5 false positives, precision would be 0.5

If the ground state truth was 12 POIs, where we correctly identified 5 but 7 were false negatives, our recall also would be 0.42

F1 is a score that balances precision against recall F1 = 2 (precision * recall) / precision + recall

In the above example F1 = 0.46

Sometimes we don't want to balance them in this way, in that the consequences of a false positive vs false negative are asymmetric. We might want a higher precision at the expense of recall or vice versa.

```
In [1]: #!/usr/bin/python
        import sys
        import pickle
        import matplotlib.pyplot as plt
        from sklearn import preprocessing
        from sklearn import grid search
        from time import time
        from sklearn import svm, cross_validation
        from sklearn.naive bayes import GaussianNB
        from sklearn.tree import DecisionTreeClassifier
        from sklearn.metrics import accuracy score, precision score, recall sc
        ore, f1 score
        from sklearn.preprocessing import MinMaxScaler
        from sklearn.neighbors import KNeighborsClassifier
        from sklearn.model selection import StratifiedShuffleSplit, GridSearch
        CV, train test split
        from sklearn.pipeline import Pipeline
        sys.path.append("../tools/")
        from feature format import featureFormat, targetFeatureSplit
        from tester import dump classifier and data
        from sklearn.feature selection import SelectKBest
        from sklearn.feature selection import chi2
        from tester import test classifier
        import numpy as np
        import pandas as pd
        from sklearn.ensemble import AdaBoostClassifier
        ### Task 1: Select what features you'll use.
        ### features list is a list of strings, each of which is a feature nam
        ### The first feature must be "poi".
        features_list = ['poi', 'salary', 'total_stock_value', 'exercised_stock_o
        ptions','from poi to this person','from this person to poi','from mess
        ages', 'to messages', 'ratio to poi', 'ratio from poi', 'ratio exercised']
        # You will need to use more features, 8 existing features, 3 new ones
        I created
        ### Load the dictionary containing the dataset
        with open("final project dataset.pkl", "r") as data file:
            data_dict = pickle.load(data_file)
```

import pprint

/Users/robertlee/anaconda/envs/py2/lib/python2.7/site-packages/sklea rn/cross_validation.py:44: DeprecationWarning: This module was depre cated in version 0.18 in favor of the model_selection module into wh ich all the refactored classes and functions are moved. Also note th at the interface of the new CV iterators are different from that of this module. This module will be removed in 0.20.

"This module will be removed in 0.20.", DeprecationWarning)
/Users/robertlee/anaconda/envs/py2/lib/python2.7/site-packages/sklea
rn/grid_search.py:43: DeprecationWarning: This module was deprecated
in version 0.18 in favor of the model_selection module into which al
l the refactored classes and functions are moved. This module will b
e removed in 0.20.

DeprecationWarning)

Task 1: Select what features you'll use.

features_list is a list of strings, each of which is a feature name.

The first feature must be "poi".

I use 8 existing features, and create 3 more (the last 3), combining them to create a list of 11 features. The first feature is the label, poi, so really there are just 10 features in play.

features_old =
['poi','salary','total_stock_value','exercised_stock_options','from_poi_to_this_person','from_this_person_to_pc
features_new = ['ratio_to_poi','ratio_from_poi','ratio_exercised']
features_list = features_old + features_new

```
features_old = ['poi','salary','total_stock_value','exercised stock op
In [2]:
        tions', 'from poi to this person', 'from this person to poi', 'from messa
        qes','to messages']
        features new = ['poi', 'ratio to poi', 'ratio from poi', 'ratio exercised
         ' 1
        features list = features_old + features_new
        features list.pop(8)
        features list
Out[2]: ['poi',
          'salary',
          'total stock value',
          'exercised stock options',
          'from poi to this person',
          'from this person to poi',
          'from messages',
          'to messages',
          'ratio to poi',
          'ratio from poi',
          'ratio exercised']
```

Here is one entry in the dictionary, key: 'ALLEN PHILLIP K'

```
pprint.pprint(data dict['ALLEN PHILLIP K'])
In [3]:
        {'bonus': 4175000,
         'deferral payments': 2869717,
         'deferred_income': -3081055,
         'director fees': 'NaN',
         'email address': 'phillip.allen@enron.com',
         'exercised stock options': 1729541,
         'expenses': 13868,
         'from messages': 2195,
         'from poi to this person': 47,
         'from this person to poi': 65,
         'loan advances': 'NaN',
         'long term incentive': 304805,
         'other': 152,
         'poi': False,
         'restricted stock': 126027,
         'restricted stock deferred': -126027,
         'salary': 201955,
         'shared receipt with poi': 1407,
         'to messages': 2902,
         'total payments': 4484442,
         'total_stock_value': 1729541}
```

Length of dictionary (number of people) and length of one of the entries == # of features

```
In [4]: print len(data_dict)
print len(data_dict['ALLEN PHILLIP K'])

146
21
```

Task 2: Remove outliers

Three of the "people" in the dictionary are not people, but bad data to be removed

"TOTAL" appears to be a summary statistic. "TRAVEL AGENCY IN THE PARK" appears to be a contractor that was paid a sum of money. "LOCKHART EUGENE E" has all NaN values. Harmless, but I go ahead and remove it.

I don't remove data outliers, because they are probably important data (the big shot executives making a lot of money and communicating a lot with each other).

```
In [5]: data_dict.pop('TOTAL', 0)
    data_dict.pop('TRAVEL AGENCY IN THE PAR', 0)
    data_dict.pop('LOCKHART EUGENE E', 0)
    print "Three entries popped"
    len(data_dict)

Three entries popped
Out[5]: 144
```

I look for 'poi = TRUE to count the number of persons of interest

18

Task 3: Create new feature(s)

The 3 new features I create are: ratio_exercised, ratio_to_poi, ratio_from_poi

- (1) ratio_exercised = exercised_stock_options / total_stock_value I created this feature to reflect how much stock the person exercised out of their total (from 0 to 1). I figured that people who were in on the scam might have exercised more of their options, knowing that it was a house of cards. On the other hand, a middle-class person who was not in on the scam might also exercise their options, given the high share price (honest desire to cash in).
- (2) ratio_to_poi = from_this_person_to_poi / to_messages Another 0 to 1 ratio. The absolute number of emails to a POI is less meaningful than the ratio of their messages to POI.
- (3) ratio_to_poi = from_poi_to_this_person / from_messages Another 0 to 1 ratio. The absolute number of emails from a POI is less meaningful than the ratio of messages from a POI.

I am aware that there might be data leakage issues regarding from_this_person_to_poi and from_poi_to_this_person, and thus any features derived from them, but it is not abssolutely clear that this is the case, so I will accept that they are OK features to use for purposes of this assignment.

