Linear Support Vector Classifier (SVC)

Declaration: The central idea and coding is abstract from Kevin mark ham youtube video seriese, Introduction to machine learning with scikit-learn video series. You can find link under resources section.

What are the features?

- trip_id: A unique number to identify each trip
- From station Number: From station number where the trip Start
- Day: Day of the trip for example Monday, Tuesday etc.
- Month: Which month trip took place
- Duration: Total trip duration in minutes
- · birthyear: Birth year of user
- · Sex: Gender identification of user
- · age: Current age of user

What is the response?

• Station Number: To Station Number where the trip ends

Question: Can we predict the End station of a bicycle trip by given trip dataset?

Libraries

```
In [260]:
```

```
import os,csv,io,mapsplotlib,time,folium,googlemaps,geopy,zipfile,requests,
warnings
import numpy as np
import pandas as pd
import datetime as dt
import seaborn as sns
import geopandas as gpd
from shapely.geometry import Point
import statsmodels.formula.api as smf
import matplotlib.pyplot as plt
import mysql.connector as sql
from sklearn.linear model import LinearRegression
model = LinearRegression()
from sklearn import metrics
from sklearn.cross validation import train test split
from sklearn.svm import LinearSVC
import numpy as np
warnings.simplefilter('ignore')
```

```
In [261]:
```

```
# display plots in the notebook
```

It is difficult to obtain better accuracy with all end stations (to_station). With all end station the model predicts 0.067 percent of accuracy, to avoid this this study focus on a circumscribed number of end station. From our following query we are supposing here that 'it does not matter where the bicycle trip will go. We are more interested in where this trip ends and to attain this goal. We have five main busiest end stations. Let's say's if a journey start from x, y and Z station what is the chances it will end on the given end stations.

In [286]:

```
#importing data from database
db connection = sql.connect(host='localhost', database='bike', user='root',
password='none')
db cursor = db connection.cursor()
db cursor.execute('SELECT trip id, tripduration minutes
Duration, from station name, from station id num, to station id, to station nam
o station id num, Day num Day, bmonth Month, year Year, age, Sex num Sex, sthours
Hours FROM bike.trip clean where to station id num in(7,9,36,37,38) order b
y to station id num;')
table rows = db cursor.fetchall()
data = pd.read sql('SELECT trip id, tripduration minutes
Duration, from station name, from station id num, to station id, to station nam
o station id num, Day num Day, bmonth Month, year Year, age, Sex num Sex, sthours
Hours FROM bike.trip clean where to station id num in(7,9,36,37,38) order b
y to station id num;', con=db connection)
df = pd.DataFrame(data)
df.tail()
```

Out[286]:

	trip_id	Duration	from_station_name	from_station_id_num	to_station_id	to_station_
25845	254994	8.23	Dexter Ave N & Aloha St	34	SLU-15	6th Ave
25846	254998	9.15	E Thomas St	14	SLU-15	6th Ave
25847	255058	7.89	E Pine St	16	SLU-15	6th Ave
25848	255087	8.03	E Pine St	16	SLU-15	6th Ave
25849	255149	5.97	Elliott Ave	1	SLU-15	6th Ave

In [287]:

```
# define X and y
feature_cols = ['Duration','from_station_id_num','Hours','age','Sex']
X = df[feature_cols]
y = df.to_station_id_num
```

In [288]:

```
clf = LinearSVC(loss = '12') # instantiated with L2 loss
print (clf)
LinearSVC(C=1.0, class weight=None, dual=True, fit intercept=True,
```

intercept_scaling=1, loss='12', max_iter=1000, multi_class='ovr',

```
In [289]:

clf = clf.fit(X, y)
```

creating an instance of the classifier. This can be done simply by calling the class name, with any arguments that the object accepts:

```
In [290]:
clf = LinearSVC(loss = '12')
```

clf is a statistical model that has parameters that control the learning algorithm (those parameters are sometimes called the *hyperparameters*). Those hyperparameters can be supplied by the user in the constructor of the model. We will discuss later how to choose a good combination using either simple empirical rules or data driven selection:

```
In [291]:
print (clf)
LinearSVC(C=1.0, class_weight=None, dual=True, fit_intercept=True,
   intercept_scaling=1, loss='12', max_iter=1000, multi_class='ovr',
   penalty='12', random_state=None, tol=0.0001, verbose=0)
```

By default the model parameters are not initialized. They will be tuned automatically from the data by calling the fit method with the data X and labels Y:

```
In [292]:
clf = clf.fit(X, y)
```

We can now see some of the fit parameters within the classifier object.

