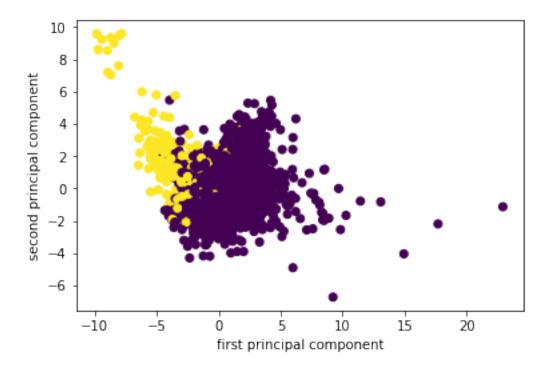
aml_hw4_task3_4

April 4, 2018

```
In [13]: X_data = pd.DataFrame(data["X"])
In [14]: from sklearn.preprocessing import StandardScaler
         scaler = StandardScaler()
         X_sc = scaler.fit_transform(X_data)
In [15]: y_sc = np.array(y_data)
In [16]: y_sc = y_sc.reshape(7200,)
In [19]: y_sc.shape
Out[19]: (7200,)
In [161]: from sklearn.decomposition import PCA
          import matplotlib.pyplot as plt
          %matplotlib
          #print(cancer.data.shape)
          pca = PCA(n_components=2)
          X_pca = pca.fit_transform(X_sc)
Using matplotlib backend: TkAgg
In [8]: print(X_pca.shape)
        print(y_data.shape)
```

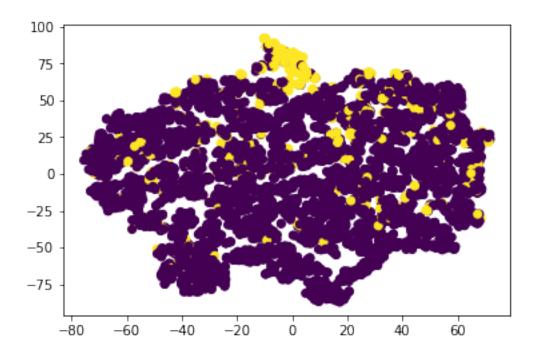
```
(7200, 2)
(7200, 1)
```



```
[ 0.374, 30.941],
[-59.743, 24.545],
[ 16.716, 49.189]], dtype=float32)
```

In []:

Out[209]: <matplotlib.collections.PathCollection at 0x7f7592fc0c18>



```
In [27]: X_sc
```

In []:

In []:

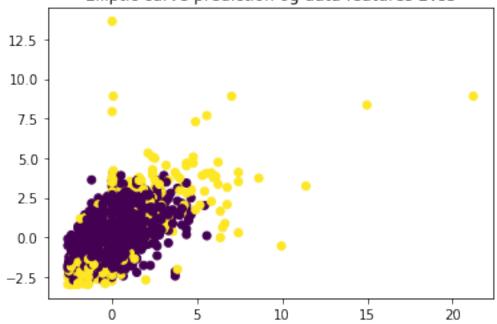
2 Task 3 Outlier Detection

2.0.1 3.1

Elliptic curve prediction og data features 2vs3

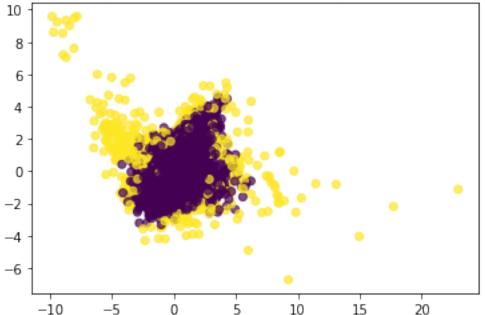
plt.scatter(X_sc[:, 2], X_sc[:, 3], c=pred_ee)

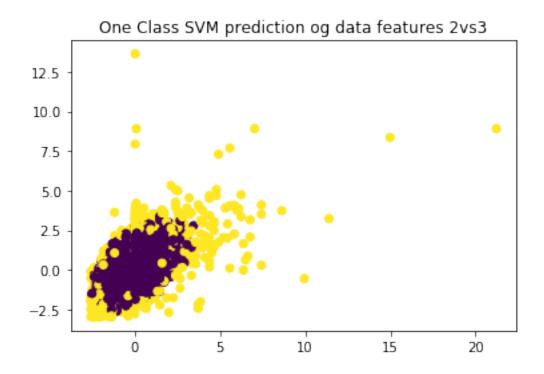
Out[211]: <matplotlib.collections.PathCollection at 0x7f7592f5bf98>