I first eliminate NaN and bad values (any of numerator < 0, any denominator <= 0)

```
new feature=[]
    for i in data dict:
        if data dict[i][numerator]=="NaN" or data dict[i][denominator]
=="NaN":
            new feature.append("NaN")
        elif data dict[i][numerator]<0 or data dict[i][denominator]<=0</pre>
            new feature.append("NaN")
        else:
            new feature.append(float(data dict[i][numerator]) / data d
ict[i][denominator])
    return new feature
### create lists of new features
ratio exercised=new feature("exercised stock options", "total stock val
ratio_to_poi=new_feature("from_poi_to_this_person","to_messages")
ratio from poi=new feature("from this person to poi", "from messages")
### insert new features into data dict
count=0
for i in data dict:
    data_dict[i]["ratio_exercised"] = ratio_exercised[count]
    data dict[i]["ratio to poi"] = ratio to poi[count]
    data dict[i]["ratio from poi"] = ratio from poi[count]
    count +=1
### store to my dataset for easy export below
my_dataset = data dict
```

Extract features and labels from dataset for local testing

data = featureFormat(my_dataset, features_list, sort_keys = True)

def new feature(numerator, denominator):

In [7]:

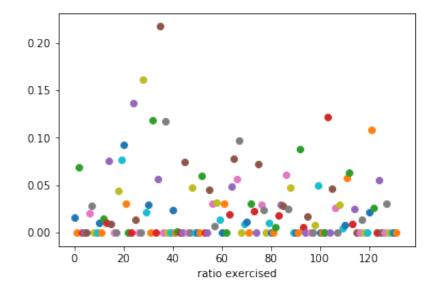
```
labels, features = targetFeatureSplit(data)
  In [8]:
           data = featureFormat(my dataset, features list, sort keys = True)
           labels, features = targetFeatureSplit(data)
```

Plot new features to get a feel for them

ratio_exercised

Ratio of stock options exercised is betwen 0 (none exercised) up to about 20%

```
In [9]: n=0
    for point in data:
        ratio_exercised = point[8]
        plt.scatter( n, ratio_exercised )
        n+=1
    plt.xlabel("ratio exercised")
    plt.show()
```



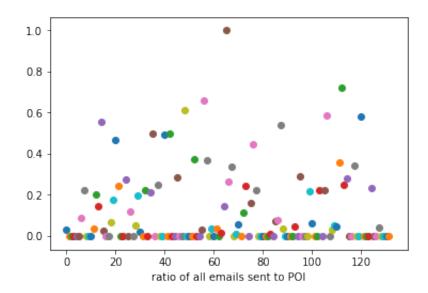
ratio_to_poi

Ratio of all emails sent to POI

Between 0 and 1.0.

The one 1.0 data point is suspiciously high, so I print one that point above the plot. This person was not a PIO, had a salary of 130K, had significant stock, sent 10 emails to POI, received 17 emails from POI, sent 17 total emails, and received 128 total emails. So this data point does not seem outlanding. In addition, since this feature is not in my top 5 best (see below), I am not too concerned with this 1.0 value.

```
n=0
In [10]:
         for point in data:
             ratio to poi = point[9]
             if ratio_to_poi == 1.0:
                 print point
             plt.scatter( n, ratio_to_poi )
         plt.xlabel("ratio of all emails sent to POI")
         plt.show()
            0.00000000e+00
                              1.30724000e+05
                                               2.28276800e+06
                                                                 2.28276800e+06
            1.00000000e+01
                              1.70000000e+01
                                               1.70000000e+01
                                                                 1.28000000e+02
            7.81250000e-02
                              1.00000000e+00
                                               1.00000000e+00]
```



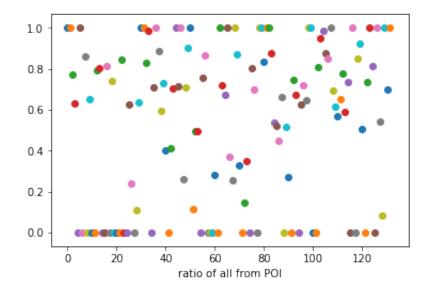
ratio_from_poi

Ratio of all emails from POI

Between 0 and 1.0.

There are a lot of 1.0 data points. So either the data is generally pretty bad, or else these 1.0 values are valid. I'll go with the latter assumption, though of course I could investigate more.

```
In [11]: n=0
    for point in data:
        ratio_from_poi = point[10]
        plt.scatter( n, ratio_from_poi )
        n+=1
    plt.xlabel("ratio of all from POI")
    plt.show()
```



I print a correlation matrix to get a sense of feature correlation with the labels (row/column 0)

Of the 3 new features, 8,9,10, only 9, ratio_from_poi has a fairly strong correlation, 0.31. Thus one could argue that is the only new feature worth looking at. But for sake of academic exercise, rather than trying to engineer additional new features, I'll continue with these.

Of the original 7 features, 1,2,3 all have correlations to the label above 0.3. These are the 3 best features, and 9, above, is the forth best. None of the other features are even close in terms of correlation. Thus one could argue that we should only use these 3 or 4 features.

Not surprisingly, the first three features, all related to money, are highly correlated to each other with 2 and 3 being correlated at 0.96. Thus 2 and 3 in particular would be amendable to PCA to reduce the number of features. For the scope of this project, though, PCA didn't seem necessary, as we are already working with a small number of features, so I did not do PCA.

1 salary 2 total_stock_value 3 exercised_stock_options 9 ratio_from_poi

```
In [12]: n=0
    for feature in features_list:
        print n, feature
        n+=1

    df = pd.DataFrame(data)
    df.corr()
    #for each in np.corrcoef(data):
        # print each
```

0 poi
1 salary
2 total_stock_value
3 exercised_stock_options
4 from_poi_to_this_person
5 from_this_person_to_poi
6 from_messages
7 to_messages
8 ratio_to_poi
9 ratio_from_poi

10 ratio_exercised

Out[12]:

	0	1	2	3	4	5	6	7
0	1.000000	0.323062	0.373375	0.378329	0.175804	0.120282	-0.043027	0.09152
1	0.323062	1.000000	0.507512	0.434179	0.382845	0.188981	0.128107	0.36915
2	0.373375	0.507512	1.000000	0.963300	0.118218	0.008916	-0.022347	0.09709
3	0.378329	0.434179	0.963300	1.000000	0.107552	-0.009687	-0.033640	0.06546
4	0.175804	0.382845	0.118218	0.107552	1.000000	0.491520	0.245550	0.60815
5	0.120282	0.188981	0.008916	-0.009687	0.491520	1.000000	0.606896	0.60073
6	-0.043027	0.128107	-0.022347	-0.033640	0.245550	0.606896	1.000000	0.50404
7	0.091524	0.369150	0.097094	0.065464	0.608153	0.600738	0.504045	1.00000
8	0.130400	0.214900	-0.008048	-0.000078	0.571588	0.070609	-0.018426	0.10368
9	0.310498	0.271439	0.101619	0.088343	0.237443	0.069135	-0.072205	0.09692
10	-0.063065	-0.020504	0.212430	0.290629	0.006517	-0.007351	0.047233	-0.01396

Trying out a few classifiers

Rather than use systematic cross-validation and classifier search with gridSearchCSV, for now I simply try some classifiers on three sizes, just to get a feel: 0.1, 0.3, 0.5

features_train, features_test, labels_train, labels_test = cross_validation.train_test_split(features, labels, test_size=X, random_state=42)

Since there are only 18 POIs, we can only go so small with the test size before we don't get any POIs. 0.1 is probably about the smallest we can use.

Trying NB with the above features and test sizes

```
def classify NB(features train, labels train):
In [14]:
             ### your code goes here--should return a trained decision tree cla
         ssifer
             clf = GaussianNB()
             clf.fit(features train, labels train)
             return clf
         def class run(features train, labels train, features test, labels test
         ):
             clf = classify NB(features train, labels train)
             pred = clf.predict(features test)
             return {"Accuracy":accuracy_score(pred,labels_test),\
                                    "Precision":precision score(pred, labels test
         ),\
                                    "Recall":recall score(pred, labels test), \
                                    "F1":f1 score(pred, labels test)}
```

```
In [15]: | print "all 10 features (plus poi label)"
         def print it(title,input):
            print title
            answer = input
            for each in answer:
                print each," : ",answer[each]
            print
         print it("NB: 0.1 test size", class run(features train1, labels train1,
         features test1, labels test1))
         print it("NB: 0.3 test size", class run(features train3, labels train3,
         features_test3, labels_test3))
         print it("NB: 0.5 test size",class run(features train5, labels train5,
         features test5, labels test5))
         all 10 features (plus poi label)
        NB: 0.1 test size
        Recall : 0.5
         F1 : 0.4
        Accuracy : 0.785714285714
        NB: 0.3 test size
```

Recall : 0.75

Precision : 0.5 Accuracy : 0.9

NB: 0.5 test size

Precision : 0.5

Recall : 0.3333333333333

Accuracy : 0.863636363636

F1 : 0.6

F1 : 0.4

Trying NB with different subsets of the features

```
features list contains all 11 features (8 original, 3 new)
I create these subset feature lists as well
Each must contain the first feature, poi
features old = poi + the original 7
features new = poi + the 3 new ones
features first = poi + 1,2,3 ($ related)
features\_second = poi + 4,5,6,7  (email related)
features_F = features_first + 'ratio_exercised' ($ related)
features_E = features_second + ratio_to_poi' + 'ratio_from_poi' (email related)
 In [16]: | features_first = ['poi', 'salary', 'total_stock_value', 'exercised stock
            options']
            features second = ['poi', 'from poi to this person', 'from this person t
            o poi', 'from messages', 'to messages', 'ratio to poi', 'ratio from poi', '
            ratio exercised']
            features F = ['poi', 'salary', 'total stock value', 'exercised stock opti
            ons', 'ratio exercised']
            features_E = ['poi','from_poi_to_this_person','from_this_person_to_poi
             ','from messages','to messages','ratio to poi','ratio from poi']
```

Using the original 7 features yields the same results as using all 10 features

```
In [17]: data = featureFormat(my dataset, features old, sort keys = True)
         labels, features o = targetFeatureSplit(data)
         features train1o, features test1o, labels train1o, labels test1o = cro
         ss validation.train test split(features o, labels, test size=0.1, rand
         om state=42)
         features train3o, features test3o, labels train3o, labels test3o = cro
         ss validation.train test split(features o, labels, test size=0.3, rand
         features train5o, features test5o, labels train5o, labels test5o = cro
         ss validation.train test split(features o, labels, test size=0.5, rand
         om state=42)
         print "features old = the original 7"
         print it("NB: 0.1 test size", class run(features train1o, labels train1
         o, features testlo, labels testlo))
         print it("NB: 0.3 test size", class run(features train3o, labels train3
         o, features_test3o, labels_test3o))
         print it("NB: 0.5 test size", class run(features train50, labels train5
         o, features test5o, labels test5o))
         features old = the original 7
         NB: 0.1 test size
         Recall: 0.5
         F1 : 0.4
         Precision : 0.3333333333333
         Accuracy : 0.785714285714
```

Using just the 3 new features doesn't work as well

NB: 0.3 test size Recall : 0.75

Precision : 0.5 Accuracy : 0.9

NB: 0.5 test size

Precision: 0.5

Recall : 0.3333333333333

Accuracy : 0.863636363636

F1 : 0.6

F1 : 0.4

data n = featureFormat(my dataset, features new, sort keys = True) In [18]: labels, features n = targetFeatureSplit(data n) features train1n, features test1n, labels train1n, labels test1n = cro ss validation.train test split(features n, labels, test size=0.1, rand om state=42) features train3n, features test3n, labels train3n, labels test3n = cro ss validation.train test split(features n, labels, test size=0.3, rand om state=42) features train5n, features test5n, labels train5n, labels test5n = cro ss_validation.train_test_split(features_n, labels, test_size=0.5, rand om state=42) print "features new = poi + 3 new features" print print it("NB: 0.1 test size",class run(features train1n, labels train1 n, features test1n, labels test1n)) print_it("NB: 0.3 test size",class run(features train3n, labels train3 n, features test3n, labels test3n)) print_it("NB: 0.5 test size",class_run(features_train5n, labels_train5 n, features test5n, labels test5n)) features new = poi + 3 new features /Users/robertlee/anaconda/envs/py2/lib/python2.7/site-packages/sklea rn/metrics/classification.py:1115: UndefinedMetricWarning: Recall is ill-defined and being set to 0.0 due to no true samples. 'recall', 'true', average, warn for) /Users/robertlee/anaconda/envs/py2/lib/python2.7/site-packages/sklea

/Users/robertlee/anaconda/envs/py2/lib/python2.7/site-packages/sklea rn/metrics/classification.py:1115: UndefinedMetricWarning: F-score is ill-defined and being set to 0.0 due to no true samples.

'recall', 'true', average, warn for)

NB: 0.1 test size
Recall : 0.0
F1 : 0.0
Precision : 0.0

Accuracy : 0.833333333333

NB: 0.3 test size

Recall : 0.3333333333333

F1 : 0.25

Precision : 0.2

Accuracy : 0.833333333333

NB: 0.5 test size Recall : 0.0 F1 : 0.0

Precision : 0.0

Neither does using just money related features

```
data f = featureFormat(my dataset, features first, sort keys = True)
In [19]:
         labels, features f = targetFeatureSplit(data f)
         features train1f, features test1f, labels train1f, labels test1f = cro
         ss validation.train test split(features f, labels, test size=0.1, rand
         om state=42)
         features train3f, features test3f, labels train3f, labels test3f = cro
         ss validation.train test split(features f, labels, test size=0.3, rand
         om state=42)
         features train5f, features test5f, labels train5f, labels test5f = cro
         ss validation.train test split(features f, labels, test size=0.5, rand
         om state=42)
         print "features first = poi + 1,2,3 ($ related)"
         print_it("NB: 0.1 test size",class_run(features_train1f, labels_train1
         f, features test1f, labels test1f))
         print it("NB: 0.3 test size", class run(features train3f, labels train3
         f, features test3f, labels test3f))
         print it("NB: 0.5 test size", class run(features train5f, labels train5
         f, features test5f, labels test5f))
         features first = poi + 1,2,3 ($ related)
         NB: 0.1 test size
         Recall : 0.0
         F1 : 0.0
         Precision : 0.0
         Accuracy : 0.846153846154
         NB: 0.3 test size
         Recall : 0.0
         F1 : 0.0
```

Accuracy: 0.871794871795

NB: 0.5 test size

Recall: 0.25

F1: 0.1818181818

Precision: 0.142857142857

Accuracy: 0.861538461538

Precision : 0.0

Nor just email

```
data s = featureFormat(my dataset, features second, sort keys = True)
In [20]:
         labels, features s = targetFeatureSplit(data s)
         features train1s, features test1s, labels train1s, labels test1s = cro
         ss validation.train test split(features s, labels, test size=0.1, rand
         om state=42)
         features train3s, features test3s, labels train3s, labels test3s = cro
         ss validation.train test split(features s, labels, test size=0.3, rand
         om state=42)
         features train5s, features test5s, labels train5s, labels test5s = cro
         ss validation.train test split(features_s, labels, test_size=0.5, rand
         om state=42)
         print "poi + 4,5,6,7 (email related)"
         print
         print it("NB: 0.1 test size", class run(features train1s, labels train1
         s, features test1s, labels test1s))
         print it("NB: 0.3 test size", class run(features train3s, labels train3
         s, features test3s, labels test3s))
         print_it("NB: 0.5 test size",class_run(features_train5s, labels_train5
         s, features test5s, labels test5s))
```

poi + 4,5,6,7 (email related)

Accuracy : 0.8333333333333

Now I'll run SelectKBest on k=1 through k=9

```
In [21]: acc = []
         precision = []
         recall = []
         f1 = []
         for n in range (9,0,-1):
             print "SelectKBest() best features, k = ", n
             data = featureFormat(my dataset, features list, sort keys = True)
             labels, features n = targetFeatureSplit(data)
             K best n = SelectKBest(k=n)
             features Kbest n = K best n.fit transform(features n, labels)
             features train1, features test1, labels train1, labels test1 = cro
         ss validation.train test split(features Kbest n, labels, test size=0.1
         , random state=42)
             features train3, features test3, labels train3, labels test3 = cro
         ss validation.train test split(features Kbest n, labels, test size=0.3
         , random state=42)
             features_train5, features_test5, labels_train5, labels_test5 = cro
         ss validation.train test split(features Kbest n, labels, test size=0.5
         , random state=42)
             print
             ts1 = class run(features train1, labels train1, features test1, la
         bels test1)
             ts3 = class run(features train3, labels train3, features test3, la
         bels test3)
             ts5 = class run(features train5, labels train5, features test5, la
         bels test5)
             print it("NB: 0.1 test size",ts1)
             print it("NB: 0.3 test size",ts3)
             print_it("NB: 0.5 test size",ts5)
             acc.append(ts5["Accuracy"])
             precision.append(ts5["Recall"])
             recall.append(ts5["Precision"])
             f1.append(ts5["F1"])
         SelectKBest() best features, k = 9
         NB: 0.1 test size
         Recall : 0.5
         F1 : 0.4
```

NB: 0.5 test size

Recall : 0.5

F1 : 0.5

Precision : 0.5

Accuracy : 0.909090909091

SelectKBest() best features, k = 8

NB: 0.1 test size

Recall : 0.5

F1 : 0.4

NB: 0.3 test size

Recall : 0.75

F1 : 0.6

Precision : 0.5 Accuracy : 0.9

NB: 0.5 test size

Recall : 0.5

F1 : 0.5

Precision : 0.5

Accuracy : 0.909090909091

SelectKBest() best features, k = 7

NB: 0.1 test size

Recall : 0.5

F1 : 0.4

NB: 0.3 test size

Recall : 0.75

F1 : 0.6

Precision : 0.5 Accuracy : 0.9

NB: 0.5 test size

Recall : 0.428571428571

F1 : 0.461538461538

Precision : 0.5

Accuracy : 0.893939393939

SelectKBest() best features, k = 6

NB: 0.1 test size

Recall : 0.5

F1 : 0.4

NB: 0.3 test size

Recall : 0.75

F1 : 0.6

Precision : 0.5 Accuracy : 0.9

NB: 0.5 test size

Recall : 0.5

F1 : 0.5

Precision : 0.5

Accuracy : 0.909090909091

SelectKBest() best features, k = 5

NB: 0.1 test size

Recall : 0.5

F1 : 0.4

Accuracy : 0.785714285714

NB: 0.3 test size

Recall : 0.75

F1 : 0.6

Precision : 0.5

Accuracy : 0.9

NB: 0.5 test size

Recall: 0.5

F1 : 0.5

Precision : 0.5

Accuracy : 0.909090909091

SelectKBest() best features, k = 4

NB: 0.1 test size

Recall: 0.5

F1 : 0.4

Accuracy : 0.785714285714

NB: 0.3 test size

Recall : 0.75

F1 : 0.6

Precision : 0.5

Accuracy : 0.9

NB: 0.5 test size

Recall: 0.5

F1 : 0.5

Precision : 0.5

```
SelectKBest() best features, k = 3
```

NB: 0.1 test size Recall : 0.5

F1 : 0.4

NB: 0.3 test size Recall : 0.75

F1 : 0.6

Precision : 0.5 Accuracy : 0.9

NB: 0.5 test size Recall : 0.5 F1 : 0.5

Precision : 0.5

Accuracy : 0.909090909091

SelectKBest() best features, k = 2

NB: 0.1 test size Recall : 0.5

F1 : 0.4

NB: 0.3 test size Recall : 0.75

F1 : 0.6

Precision : 0.5 Accuracy : 0.9

NB: 0.5 test size

Recall : 0.428571428571

F1 : 0.461538461538

Precision : 0.5

Accuracy : 0.893939393939

SelectKBest() best features, k = 1

NB: 0.1 test size Recall : 1.0

F1 : 0.5

NB: 0.3 test size Recall : 1.0

F1 : 0.66666666667

Precision : 0.5 Accuracy : 0.925

```
NB: 0.5 test size
Recall : 0.5
F1 : 0.5
Precision : 0.5
Accuracy : 0.909090909091
```

Above feature investigations

I stored the scores for test_size=0.5 and print them below. We get pretty much the same results whether we use 9 features or 1. The best feature is the original feature of salary, perhaps this is enough. It sounds intuitively reasonable that salary predicts POI.

```
In [22]: print "Accuracies"
    print acc
    print "Precisions"
    print precision
    print "Recalls"
    print recall
    print f1

Accuracies
    [0.909090909090906, 0.9090909090906, 0.89393939393939392, 0.90
    909090909090906, 0.9090909090906, 0.90909090909090909
```

Task 4: Try a varity of classifiers

Please name your classifier clf for easy export below.

Note that if you want to do PCA or other multi-stage operations,

you'll need to use Pipelines. For more info:

http://scikit-learn.org/stable/modules/pipeline.html (http://scikit-learn.org/stable/modules/pipeline.html)

Trying KNN() with 3 neighbors

I stored the scores for test_size=0.5 and print them below. Again, we get similar results whether we use 10 features all the way to just 1, salary.

```
In [23]: def classify KNN(features train, labels train):
             ### your code goes here--should return a trained decision tree cla
         ssifer
             clf = KNeighborsClassifier(3)
             clf.fit(features train, labels train)
             return clf
         def class run KNN(features train, labels train, features test, labels
         test):
             clf = classify KNN(features train, labels train)
             pred = clf.predict(features test)
             return {"Accuracy":accuracy score(pred, labels test),\
                                  "Precision":precision score(pred, labels test
         ),\
                                   "Recall":recall score(pred, labels_test), \
                                  "F1":f1 score(pred, labels test)}
         print "All 10 features, KNN, k=3"
         print
         print it("KNN: 0.1 test size", class run KNN(features train1, labels tr
         ain1, features test1, labels test1))
         print_it("KNN: 0.3 test size",class run KNN(features train3, labels tr
         ain3, features_test3, labels_test3))
         print it("KNN: 0.5 test size", class run KNN(features train5, labels tr
         ain5, features_test5, labels_test5))
         All 10 features, KNN, k=3
         KNN: 0.1 test size
         Recall : 1.0
         F1 : 0.5
         Accuracy : 0.857142857143
         KNN: 0.3 test size
         Recall: 0.0
         F1 : 0.0
         Precision: 0.0
         Accuracy : 0.825
         KNN: 0.5 test size
         Recall : 0.375
```

F1 : 0.428571428571 Precision : 0.5

```
In [24]: | acc = []
         precision = []
         recall = []
         f1 = []
         for n in range (9,0,-1):
             print "SelectKBest() best features, k = ", n
             data = featureFormat(my dataset, features list, sort keys = True)
             labels, features n = targetFeatureSplit(data)
             K best n = SelectKBest(k=n)
             features Kbest n = K best n.fit transform(features n, labels)
             features train1, features test1, labels train1, labels test1 = cro
         ss validation.train test split(features Kbest n, labels, test size=0.1
         , random state=42)
             features train3, features test3, labels train3, labels test3 = cro
         ss validation.train test split(features Kbest n, labels, test size=0.3
         , random state=42)
             features_train5, features_test5, labels_train5, labels_test5 = cro
         ss validation.train test split(features Kbest n, labels, test size=0.5
         , random state=42)
             print
             ts1 = class run KNN(features train1, labels train1, features test1
         , labels test1)
             ts3 = class run KNN(features train3, labels train3, features test3
         , labels test3)
             ts5 = class run KNN(features train5, labels train5, features test5
         , labels test5)
             print it("KNN: 0.1 test size",ts5)
             print it("KNN: 0.3 test size",ts5)
             print it("KNN: 0.5 test size",ts5)
             acc.append(ts5["Accuracy"])
             precision.append(ts5["Recall"])
             recall.append(ts5["Precision"])
             f1.append(ts5["F1"])
         SelectKBest() best features, k = 9
         KNN: 0.1 test size
```

Recall: 0.6
F1: 0.545454545455
Precision: 0.5
Accuracy: 0.924242424242

KNN: 0.3 test size
Recall: 0.6
F1: 0.545454545455
Precision: 0.5
Accuracy: 0.924242424242

KNN: 0.5 test size

Recall : 0.6

F1 : 0.5454545455

Precision : 0.5

Accuracy : 0.924242424242

SelectKBest() best features, k = 8

KNN: 0.1 test size

Recall : 0.6

F1 : 0.545454545455

Precision : 0.5

Accuracy : 0.924242424242

KNN: 0.3 test size

Recall : 0.6

F1 : 0.5454545455

Precision : 0.5

Accuracy : 0.924242424242

KNN: 0.5 test size

Recall : 0.6

F1 : 0.545454545455

Precision : 0.5

Accuracy : 0.924242424242

SelectKBest() best features, k = 7

KNN: 0.1 test size

Recall : 0.6

F1 : 0.5454545455

Precision : 0.5

Accuracy : 0.924242424242

KNN: 0.3 test size

Recall : 0.6

F1 : 0.5454545455

Precision : 0.5

Accuracy : 0.924242424242

KNN: 0.5 test size

Recall : 0.6

F1 : 0.5454545455

Precision : 0.5

Accuracy : 0.924242424242

SelectKBest() best features, k = 6

KNN: 0.1 test size

Recall : 0.6

F1 : 0.5454545455

Precision : 0.5

KNN: 0.3 test size

Recall : 0.6

F1 : 0.5454545455

Precision : 0.5

Accuracy : 0.924242424242

KNN: 0.5 test size

Recall : 0.6

F1 : 0.545454545455

Precision : 0.5

Accuracy : 0.924242424242

SelectKBest() best features, k = 5

KNN: 0.1 test size

Recall : 0.6

F1 : 0.545454545455

Precision : 0.5

Accuracy : 0.924242424242

KNN: 0.3 test size

Recall : 0.6

F1 : 0.5454545455

Precision : 0.5

Accuracy : 0.924242424242

KNN: 0.5 test size

Recall: 0.6

F1 : 0.5454545455

Precision : 0.5

Accuracy : 0.924242424242

SelectKBest() best features, k = 4

KNN: 0.1 test size

Recall: 0.6

F1 : 0.545454545455

Precision : 0.5

Accuracy : 0.924242424242

KNN: 0.3 test size

Recall : 0.6

F1 : 0.545454545455

Precision : 0.5

Accuracy : 0.924242424242

KNN: 0.5 test size

Recall: 0.6

F1 : 0.545454545455

Precision : 0.5

SelectKBest() best features, k = 3

KNN: 0.1 test size

Recall : 0.6

F1 : 0.5454545455

Precision : 0.5

Accuracy : 0.924242424242

KNN: 0.3 test size

Recall : 0.6

F1 : 0.545454545455

Precision : 0.5

Accuracy : 0.924242424242

KNN: 0.5 test size

Recall: 0.6

F1 : 0.5454545455

Precision : 0.5

Accuracy : 0.924242424242

SelectKBest() best features, k = 2

KNN: 0.1 test size

Recall : 0.6

F1 : 0.545454545455

Precision : 0.5

Accuracy : 0.924242424242

KNN: 0.3 test size

Recall : 0.6

F1 : 0.5454545455

Precision : 0.5

Accuracy : 0.924242424242

KNN: 0.5 test size

Recall : 0.6

F1 : 0.5454545455

Precision : 0.5

Accuracy : 0.924242424242

SelectKBest() best features, k = 1

KNN: 0.1 test size

Recall : 0.375

F1 : 0.428571428571

Precision : 0.5

Accuracy : 0.8787878788

KNN: 0.3 test size

Recall : 0.375

F1 : 0.428571428571

Precision : 0.5

KNN: 0.5 test size
Recall : 0.375

F1 : 0.428571428571

Precision : 0.5
Accuracy : 0.8787878788

In [25]: print "Accuracies"

print acc

print "Precisions"
print precision
print "Recalls"
print recall

print "F1s"

print f1

Accuracies

Precisions

Recalls

[0.5, 0.5, 0.5, 0.5, 0.5, 0.5, 0.5, 0.5]

Fls

[0.54545454545454541, 0.5454545454545454541, 0.54545454545454545454545 545454545454541, 0.5454545454545454541, 0.54545454545454545 4545454541, 0.5454545454545454545454545454545454571428571 Task 5: Tune your classifier to achieve better than .3 precision and recall using our testing script. Check the tester.py script in the final project folder for details on the evaluation method, especially the test_classifier function. Because of the small size of the dataset, the script uses stratified shuffle split cross validation. For more info:

http://scikit-

<u>learn.org/stable/modules/generated/sklearn.cross_validation.StratifiedShuffleSp</u> (http://scikit-

learn.org/stable/modules/generated/sklearn.cross_validation.StratifiedShuffleSp

Example starting point. Try investigating other evaluation techniques!

10 gridSearchCSVs on KNN, evaluating by F1 score

Of these 10 searches, I got the following two classifiers, 5 times each Classifier 1

KNeighborsClassifier(algorithm='auto', leaf_size=3, metric='minkowski', metric_params=None, n_jobs=1, n_neighbors=1, p=1, weights='uniform')

F1 = 0.45

Classifier 2 KNeighborsClassifier(algorithm='auto', leaf_size=3, metric='minkowski', metric_params=None, n_jobs=1, n_neighbors=1, p=1, weights='uniform')
F1 = 0.37

All 10 features, test size 0.3

```
In [26]: data = featureFormat(my dataset, features list, sort keys = True)
         labels, features = targetFeatureSplit(data)
         knn = KNeighborsClassifier()
         scaler = MinMaxScaler()
         sss = StratifiedShuffleSplit(n splits = 100, test size=0.3, random stat
         e = 42)
         pipe = Pipeline(steps=[('scaling', scaler), ("knn", knn)])
         param grid = {"knn n neighbors": [1, 3, 5, 8, 10, 12, 14],
                       "knn__algorithm": ["auto", "ball_tree", "kd_tree", "brute
         "],
                       "knn leaf size": range(3,10,1),
                       "knn p": [1,2]
                      }
         knnclf = GridSearchCV(pipe, param grid, scoring='f1', cv=sss)
         knnclf.fit(features, labels)
         bestknn = knnclf.best estimator
         bestparams = knnclf.best params
         bestscore = knnclf.best score
         bestscorer = knnclf.scorer
         bestcv = knnclf.cv results
         print "Best KNN"
         print bestknn
         print bestparams
         print bestscore
         print bestscorer
         /Users/robertlee/anaconda/envs/py2/lib/python2.7/site-packages/sklea
         rn/metrics/classification.py:1113: UndefinedMetricWarning: F-score i
         s ill-defined and being set to 0.0 due to no predicted samples.
           'precision', 'predicted', average, warn for)
         Best KNN
         Pipeline(steps=[('scaling', MinMaxScaler(copy=True, feature range=(0
         , 1))), ('knn', KNeighborsClassifier(algorithm='auto', leaf size=3,
         metric='minkowski',
                    metric params=None, n jobs=1, n neighbors=1, p=1,
                    weights='uniform'))])
```

{'knn leaf size': 3, 'knn algorithm': 'auto', 'knn n neighbors':

1, 'knn_p': 1} 0.371892385392

make scorer(f1 score)

test_classifier issue

I then tried using test_classifier method from tester.py. I reduced the number of folds from 1000 to 10, but it still was running after 5 minutes, so I think something is wrong.

print test_classifier(knnclf, my_dataset, features_list)

So instead below I test the classifier three times with different test sizes. The results look adequate, but I guess I need you to run the tester on my final classifier.

```
In [56]: best = KNeighborsClassifier(algorithm='auto', leaf size=3, metric='min
         kowski',
                    metric params=None, n jobs=1, n neighbors=1, p=1,
                    weights='uniform')
         data = featureFormat(my dataset, features list, sort keys = True)
         labels, features = targetFeatureSplit(data)
         features train1, features test1, labels train1, labels test1 = cross v
         alidation.train test split(features, labels, test size=0.1, random sta
         te=42)
         features_train3, features_test3, labels_train3, labels_test3 = cross_v
         alidation.train test split(features, labels, test size=0.3, random sta
         te=42)
         features train5, features test5, labels train5, labels test5 = cross v
         alidation.train test split(features, labels, test size=0.5, random sta
         te=42)
         print "Testing Above Classifier"
         print
         best.fit(features train1, labels train1)
         pred = best.predict(features test1)
         print "Testing on 0.1 of sample"
         print "accuracy",accuracy score(pred,labels test1)
         print "precision",precision_score(pred,labels_test1)
         print "recall", recall score(pred, labels test1)
         print "F1",f1_score(pred,labels_test1)
         print
         best.fit(features train3, labels train3)
         pred = best.predict(features test3)
         print "Testing on 0.3 of sample"
         print "accuracy",accuracy score(pred,labels test3)
         print "precision", precision score(pred, labels test3)
         print "recall", recall score(pred, labels test3)
         print "F1",f1 score(pred,labels test3)
         print
         best.fit(features train5, labels train5)
         pred = best.predict(features test5)
         print "Testing on 0.5 of sample"
         print "accuracy",accuracy score(pred,labels test5)
         print "precision",precision score(pred,labels test5)
         print "recall", recall_score(pred, labels_test5)
         print "F1",f1 score(pred,labels test5)
```

Testing Above Classifier

Testing on 0.3 of sample accuracy 0.825 precision 0.5 recall 0.428571428571 F1 0.461538461538

Testing on 0.5 of sample accuracy 0.818181818182 precision 0.5 recall 0.25 F1 0.3333333333333

All 10 features, test size 0.1

```
knn = KNeighborsClassifier()
In [27]:
         scaler = MinMaxScaler()
         sss = StratifiedShuffleSplit(n splits = 100, test size=0.1, random stat
         pipe = Pipeline(steps=[('scaling', scaler), ("knn", knn)])
         param grid = {"knn n neighbors": [1, 3, 5, 8, 10, 12, 14],
                       "knn algorithm": ["auto", "ball tree", "kd tree", "brute
         "],
                       "knn leaf_size": range(3,10,1),
                       "knn__p": [1,2]
                      }
         knnclf = GridSearchCV(pipe, param grid, scoring='f1', cv=sss)
         knnclf.fit(features, labels)
         bestknn = knnclf.best estimator
         bestparams = knnclf.best params
         bestscore = knnclf.best_score_
         bestscorer = knnclf.scorer
         bestcv = knnclf.cv results
         print "Best KNN"
         print bestknn
         print bestparams
         print bestscore
         print bestscorer
```

```
In [57]: best = KNeighborsClassifier(algorithm='auto', leaf size=3, metric='min
         kowski',
                    metric params=None, n jobs=1, n neighbors=1, p=2,
                    weights='uniform')
         data = featureFormat(my dataset, features list, sort keys = True)
         labels, features = targetFeatureSplit(data)
         features train1, features test1, labels train1, labels test1 = cross v
         alidation.train test split(features, labels, test size=0.1, random sta
         te=42)
         features_train3, features_test3, labels_train3, labels_test3 = cross_v
         alidation.train test split(features, labels, test size=0.3, random sta
         te=42)
         features train5, features test5, labels train5, labels test5 = cross v
         alidation.train test split(features, labels, test size=0.5, random sta
         te=42)
         print "Testing Above Classifier"
         print
         best.fit(features train1, labels train1)
         pred = best.predict(features test1)
         print "Testing on 0.1 of sample"
         print "accuracy",accuracy score(pred,labels test1)
         print "precision",precision_score(pred,labels_test1)
         print "recall", recall score(pred, labels test1)
         print "F1",f1_score(pred,labels_test1)
         print
         best.fit(features train3, labels train3)
         pred = best.predict(features test3)
         print "Testing on 0.3 of sample"
         print "accuracy",accuracy score(pred,labels test3)
         print "precision", precision score(pred, labels test3)
         print "recall", recall score(pred, labels test3)
         print "F1",f1 score(pred,labels test3)
         print
         best.fit(features train5, labels train5)
         pred = best.predict(features test5)
         print "Testing on 0.5 of sample"
         print "accuracy",accuracy score(pred,labels test5)
         print "precision", precision score(pred, labels test5)
         print "recall", recall_score(pred, labels_test5)
         print "F1",f1 score(pred,labels test5)
```

Testing Above Classifier

Testing on 0.3 of sample accuracy 0.825 precision 0.5 recall 0.428571428571 F1 0.461538461538

Testing on 0.5 of sample accuracy 0.818181818182 precision 0.5 recall 0.25 F1 0.3333333333333

5 best features, test size 0.3

```
data = featureFormat(my dataset, features list, sort keys = True)
In [28]:
         labels, features = targetFeatureSplit(data)
         K best = SelectKBest(k=5)
         features kbest = K best.fit transform(features, labels)
         knn = KNeighborsClassifier()
         scaler = MinMaxScaler()
         sss = StratifiedShuffleSplit(n splits = 100, test size=0.3, random stat
         e = 42)
         pipe = Pipeline(steps=[('scaling', scaler), ("knn", knn)])
         param grid = {"knn n neighbors": [1, 3, 5, 8, 10, 12, 14],
                       "knn algorithm": ["auto", "ball tree", "kd tree", "brute
         "],
                       "knn leaf size": range(3,10,1),
                       "knn p": [1,2]
         knnclf = GridSearchCV(pipe, param grid, scoring='f1', cv=sss)
         knnclf.fit(features, labels)
         bestknn = knnclf.best estimator
         bestparams = knnclf.best params
         bestscore = knnclf.best score
         bestscorer = knnclf.scorer
         bestcv = knnclf.cv results
         print "Best KNN"
         print bestknn
         print bestparams
         print bestscore
         print bestscorer
         Best KNN
         Pipeline(steps=[('scaling', MinMaxScaler(copy=True, feature range=(0
         , 1))), ('knn', KNeighborsClassifier(algorithm='auto', leaf size=3,
         metric='minkowski',
                    metric params=None, n jobs=1, n neighbors=1, p=1,
                    weights='uniform'))))
         {'knn__leaf_size': 3, 'knn__algorithm': 'auto', 'knn__n_neighbors':
         1, 'knn p': 1}
         0.371892385392
```

```
In [29]: data = featureFormat(my dataset, features list, sort keys = True)
         labels, features = targetFeatureSplit(data)
         K best = SelectKBest(k=5)
         features kbest = K best.fit transform(features, labels)
         knn = KNeighborsClassifier()
         scaler = MinMaxScaler()
         sss = StratifiedShuffleSplit(n splits = 100, test size=0.1, random stat
         e = 42)
         pipe = Pipeline(steps=[('scaling', scaler), ("knn", knn)])
         param grid = {"knn n neighbors": [1, 3, 5, 8, 10, 12, 14],
                       "knn algorithm": ["auto", "ball tree", "kd tree", "brute
         "],
                       "knn leaf size": range(3,10,1),
                       "knn p": [1,2]
         knnclf = GridSearchCV(pipe, param grid, scoring='f1', cv=sss)
         knnclf.fit(features, labels)
         bestknn = knnclf.best estimator
         bestparams = knnclf.best params
         bestscore = knnclf.best score
         bestscorer = knnclf.scorer
         bestcv = knnclf.cv results
         print "Best KNN"
         print bestknn
         print bestparams
         print bestscore
         print bestscorer
         Best KNN
         Pipeline(steps=[('scaling', MinMaxScaler(copy=True, feature range=(0
         , 1))), ('knn', KNeighborsClassifier(algorithm='auto', leaf size=3,
         metric='minkowski',
                    metric params=None, n jobs=1, n neighbors=1, p=2,
                    weights='uniform'))))
         {'knn__leaf_size': 3, 'knn__algorithm': 'auto', 'knn__n_neighbors':
         1, 'knn p': 2}
         0.446904761905
```

```
In [53]: data = featureFormat(my dataset, features list, sort keys = True)
         labels, features = targetFeatureSplit(data)
         K best = SelectKBest(k=3)
         features kbest = K best.fit transform(features, labels)
         knn = KNeighborsClassifier()
         scaler = MinMaxScaler()
         sss = StratifiedShuffleSplit(n splits = 100, test size=0.3,random stat
         e = 42)
         pipe = Pipeline(steps=[('scaling', scaler), ("knn", knn)])
         param grid = {"knn n neighbors": [1, 3, 5, 8, 10, 12, 14],
                       "knn algorithm": ["auto", "ball tree", "kd tree", "brute
         "],
                       "knn leaf size": range(3,10,1),
                       "knn p": [1,2]
         knnclf = GridSearchCV(pipe, param grid, scoring='f1', cv=sss)
         knnclf.fit(features, labels)
         bestknn = knnclf.best estimator
         bestparams = knnclf.best params
         bestscore = knnclf.best score
         bestscorer = knnclf.scorer
         bestcv = knnclf.cv results
         print "Best KNN"
         print bestknn
         print bestparams
         print bestscore
         print bestscorer
         Best KNN
         Pipeline(steps=[('scaling', MinMaxScaler(copy=True, feature range=(0
         , 1))), ('knn', KNeighborsClassifier(algorithm='auto', leaf size=3,
         metric='minkowski',
                    metric params=None, n jobs=1, n neighbors=1, p=1,
                    weights='uniform'))))
         {'knn__leaf_size': 3, 'knn__algorithm': 'auto', 'knn__n_neighbors':
         1, 'knn p': 1}
         0.371892385392
```

```
In [52]: data = featureFormat(my dataset, features list, sort keys = True)
         labels, features = targetFeatureSplit(data)
         K best = SelectKBest(k=3)
         features kbest = K best.fit transform(features, labels)
         knn = KNeighborsClassifier()
         scaler = MinMaxScaler()
         sss = StratifiedShuffleSplit(n splits = 100, test size=0.1, random stat
         e = 42)
         pipe = Pipeline(steps=[('scaling', scaler), ("knn", knn)])
         param grid = {"knn n neighbors": [1, 3, 5, 8, 10, 12, 14],
                       "knn algorithm": ["auto", "ball tree", "kd tree", "brute
         "],
                       "knn leaf size": range(3,10,1),
                       "knn p": [1,2]
         knnclf = GridSearchCV(pipe, param grid, scoring='f1', cv=sss)
         knnclf.fit(features, labels)
         bestknn = knnclf.best estimator
         bestparams = knnclf.best params
         bestscore = knnclf.best score
         bestscorer = knnclf.scorer
         bestcv = knnclf.cv results
         print "Best KNN"
         print bestknn
         print bestparams
         print bestscore
         print bestscorer
         Best KNN
         Pipeline(steps=[('scaling', MinMaxScaler(copy=True, feature range=(0
         , 1))), ('knn', KNeighborsClassifier(algorithm='auto', leaf size=3,
         metric='minkowski',
                    metric params=None, n jobs=1, n neighbors=1, p=2,
                    weights='uniform'))))
         {'knn__leaf_size': 3, 'knn__algorithm': 'auto', 'knn__n_neighbors':
         1, 'knn p': 2}
         0.446904761905
```

```
In [30]: data = featureFormat(my dataset, features list, sort keys = True)
         labels, features = targetFeatureSplit(data)
         K best = SelectKBest(k=1)
         features kbest = K best.fit transform(features, labels)
         knn = KNeighborsClassifier()
         scaler = MinMaxScaler()
         sss = StratifiedShuffleSplit(n splits = 100, test size=0.3, random stat
         e = 42)
         pipe = Pipeline(steps=[('scaling', scaler), ("knn", knn)])
         param grid = {"knn n neighbors": [1, 3, 5, 8, 10, 12, 14],
                       "knn algorithm": ["auto", "ball tree", "kd tree", "brute
         "],
                       "knn leaf size": range(3,10,1),
                       "knn p": [1,2]
         knnclf = GridSearchCV(pipe, param grid, scoring='f1', cv=sss)
         knnclf.fit(features, labels)
         bestknn = knnclf.best estimator
         bestparams = knnclf.best params
         bestscore = knnclf.best score
         bestscorer = knnclf.scorer
         bestcv = knnclf.cv results
         print "Best KNN"
         print bestknn
         print bestparams
         print bestscore
         print bestscorer
         Best KNN
         Pipeline(steps=[('scaling', MinMaxScaler(copy=True, feature range=(0
         , 1))), ('knn', KNeighborsClassifier(algorithm='auto', leaf size=3,
         metric='minkowski',
                    metric params=None, n jobs=1, n neighbors=1, p=1,
                    weights='uniform'))))
         {'knn__leaf_size': 3, 'knn__algorithm': 'auto', 'knn__n_neighbors':
         1, 'knn p': 1}
         0.371892385392
```

```
In [31]: data = featureFormat(my dataset, features list, sort keys = True)
         labels, features = targetFeatureSplit(data)
         K best = SelectKBest(k=1)
         features kbest = K best.fit transform(features, labels)
         knn = KNeighborsClassifier()
         scaler = MinMaxScaler()
         sss = StratifiedShuffleSplit(n splits = 100, test size=0.1,random stat
         e = 42)
         pipe = Pipeline(steps=[('scaling', scaler), ("knn", knn)])
         param grid = {"knn n neighbors": [1, 3, 5, 8, 10, 12, 14],
                       "knn algorithm": ["auto", "ball tree", "kd tree", "brute
         "],
                       "knn leaf size": range(3,10,1),
                       "knn p": [1,2]
         knnclf = GridSearchCV(pipe, param grid, scoring='f1', cv=sss)
         knnclf.fit(features, labels)
         bestknn = knnclf.best estimator
         bestparams = knnclf.best params
         bestscore = knnclf.best score
         bestscorer = knnclf.scorer
         bestcv = knnclf.cv results
         print "Best KNN"
         print bestknn
         print bestparams
         print bestscore
         print bestscorer
         Best KNN
         Pipeline(steps=[('scaling', MinMaxScaler(copy=True, feature range=(0
         , 1))), ('knn', KNeighborsClassifier(algorithm='auto', leaf size=3,
         metric='minkowski',
                    metric params=None, n jobs=1, n neighbors=1, p=2,
                    weights='uniform'))))
         {'knn__leaf_size': 3, 'knn__algorithm': 'auto', 'knn__n_neighbors':
```

AdaBoostClassifier() with DecisionTreeClassifier()

1, 'knn_p': 2} 0.446904761905

```
DTC = DecisionTreeClassifier(random state = 42, max features = "auto",
In [49]:
         class weight = "balanced", max depth = None)
         ada = AdaBoostClassifier(base estimator=DTC)
         scaler = MinMaxScaler()
         sss = StratifiedShuffleSplit(n splits = 100, test size=0.1, random stat
         e = 42)
         pipe = Pipeline(steps=[('scaling',scaler),("ada", ada)])
         param grid = {"ada base estimator criterion" : ["gini", "entropy"],
                       "ada__base_estimator__splitter" : ["best", "random"],
                       "ada n estimators": [1, 10, 50, 100, 200]
                      }
         adaclf = GridSearchCV(pipe, param grid, scoring='f1', cv=sss)
         adaclf.fit(features, labels)
         bestada = adaclf.best estimator
         bestparams = adaclf.best_params_
         bestscore = adaclf.best score
         bestscorer = adaclf.scorer
         bestcv = adaclf.cv results
         print "Best Ada"
         print bestada
         print bestparams
         print bestscore
         print bestscorer
         Best Ada
         Pipeline(steps=[('scaling', MinMaxScaler(copy=True, feature range=(0
         , 1))), ('ada', AdaBoostClassifier(algorithm='SAMME.R',
         base estimator=DecisionTreeClassifier(class weight='balanced', crite
         rion='entropy',
                     max depth=None, max features='auto', max leaf nodes=None
                     min impur...dom state=42, splitter='random'),
```

learning rate=1.0, n estimators=10, random state=None))])

{'ada__base_estimator__splitter': 'random', 'ada__n_estimators': 10,

AdaBoostClassifier() with GaussianNB()

0.324214285714

make scorer(f1 score)

'ada base estimator criterion': 'entropy'}

```
In [51]: | NB = GaussianNB()
         adaN = AdaBoostClassifier(base_estimator=NB)
         scaler = MinMaxScaler()
         sss = StratifiedShuffleSplit(n splits = 100, test size=0.1, random stat
         e = 42)
         pipe = Pipeline(steps=[('scaling',scaler),("ada", adaN)])
         param_grid = {
                        "ada n estimators": [1, 10, 50, 100, 200]
         adaclf = GridSearchCV(pipe, param grid, scoring='f1', cv=sss)
         adaclf.fit(features, labels)
         bestada = adaclf.best estimator
         bestparams = adaclf.best_params_
         bestscore = adaclf.best_score_
         bestscorer = adaclf.scorer_
         bestcv = adaclf.cv results
         print "Best Ada"
         print bestada
         print bestparams
         print bestscore
         print bestscorer
```

```
/Users/robertlee/anaconda/envs/py2/lib/python2.7/site-packages/sklea
rn/naive bayes.py:232: RuntimeWarning: invalid value encountered in
divide
 new mu = np.average(X, axis=0, weights=sample weight / n new)
/Users/robertlee/anaconda/envs/py2/lib/python2.7/site-packages/sklea
rn/naive bayes.py:234: RuntimeWarning: invalid value encountered in
divide
 weights=sample_weight / n_new)
/Users/robertlee/anaconda/envs/py2/lib/python2.7/site-packages/sklea
rn/naive bayes.py:427: RuntimeWarning: divide by zero encountered in
  jointi = np.log(self.class prior [i])
/Users/robertlee/anaconda/envs/py2/lib/python2.7/site-packages/sklea
rn/ensemble/weight boosting.py:519: RuntimeWarning: invalid value en
countered in less
 proba[proba < np.finfo(proba.dtype).eps] = np.finfo(proba.dtype).e</pre>
/Users/robertlee/anaconda/envs/py2/lib/python2.7/site-packages/sklea
rn/ensemble/weight boosting.py:288: RuntimeWarning: invalid value en
countered in less
 proba[proba < np.finfo(proba.dtype).eps] = np.finfo(proba.dtype).e</pre>
ps
/Users/robertlee/anaconda/envs/py2/lib/python2.7/site-packages/sklea
rn/ensemble/weight boosting.py:603: RuntimeWarning: invalid value en
countered in greater
  return self.classes .take(pred > 0, axis=0)
/Users/robertlee/anaconda/envs/py2/lib/python2.7/site-packages/sklea
rn/ensemble/weight boosting.py:531: RuntimeWarning: invalid value en
countered in less
  (estimator weight < 0)))</pre>
/Users/robertlee/anaconda/envs/py2/lib/python2.7/site-packages/sklea
rn/ensemble/weight boosting.py:530: RuntimeWarning: invalid value en
countered in greater
  ((sample weight > 0)
Best Ada
Pipeline(steps=[('scaling', MinMaxScaler(copy=True, feature_range=(0)
, 1))), ('ada', AdaBoostClassifier(algorithm='SAMME.R',
          base estimator=GaussianNB(priors=None), learning rate=1.0,
          n estimators=1, random state=None))])
{'ada__n_estimators': 1}
0.253238095238
```

Dumping classifer, dataset, and feature_list to .pkl

```
In [ ]: | clf = best
        CLF PICKLE FILENAME = "my classifier.pkl"
        DATASET_PICKLE_FILENAME = "my_dataset.pkl"
        FEATURE LIST FILENAME = "my feature list.pkl"
        def dump classifier and data(clf, dataset, feature list):
            with open(CLF PICKLE_FILENAME, "w") as clf_outfile:
                pickle.dump(clf, clf outfile)
            with open(DATASET PICKLE FILENAME, "w") as dataset outfile:
                pickle.dump(dataset, dataset_outfile)
            with open(FEATURE LIST FILENAME, "w") as featurelist outfile:
                pickle.dump(feature list, featurelist outfile)
        dump classifier and data(bestknn, data dict, features list)
In [ ]: ## Checking the written files
In [ ]: with open("my_classifier.pkl", "r") as file:
            clf = pickle.load(file)
        print clf
In [ ]: | with open("my_dataset.pkl", "r") as file:
            data = pickle.load(file)
        for person in data:
            print person, data[person]
            break
In [ ]: with open("my feature list.pkl", "r") as file:
            features = pickle.load(file)
        print features
In [ ]:
```