In scikit-learn, parameters defined by training have a trailing underscore.

Once the model is trained, it can be used to predict the most likely outcome on unseen data.

```
In [297]:
```

```
#here is the from station number. What end station does this get predicted
as?
X_new = [[22, 31, 35, 7, 9]]
clf.predict(X_new) #end station number 37
Out[297]:
array([37], dtype=int64)
```

All classification tasks involve predicting an unknown category based on observed features.

Probabilistic Prediction

```
In [298]:
```

```
from sklearn.linear_model import LogisticRegression
clf2 = LogisticRegression()
clf2.fit(X, y)
print (clf2.predict_proba(X_new))

[[ 0.26817904   0.11485692   0.0005781   0.56730664   0.0490793 ]]

In [300]:
clf2.predict(X_new) #end station number 37

Out[300]:
array([37], dtype=int64)
```

The result gives the probability (between zero and one) that the test point comes from any of the three classes.

This means that the model estimates that the sample in X_new has:

- 90% likelyhood to belong to the 'to station' class (end station = 37)
- 9% likelyhood to belong to the 'to station' class (end station = 36)
- < 1% likelyhood to belong to the 'to station' class (end station = 38)

Evaluating the Model

Let's get a rough evaluation our model by using it to predict the values of the training data:

```
In [301]:
y_model = clf2.predict(X)

In [302]:
print (y_model == y)

0     False
```

\perp	raise	
2	False	
3	False	
4	False	
5	False	
6		
	False	
7	False	
8	True	
9	False	
10	False	
11	False	
12	False	
13	False	
14	False	
15	True	
16	False	
17	True	
18	False	
19	False	
20	False	
21	False	
22	False	
23	False	
24	False	
25	False	
26	False	
27	False	
28	False	
29	False	
	•••	
25820	False	
25821	False	
25822	False	
25823	False	
25824	False	
25825	False	
25826	False	
25827	False	
25828	False	
25829	False	
25830	False	
25831	False	
25832	False	
25833	False	
25834	False	
25835	False	
25836	False	
25837	False	
25838	False	
25839	False	
25840	True	
25841	True	
25842	False	
25843	False	
25844	False	
25845	False	
25846	False	
25847	False	
25848	True	
25849	False	
Nama. +		

We see that most of the predictions are incorrect!

Measuring Classification Performance: Validation & Testing

Cross-Validation

Learning the parameters of a prediction function and testing it on the same data is a methodological mistake: a model that would just repeat the labels of the samples that it has just seen would have a perfect score but would fail to predict anything useful on yet-unseen data.

To avoid over-fitting, we have to define two different sets:

- a training set X_train, y_train which is used for learning the parameters of a predictive model
- a testing set X_test, y_test which is used for evaluating the fitted predictive model

In scikit-learn such a random split can be quickly computed with the <code>train_test_split</code> helper function. It can be used this way:

In [306]:

```
from sklearn import cross_validation
X_train, X_test, y_train, y_test = cross_validation.train_test_split(X, y, t
est_size=0.25, random_state=0) # 25% of test data
print (X.shape, X_train.shape, X_test.shape) # Training data => (6463, 5)
(25850, 5) (19387, 5) (6463, 5)
```

Now we train on the training data, and test on the testing data:

In [307]:

```
clf = LinearSVC(loss='12').fit(X train, y train)
y pred = clf.predict(X test)
print (y pred == y test)
15014 False
15095
        True
8653
        True
15333 False
       False
22750
       False
10281
10227
       False
3861
       False
19032
       False
       False
6662
10139
        True
1799
        True
21841
       False
       False
4043
4725
        True
24434
       False
```

```
21407
       False
21838
       False
13339
       False
20510
       False
4285
      False
12461
      False
19098
      False
5619
      False
379
       False
21396
       False
5192
      False
3822
      False
15548
       False
25161
       False
        . . .
25166
       False
5274
        True
20725
       False
11758
       False
       False
25186
14466
       False
24563
       False
13321
       False
20033
      False
       False
20672
1143
        True
23027
       False
        True
6331
25178
       False
16581
      False
19207
       False
15234
       False
13045
       False
        True
6644
16760
      False
23173 False
        True
7597
16136
        True
17487
      False
        True
878
467
       False
7016
      False
5536
      False
15761
       False
5583
       False
Name: to station id num, dtype: bool
```

There is an issue here, however: by defining these two sets, we drastically reduce the number of samples which can be used for learning the model, and the results can depend on a particular random choice for the pair of (train, test) sets.

A solution is to split the whole data several consecutive times in different train set and test set, and to return the averaged value of the prediction scores obtained with the different sets. Such a procedure is called **cross-validation**. This approach can be computationally expensive, but does not waste too much data (as it is the case when fixing an arbitrary test set), which is a major advantage in problem such as inverse inference where the number of samples is very small.

We'll explore cross-validation, but for more information on cross-validation in scikit-learn here:

Cross-validation for parameter tuning, model selection, and feature selection

In [308]:

```
from sklearn.cross_validation import train_test_split
from sklearn.neighbors import KNeighborsClassifier
from sklearn import metrics
# use train/test split with different random_state values
X_train, X_test, y_train, y_test = train_test_split(X, y, random_state=4)
# check classification accuracy of KNN with K=5
knn = KNeighborsClassifier(n_neighbors=5)
knn.fit(X_train, y_train)
y_pred = knn.predict(X_test)
print ((metrics.accuracy_score(y_test, y_pred))*100)
```

72.1646294291

we created a bunch of train/test splits, calculated the testing accuracy for each, and averaged the results together and That's the essense of **cross-validation!**

Method Of K-cross Validation

- 1. Split the dataset into K equal partitions (or "folds").
- 2. Use fold 1 as the **testing set** and the union of the other folds as the **training set**.
- 3. Calculate **testing accuracy**.
- 4. Repeat steps 2 and 3 K times, using a different fold as the testing set each time.
- 5. Use the average testing accuracy as the estimate of out-of-sample accuracy.

Parameter tuning

Goal: Select the best tuning parameters (aka "hyperparameters") for KNN

```
In [315]:
```

```
from sklearn.cross_validation import cross_val_score
```

```
In [316]:
```

```
# 10-fold cross-validation with K=5 for KNN (the n_neighbors parameter)
knn = KNeighborsClassifier(n_neighbors=5)
scores = cross_val_score(knn, X, y, cv=10, scoring='accuracy')
print (scores)
```

```
[ 0.76352396  0.76334107  0.76836814  0.65390565  0.68201161  0.70085139  0.68846749  0.70123839  0.68537152  0.59349593]
```

```
In [317]:
```

```
# use average accuracy as an estimate of out-of-sample accuracy
```

```
print ((scores.mean())*100)
70.0057513886
In [318]:
# search for an optimal value of K for KNN
```

```
# search for an optimal value of K for KNN
k_range = range(1, 31)
k_scores = []
for k in k_range:
    knn = KNeighborsClassifier(n_neighbors=k)
    scores = cross_val_score(knn, X, y, cv=10, scoring='accuracy')
    k_scores.append(scores.mean())
print (k_scores)
```

```
 \begin{bmatrix} 0.77018611022881145, & 0.71980910690268751, & 0.70778630842517321, \\ 0.70191100103433901, & 0.70005751388561444, & 0.69456571797540645, \\ 0.68775720442010313, & 0.68625010043589507, & 0.68338717619918021, \\ 0.67882406712522447, & 0.67503423561828779, & 0.67031535609937565, \\ 0.66540247068146352, & 0.66145728601652398, & 0.65801557129775445, \\ 0.65600341199435996, & 0.65279269844244692, & 0.65112880600595235, \\ 0.65074210827620882, & 0.64799576192400632, & 0.6468742639994598, \\ 0.64521028196917174, & 0.64385661520741366, & 0.641302869332942, \\ 0.63863440041152075, & 0.63697047844981969, & 0.6358491001533868, \\ 0.63352840515897724, & 0.63105301602567088, & 0.62938949864783933 \end{bmatrix}
```

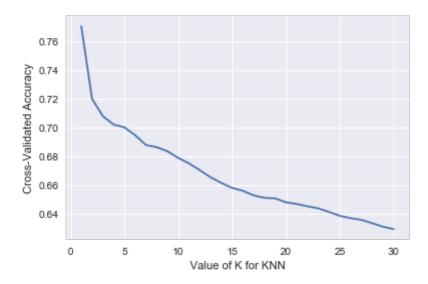
In [319]:

```
import matplotlib.pyplot as plt
%matplotlib inline

# plot the value of K for KNN (x-axis) versus the cross-validated accuracy
(y-axis)
plt.plot(k_range, k_scores)
plt.xlabel('Value of K for KNN')
plt.ylabel('Cross-Validated Accuracy')
```

Out[319]:

<matplotlib.text.Text at 0x1bcf1abe4e0>



The following graphs display the lower value of k is better for our model accuracy and with K=1 we will get max of 76 percent of accuracy which is ok.

Model selection

Goal: Compare the best KNN model with logistic regression

```
In [322]:
```

```
# 10-fold cross-validation with the best KNN model
knn = KNeighborsClassifier(n_neighbors=2)
print ((cross_val_score(knn, X, y, cv=10, scoring='accuracy').mean())*100)
```

71.9809106903

```
In [323]:
```

```
# 10-fold cross-validation with logistic regression
from sklearn.linear_model import LogisticRegression
logreg = LogisticRegression()
print ((cross_val_score(logreg, X, y, cv=10, scoring='accuracy').mean())*10
0)
```

26.9206953717

From the above result we are selecting KNN Neighbours with K=1 which is produce a better mean accuracy which is 72 percent aproximately. As compare to logistic regression. The logistic regression offer mean accuracy of 26.92 which is approximately 27 percent that is very low. So, because of that result, this study will use KNN Neighbours with k=1 for our predictions.

Improvements to cross-validation

Repeated cross-validation

- Repeat cross-validation multiple times (with different random splits of the data) and average the results
- More reliable estimate of out-of-sample performance by reducing the variance associated with a single trial of cross-validation

Creating a hold-out set

- "Hold out" a portion of the data **before** beginning the model building process
- Locate the best model using cross-validation on the remaining data, and test it using the hold-out set
- More reliable estimate of out-of-sample performance since hold-out set is truly out-of-sample

Feature engineering and selection within cross-validation iterations

Normally, feature engineering and selection occurs before cross-validation

- Instead, perform all reature engineering and selection within each cross-validation iteration
- More reliable estimate of out-of-sample performance since it better mimics the application of the model to out-of-sample data

Resources

References: From the video series: Introduction to machine learning with scikit-learn

- scikit-learn documentation: Cross-validation, Model evaluation
- scikit-learn issue on GitHub: MSE is negative when returned by cross_val_score
- Section 5.1 of <u>An Introduction to Statistical Learning</u> (11 pages) and related videos: <u>K-fold and leave-one-out cross-validation</u> (14 minutes), <u>Cross-validation the right and wrong ways</u> (10 minutes)
- Scott Fortmann-Roe: <u>Accurately Measuring Model Prediction Error</u>
- Machine Learning Mastery: An Introduction to Feature Selection
- Harvard CS109: <u>Cross-Validation: The Right and Wrong Way</u>
- Journal of Cheminformatics: <u>Cross-validation pitfalls when selecting and assessing</u> regression and classification models