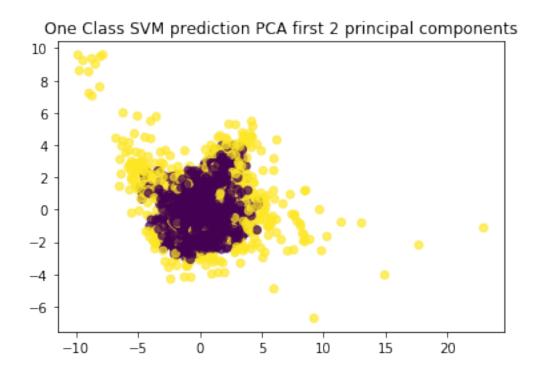
Out[212]: <matplotlib.collections.PathCollection at 0x7f7592efaac8>



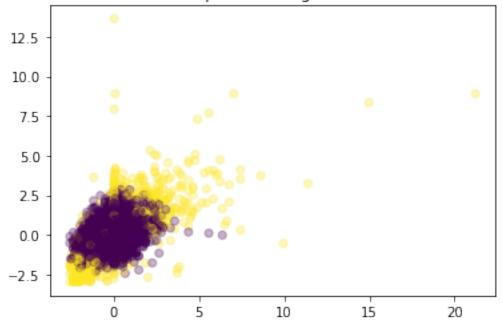




Out[217]: <matplotlib.collections.PathCollection at 0x7f7592e052b0>

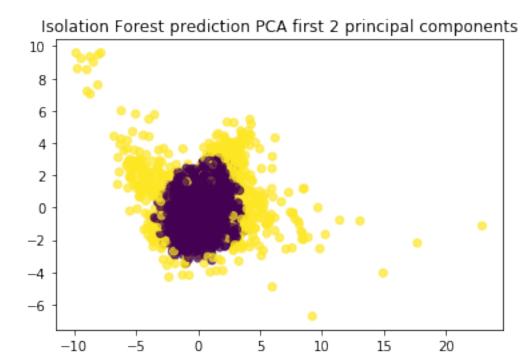


Isolation Forest prediction og data features 2vs3



```
In [218]: plt.title("Isolation Forest prediction PCA first 2 principal components")
    plt.scatter(X_pca[:, 0], X_pca[:, 1], c=pred_if, alpha = 0.7)
```

Out[218]: <matplotlib.collections.PathCollection at Ox7f7592e05cc0>



2.0.2 Without using the ground truth we cannot really gauge how good an outlier detection algorithm works as the visualizing the predictions of all three models on X_pca leads to similar looking plots. Not much can be inferred from them

As we can see only the Isolation Forest decision function provides output values between 0 and 1

2.0.3 3.2

```
In [65]: from sklearn.metrics import average_precision_score
In []: for i in range(len(y_sc)):
            #print(y_sc[i])
        for i in range(len(y_sc)):
            if(pred_ee[i] == 1):
                pred_ee[i] = 0
            else:
                pred_ee[i] = 1
            print(pred_ee[i])
        for i in range(len(y_sc)):
            if(pred_oc[i] == 1):
                pred_oc[i] = 0
            else:
                pred_oc[i] = 1
            print(pred_oc[i])
        for i in range(len(y_sc)):
            if(pred_if[i] == 1):
                pred_if[i] = 0
            else:
                pred_if[i] = 1
            print(pred_if[i])
In [230]: average_precision_ee = average_precision_score(y_sc, pred_ee)
          average_precision_oc = average_precision_score(y_sc, pred_oc)
          average_precision_if = average_precision_score(y_sc, pred_if)
          print('Elliptic Envelope Average precision-recall score: {0:0.2f}'.format(
                average_precision_ee))
          print('One Class SVM Average precision-recall score: {0:0.2f}'.format(
                average_precision_oc))
          print('Isolation Forest Average precision-recall score: {0:0.2f}'.format(
                average_precision_if))
Elliptic Envelope Average precision-recall score: 0.25
One Class SVM Average precision-recall score: 0.11
Isolation Forest Average precision-recall score: 0.15
```

```
In [121]: average_precision_ee = average_precision_score(y_sc, pred_ee)
         average_precision_oc = average_precision_score(y_sc, pred_oc)
         average_precision_if = average_precision_score(y_sc, pred_if)
         print('Elliptic Envelope Average precision-recall score: {0:0.2f}'.format(
               average_precision_ee))
         print('One Class SVM Average precision-recall score: {0:0.2f}'.format(
               average_precision_oc))
         print('Isolation Forest Average precision-recall score: {0:0.2f}'.format(
               average_precision_if))
Elliptic Envelope Average precision-recall score: 0.25
One Class SVM Average precision-recall score: 0.11
Isolation Forest Average precision-recall score: 0.15
In [122]: from sklearn.metrics import roc_auc_score
         from sklearn.metrics import zero_one_loss
         ee_auc = roc_auc_score(y_sc, pred_ee)
         oc_auc = roc_auc_score(y_sc, pred_oc)
         if_auc = roc_auc_score(y_sc, pred_if)
         ee_loss = zero_one_loss(y_sc, pred_ee)
         oc_loss = zero_one_loss(y_sc, pred_oc)
         if_loss = zero_one_loss(y_sc, pred_if)
         print("AUC for Elliptic Envelope is ", ee_auc)
         print("AUC for One Class SVM: ", oc_auc)
         print("AUC for Isolated Forest: ", if_auc)
         print("Zero-one loss for Elliptic Envelope is ", ee_loss)
         print("Zero-one loss for One Class SVM: ", oc_loss)
         print("Zero-one loss for Isolated Forest: ", if_loss)
AUC for Elliptic Envelope is 0.7077235813468987
AUC for One Class SVM: 0.5841154340153115
AUC for Isolated Forest: 0.6288392884232243
Zero-one loss for Elliptic Envelope is 0.080277777777778
Zero-one loss for One Class SVM: 0.114583333333333333
```

2.0.4 The AUC and average precision readings suggest that EllipticEnvelope is the best

The clustering approaches from task 2 yield poor results with in some cases 3 clusters giving the best results. Thus, the approaches in Task 3 which assume the additional information that you know the number of classes and their proportion gives better results

However comaparing the visualizations in Task 3 with the ground truth clustering of the PCA visualization in Task 1 we notice that even Task 3 is quite off. Therefore there is no tangible way of comparing one model from another without knowing the ground truth and just via visualization. This could be the result of the data being highly non Gaussian

3 Task 4

```
In [191]: from sklearn.model_selection import train_test_split
          X_train, X_test, y_train, y_test = train_test_split(X_sc, y_sc, stratify = y_sc)
Imbalanced Learning
In [195]: from sklearn.model_selection import cross_validate
          from sklearn.linear_model import LogisticRegression
          scores = cross_validate(LogisticRegression(),
                                  X_sc, y_sc, cv=10, scoring=('roc_auc', 'average_precision'))
          scores['test_roc_auc'].mean(), scores['test_average_precision'].mean()
Out [195]: (0.9726447870177892, 0.7637355829525652)
In [196]: from sklearn.ensemble import RandomForestClassifier
          scores = cross_validate(RandomForestClassifier(n_estimators=100),
                                  X_sc, y_sc, cv=10, scoring=('roc_auc', 'average_precision'))
          scores['test_roc_auc'].mean(), scores['test_average_precision'].mean()
Out [196]: (0.9957853471700322, 0.9321881275718397)
Balanced weights
In [197]: scores = cross_validate(LogisticRegression(class_weight='balanced'),
                                  X_sc, y_sc, cv=10, scoring=('roc_auc', 'average_precision'))
         print("Test ROC AUC for Logistic :", scores['test_roc_auc'].mean(),"| Average Precision
Test ROC AUC for Logistic: 0.9884794372602356 | Average Precision for Logistic: 0.828698915304
In [198]: scores = cross_validate(RandomForestClassifier(n_estimators=100, class_weight='balance
                                  X_sc, y_sc, cv=10, scoring=('roc_auc', 'average_precision'))
         print("Test ROC AUC for Random Forest :", scores['test_roc_auc'].mean(),"| Average Pre
```

Test ROC AUC for Random Forest: 0.9957258147797312 | Average Precision for Random Forest: 0.92

Grid Searching C for LogReg and other pruning parameters for the trees

```
In [133]: from sklearn.model_selection import GridSearchCV
In [134]: param_grid_rf = {
              'max_depth': [80, 90, 100, 110],
              'max_features': [2, 3],
              'n_estimators': [100, 200, 300, 1000]
          }
In [135]: rf_grid = GridSearchCV(RandomForestClassifier(), param_grid=param_grid_rf, cv=3)
In [136]: rf_grid.fit(X_train,y_train)
Out[136]: GridSearchCV(cv=3, error_score='raise',
                 estimator=RandomForestClassifier(bootstrap=True, class_weight=None, criterion='
                      max_depth=None, max_features='auto', 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=10, n_jobs=1,
                      oob_score=False, random_state=None, verbose=0,
                      warm_start=False),
                 fit_params=None, iid=True, n_jobs=1,
                 param_grid={'max_depth': [80, 90, 100, 110], 'max_features': [2, 3], 'n_estimat
                 pre_dispatch='2*n_jobs', refit=True, return_train_score='warn',
                 scoring=None, verbose=0)
In [137]: rf_grid.score(X_test,y_test)
Out[137]: 0.98333333333333333
In [143]: print("AUC for RF is ", roc_auc_score(y_test, rf_grid.predict(X_test)))
          print("Avg Precision for RF: ", average_precision_score(y_test, rf_grid.predict(X_tes
AUC for RF is 0.9737045072188569
Avg Precision for RF: 0.8079279080053073
In [199]: param_grid_lr = {
             'C': [ 0.1, 1,2,5 ,10,100]
In [200]: lr_grid = GridSearchCV(LogisticRegression(), param_grid=param_grid_lr, cv=10)
In [201]: lr_grid.fit(X_train,y_train)
```

3.0.1 Changing to balanced weights gives the best avg precision and recall scores. So yes. It helps

In []: