# **Assignment 3**

In this assignment, you'll explore linear regression and logistic regression, the perceptron, and support vector machines: popular approaches which share a similar form of producing output in the form of a weighted combination of attributes, but with very different approaches to learning the underyling model. The assignment explores two of the general data mining tasks: regression and classification, and takes a closer look at multiclass classification with these techniques.

Similar to previous assignments, you're expected to respond to each question with your answer in a Markdown cell and clearly labeled code supporting your answer in a code cell. When you submit the assignment, you should upload the two notebooks (.ipynb files) corresponding to your solutions and also generate a PDF of each notebook that includes the answers, code, and all intermediate output. In total, you will submit four files: two notebooks and two PDFs generated from those notebooks. This assignment is due on 2/16/16 at 11:59pm.

# Part 1: Linear and Logistic Regression, Perceptrons (40 points)

We'll start by looking at a simple example of linear regression. As with many of the other data mining algorithms we've encountered, we'll be using the implementation available in the <code>linear\_model</code> module of scikit-learn. Now would be a good time to go read the documentation (<a href="http://scikit-learn.org/stable/modules/linear\_model.html">http://scikit-learn.org/stable/modules/linear\_model.html</a>) on linear models and refresh your memory on the high-level ideas. Once you've done that, we'll go through a simple example.

```
In [15]: ## Preliminaries
         #Show plots in the notebook
         %matplotlib inline
         from sklearn import datasets, preprocessing, cross validation, feature
         from sklearn import linear model, svm, metrics, ensemble
         import numpy as np
         import pandas as pd
         import matplotlib.pyplot as plt
         import urllib2
         # Helper functions
         def folds to split(data, targets, train, test):
             data tr = pd.DataFrame(data).iloc[train]
             data te = pd.DataFrame(data).iloc[test]
             labels tr = pd.DataFrame(targets).iloc[train]
             labels te = pd.DataFrame(targets).iloc[test]
             return [data tr, data te, labels tr, labels te]
```

# **Linear Regression on the Diabetes Dataset**

The first dataset we'll explore is the Diabetes dataset (http://scikit-

learn.org/stable/modules/generated/sklearn.datasets.load diabetes.html#sklearn.datasets.load one of the pre-packaged datasets available in scikit-learn. There aren't too many details about t dataset out there, but the common description is that it contains 10 physiological variables (age weight, blood pressure) for 442 patients, and the target is an indication of disease progression a year. Like any good data miners, let's poke around and check out the data before we get starter

```
In [16]: # Load the data
diabetes = datasets.load_diabetes();
# Put it into pandas DataFrames
diabetes_data_df = pd.DataFrame(diabetes.data);
diabetes_target_df = pd.DataFrame(diabetes.target)

# How many attributes and records are there?
print diabetes_data_df.shape

# What are the descriptive statistics?
print diabetes_data_df.describe()
# How are the labels distributed?
diabetes_target_df.hist()

# How are the attributes distributed?
plt.figure()
axes = diabetes_data_df.boxplot()
```

(442, 10)0 1 2 3 count 4.420000e+02 4.420000e+02 4.420000e+02 4.420000e+02 4.420 000e+02 mean -3.634285e-16 1.308343e-16 -8.045349e-16 1.281655e-16 -8.835 316e-17 4.761905e-02 4.761905e-02 4.761905e-02 4.761905e-02 4.761 std 905e-02 min -1.072256e-01 -4.464164e-02 -9.027530e-02 -1.123996e-01 -1.267807e-01 25% -3.729927e-02 -4.464164e-02 -3.422907e-02 -3.665645e-02 -3.42450% 5.383060e-03 -4.464164e-02 -7.283766e-03 -5.670611e-03 -4.320866e-03 75% 3.807591e-02 5.068012e-02 3.124802e-02 3.564384e-022.835 801e-02 5.068012e-02 1.705552e-01 1.320442e-01 max 1.107267e-01 137e-01 5 6 7 8 9 count 4.420000e+02 4.420000e+02 4.420000e+02 4.420000e+02 4.420 000e+02 mean 1.327024e-16 -4.574646e-16 3.777301e-16 -3.830854e-16 -3.412 882e-16 4.761905e-02 4.761905e-02 4.761905e-02 4.761905e-02 4.761 std 905e-02

min -1.156131e-01 -1.023071e-01 -7.639450e-02 -1.260974e-01 -1.377 672e-01 25% -3.035840e-02 -3.511716e-02 -3.949338e-02 -3.324879e-02 -3.317 903e-02 50% -3.819065e-03 -6.584468e-03 -2.592262e-03 -1.947634e-03 -1.077 698e-03

75% 2.984439e-02 2.931150e-02 3.430886e-02 3.243323e-02 2.791 705e-02

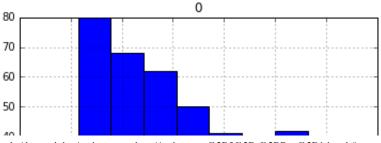
max 1.987880e-01 1.811791e-01 1.852344e-01 1.335990e-01 1.356 118e-01

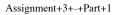
/Users/viktorjankov/anaconda/lib/python2.7/site-packages/ipykernel/\_ \_main\_\_.py:17: FutureWarning:

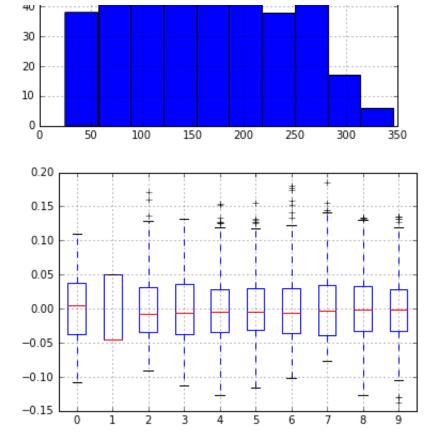
The default value for 'return\_type' will change to 'axes' in a futur e release.

To use the future behavior now, set return type='axes'.

To keep the previous behavior and silence this warning, set return\_type='dict'.







Notice anything interesting about the data? The attributes have been normalized - each has a similar mean and standard deviation. Now let's train our first regression model.

```
In [17]: | # First we'll do a simple train-test split:
         [dbt tr data, dbt te data,
          dbt tr target, dbt te target] = cross validation.train test split(diab
         # Create the LinearRegression classifier
         lr = linear model.LinearRegression()
         # Learn the linear regression model
         lr.fit(dbt tr data, dbt tr target)
         # Print out the coefficient of determination (R^2)
         print "R^2:\t", lr.score(dbt te data, dbt te target)
         # Peek at the predictions
         dbt_te_predict = lr.predict(dbt_te_data)
         print dbt te predict
         # And also the mean squared error:
         print "MSE:\t", metrics.mean squared error(dbt te target, dbt te predic
         # Which attributes were important to this prediction? We can find out b
         print lr.coef
```

R^2: 0.53371	.6155537			
[ 124.87699265 35	241.04779116	239.55002222	262.61331893	245.760759
215.38130957 48	98.25024738	167.10341193	86.15467868	149.231306
139.44688429	121.07843815	145.93392542	113.89242688	137.000019
59 106.07596308	188.8720697	151.64642212	147.17548161	193.976209
2 193.24395545	123.38347892	244.93128416	221.35017651	73.231771
37 175.0689274	270.81318019	76.43093304	147.64128689	106.932261
73 72.18329933	246.22062792	155.49372733	132.93631466	122.141629
70.14373998	182.1432308	259.16039335	119.69523429	217.322246
65 156.87184212		152.68147644	97.91484124	99.479911
66	209.45289148			
60.35999351	163.47427462	175.63241326	165.73475646	120.171374
143.1380446 62	65.50402188	161.25927147	165.59617116	127.563187
145.47983532 41	119.2772516	140.89228502	138.44957973	234.270316
70.42410076	219.75216763	96.49453974	169.97467564	75.491765
72 141.43569069	146.08650475	97.22882088	235.72934302	115.606263
81 140.86586191	83.8398695	226.79822353	232.14151232	153.569348
61 85.00662969	185.94162205	100.59978839	80.74525585	160.501532
47				
136.07622533 09		150.66490274	181.52512021	
205.67406873 13	142.71143086	127.42691032	139.56003045	151.354371
242.54261548 86	206.12199133	145.09312121	120.10821335	248.910441
177.46641126 24	145.28356072	79.89910472	155.73208421	113.432393
162.18337986	99.99524699	191.84896583	154.24943403	166.819089
55 104.90171597	220.98892005	134.51002344	38.0086503	203.293999
74 181.12997599]				
MSE: 2919.42				
[ 11.18121228 56	-219.22917848	492.74550531	360.62854712	-880.756963
560.47207888	138.68964051	208.07342467	723.26859805	47.007418
34]				

Hopefully you can see the similarities between how to build a <u>LinearRegression model (http://sclearn.org/stable/modules/generated/sklearn.linear model.LinearRegression.html#sklearn.linear and the previous assignment on DecisionTrees. The big difference is that the predictions are no values (much like the training targets). One consequence is that we can't use the same metrics classification. One metric that we've discussed (and that linear regression attempts to minimize <u>squared error (https://en.wikipedia.org/wiki/Mean squared error)</u> or MSE. If you understand how leap to the default metric that regression reports (via the <code>.score()</code> function) -- the <u>coefficient coefficient of determination</u> (R^2) -- shouldn't be too hard to unot the R^2 metric and make sure you understand what it measures.</u>

The above example used a simple train-test split, but we can be more confident of our conclusivalidation. Let's use cross-validation to experiment with the same dataset and take a look at the example provides you with some tricks that you might have missed in the last assignment, so up below and learning to use it should be helpful

```
In [18]: | foldnum = 0
         fold results = pd.DataFrame()
         for train, test in cross validation.KFold(len(diabetes.data), n folds=1
             foldnum+=1
             [dbt tr data, dbt te data,
              dbt tr target, dbt te target] = folds to split(diabetes.data,diabe
             lr = linear model.LinearRegression()
             lr.fit(dbt tr data, dbt tr target)
             # We could print out our results
             print "Fold %d\t\t R^2 metric = %03.3f \t\t MSE = %03.1f " % (foldn
                                                                             lr.sc
                                                                             metri
                                                                            )
             # But a nicer way to store them is in a DataFrame
             fold results.loc[foldnum, 'R^2'] = lr.score(dbt te data, dbt te tar
             fold results.loc[foldnum, 'MSE'] = metrics.mean squared error(dbt t
             # By the way, if you were searching over parameters in an inner for
             # you could store those in your results DataFrame just as easily, f
             # for param in params.keys():
             #
                  for paramVal in params[param]:
             #
                      paramDict={};
             #
                      paramDict[param]=paramVal
             #
                      dtree = tree.DecisionTreeClassifier(random state=20160121,
             #
                      dtree.fit(mushroom train, mushroom train labels)
                      fold results.loc[foldnum,'%s=%s' % (param, paramVal)]=dtre
         #Now let's look at the results:
         print fold results
         #And compute the mean error across folds:
         print fold results.mean()
```

```
Fold 1
                 R^2 metric = 0.556
                                                  MSE = 2533.8
Fold 2
                 R^2 metric = 0.231
                                                  MSE = 2870.8
Fold 3
                 R^2 metric = 0.354
                                                  MSE = 3512.7
Fold 4
                 R^2 metric = 0.622
                                                  MSE = 2759.2
Fold 5
                 R^2 metric = 0.266
                                                  MSE = 3555.7
Fold 6
                 R^2 metric = 0.618
                                                  MSE = 2900.4
Fold 7
                 R^2 metric = 0.418
                                                  MSE = 3696.3
Fold 8
                 R^2 metric = 0.435
                                                  MSE = 2282.3
Fold 9
                 R^2 metric = 0.434
                                                  MSE = 4122.9
Fold 10
                 R^2 metric = 0.686
                                                  MSE = 1769.7
         R^2
                      MSE
1
    0.556144
              2533.848109
2
    0.230561
              2870.767711
3
              3512.723509
    0.353578
4
    0.621905 2759.227129
5
    0.265876 3555.677943
6
              2900.380412
    0.618193
7
    0.418159
              3696.281878
8
    0.435152
              2282.279598
    0.434370 4122.939981
10 0.685685 1769.684057
R^2
          0.461962
MSE
       3000.381033
dtype: float64
```

# Question 1: Linear Regression, Ridge Regression Lasso (20 points)

Now it's your turn to perform a linear regression experiment! Use the <u>Boston Housing Prices Da</u> (<a href="http://scikit-">http://scikit-</a>

<u>learn.org/stable/modules/generated/sklearn.datasets.load\_boston.html#sklearn.datasets.load\_t</u> built into scikit-learn to better understand regression. You can load it with datasets.load\_bo and use the .DESCR field to learn more about the dataset.

- 1. Perform exploratory data analysis on the Boston Housing data.
  - From your exploratory data analysis, what do you notice about the attribute values?
  - Which parameter to the LinearRegression model could you use to deal with this issue?
  - Perform 10-fold cross-validation (with shuffle=True and random\_state=2016020 with a LinearRegression model that uses the parameter identified above. Report the average coefficient of determination and mean squared error metric on the test set (averaged across folds).
- 2. Two very popular options in Linear Regression are the <u>Lasso method</u> (<a href="https://en.wikipedia.org/wiki/Lasso">https://en.wikipedia.org/wiki/Lasso</a> (statistics)) and <u>Ridge Regression</u> (<a href="https://en.wikipedia.org/wiki/Tikhonov regularization">https://en.wikipedia.org/wiki/Tikhonov regularization</a>). These are implemented in the <u>Lasso (http://scikit-</u>

learn.org/stable/modules/generated/sklearn.linear model.Lasso.html#sklearn.linear model.

and Ridge (http://scikit-

<u>learn.org/stable/modules/generated/sklearn.linear\_model.Ridge.html#sklearn.linear\_model.</u> models in sklearn.

- What one word captures what Lasso and Ridge do? (It's in the first sentence of both Wikipedia articles and sklearn documentation pages)
- Perform 10-fold CV with normal LinearRegression, Lasso, and Ridge. Compare the attribute weights (coef) of each of these methods. What do you observe?
- Report the average R^2 and MSE for each method on the test set across folds. How d these variants of linear regression perform?

Hint: If you'd like to understand Ridge Regression better, you might want to look at this <u>exa</u> (http://scikit-learn.org/stable/auto\_examples/linear\_model/plot\_ols\_ridge\_variance.html)

# **Answers 1:**

## 1.

- a) From the exploratory data analysis, we noticed that the range of the values is huge, which is not good. For example, most of the attributes have a value below 100, while the 9th and 11th attribute have values in the hundreds. Betewen 300 and 700 for the 9th attribute [TAX] and around 400 for the 11th attribute [B]
- **b)** To deal with this issue, we'll use the normalize=True parameter to the LinearRegression model
- c) Averages:

R^2 0.713643

MSE 23.657502

## 2.

- **a)** Regularization. Both Lasso and Ridge regression are used for regularization to prevent overfitting
- **b)** Looking at the attribute weights, we see several differences between the different models. First, Lasso gets completly rid of most features, by settings the weights of 0. It only keeps 2 or 3 features in most cases as the only relevant once to predict the median house price. The Ridge regression model on the other hand, keeps most features, but we can notice that the weights are slightly different to the weights of the original linear regresion (i.e there are only small changes to the weights).

## c)

Averages:

Linear Regression:

R^2 0.713643

MSE 23.657502

Lasso Regularization: R^2 0.580869 MSE 35.198977

Ridge Regression: R^2 0.708486 MSE 24.133678

TTest for LinearRegression vs RidgeRegression
Ttest\_relResult(statistic=1.2154248843325626, pvalue=0.25512413165702003)

TTest for Lasso vs RidgeRegression
Ttest\_relResult(statistic=-9.4005571317290304, pvalue=5.9711873467613727e-06)

TTest for LinearRegression vs Lasso Ttest\_relResult(statistic=8.7799598190605064, pvalue=1.0449626130453215e-05)

The Lasso regression model performs significantly worse than the linear regression and the ridge regression model. This means that the model is not overfitted, and most features are relevant to predicting the median house prices. So when Lasso gets rid of some features, the model performs worse.

The Ridge regression performs slightly worse than the regular linear regression, but according to the ttest, this difference is not significant.

```
In [19]: # Load the data
boston = datasets.load_boston();

# Brief description of the data
print boston.DESCR
```

Boston House Prices dataset

#### Notes

\_\_\_\_\_

Data Set Characteristics:

:Number of Instances: 506

:Number of Attributes: 13 numeric/categorical predictive

:Median Value (attribute 14) is usually the target

:Attribute Information (in order):

- CRIM per capita crime rate by town
- ZN proportion of residential land zoned for lots ove r 25,000 sq.ft.
  - INDUS proportion of non-retail business acres per town
- CHAS Charles River dummy variable (= 1 if tract bounds river; 0 otherwise)
- NOX nitric oxides concentration (parts per 10 millio n)
  - RM average number of rooms per dwelling
  - AGE proportion of owner-occupied units built prior to

1940

- DIS weighted distances to five Boston employment cent

res

- RAD index of accessibility to radial highways
- TAX full-value property-tax rate per \$10,000
- PTRATIO pupil-teacher ratio by town
- -B 1000(Bk 0.63)^2 where Bk is the proportion of b

lacks by town

- LSTAT % lower status of the population
- MEDV Median value of owner-occupied homes in \$1000's

:Missing Attribute Values: None

:Creator: Harrison, D. and Rubinfeld, D.L.

This is a copy of UCI ML housing dataset.

http://archive.ics.uci.edu/ml/datasets/Housing (http://archive.ics.u ci.edu/ml/datasets/Housing)

This dataset was taken from the StatLib library which is maintained at Carnegie Mellon University.

The Boston house-price data of Harrison, D. and Rubinfeld, D.L. 'Hed

prices and the demand for clean air', J. Environ. Economics & Manage ment,

vol.5, 81-102, 1978. Used in Belsley, Kuh & Welsch, 'Regression di agnostics

...', Wiley, 1980. N.B. Various transformations are used in the table on

pages 244-261 of the latter.

The Boston house-price data has been used in many machine learning p apers that address regression problems.

### \*\*References\*\*

- Belsley, Kuh & Welsch, 'Regression diagnostics: Identifying Influential Data and Sources of Collinearity', Wiley, 1980. 244-261.
- Quinlan,R. (1993). Combining Instance-Based and Model-Based Lea rning. In Proceedings on the Tenth International Conference of Machine Learning, 236-243, University of Massachusetts, Amherst. Morgan K aufmann.
- many more! (see http://archive.ics.uci.edu/ml/datasets/Housing)
  (http://archive.ics.uci.edu/ml/datasets/Housing))

```
In [20]: # Part la) Explore the dataset
    # Put it into pandas DataFrames
    boston_data_df = pd.DataFrame(boston.data);
    boston_target_df = pd.DataFrame(boston.target)

# How many attributes and records are there?
print boston_data_df.shape

# What are the descriptive statistics?
print boston_data_df.describe()

# How are the labels distributed?
boston_target_df.hist()

# How are the attributes distributed?
plt.figure()
axes = boston_data_df.boxplot()
```

(506, 13)					
0	1	2	3	4	
5 \					
count 506.000000	506.000000	506.000000	506.000000	506.000000	5
06.000000					
mean 3.593761	11.363636	11.136779	0.069170	0.554695	
6.284634 std 8.596783	22 222452	6.860353	0 252004	0 115070	
std 8.596783 0.702617	23.322453	0.800333	0.253994	0.115878	
min 0.006320	0.00000	0.460000	0.000000	0.385000	
3.561000	0.00000	0.400000	0.000000	0.303000	
25% 0.082045	0.000000	5.190000	0.000000	0.449000	
5.885500					
50% 0.256510	0.000000	9.690000	0.000000	0.538000	
6.208500					
75% 3.647423	12.500000	18.100000	0.000000	0.624000	
6.623500					
max 88.976200	100.000000	27.740000	1.000000	0.871000	
8.780000					
	_				
6	7	8	9	10	
11 \ count 506.000000	506.000000	506.000000	506.000000	506.000000	5
count 506.000000 06.000000	306.000000	300.000000	300.000000	300.000000	5
mean 68.574901	3.795043	9.549407	408.237154	18.455534	3
56.674032	00,75010	30313107	100020,131	101133301	
std 28.148861	2.105710	8.707259	168.537116	2.164946	
91.294864					
min 2.900000	1.129600	1.000000	187.000000	12.600000	
0.320000					
25% 45.025000	2.100175	4.000000	279.000000	17.400000	3
75.377500					_
50% 77.500000	3.207450	5.000000	330.000000	19.050000	3
91.440000	E 100/2E	24 000000	666 000000	20 200000	2
75% 94.075000 96.225000	5.188425	24.000000	666.000000	20.200000	3
max 100.000000	12.126500	24.000000	711.000000	22.000000	3
96.900000	12.120500	24.000000	711.000000	22.000000	J
12					
count 506.000000					
mean 12.653063					
std 7.141062					
min 1.730000					
25% 6.950000					
50% 11.360000					
75% 16.955000					
max 37.970000					

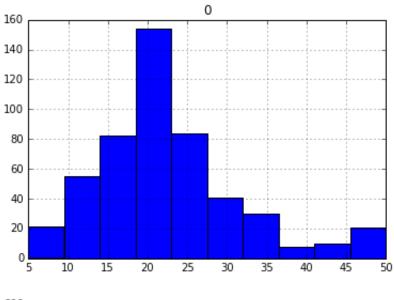
<sup>/</sup>Users/viktorjankov/anaconda/lib/python2.7/site-packages/ipykernel/\_ \_main\_\_.py:17: FutureWarning:

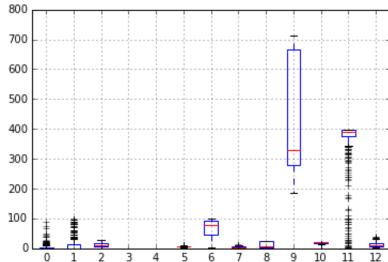
The default value for 'return\_type' will change to 'axes' in a futur e release.

To use the future behavior now, set return\_type='axes'.

ma been the measurest behavior and dilense this receive and return

To keep the previous behavior and silence this warning, set return\_type='dict'.





```
In [21]: # Part 1b) Cross Validation
         foldnum = 0
         fold results = pd.DataFrame()
         for train, test in cross validation. KFold(len(boston.data), n folds=10,
             foldnum+=1
             [boston_tr_data, boston_te_data,
              boston tr target, boston te target] = folds to split(boston.data,b
             lr = linear model.LinearRegression(normalize=True)
             lr.fit(boston tr data, boston tr target)
             # But a nicer way to store them is in a DataFrame
             fold results.loc[foldnum, 'R^2'] = lr.score(boston_te_data, boston_
             fold results.loc[foldnum, 'MSE'] = metrics.mean squared error(bosto
         #Now let's look at the results:
         print fold results
         #And compute the mean error across folds:
         print fold results.mean()
```

```
R^2
                    MSE
1
    0.717454 23.926098
2
    0.620171 34.779133
3
    0.783767 18.841212
4
    0.556936 30.379589
5
    0.769709 16.603228
    0.693492 22.287915
6
7
   0.828865 19.256283
    0.631266 33.237697
9
    0.796774 16.961117
10 0.737993 20.302751
R^2
        0.713643
MSE
       23.657502
dtype: float64
```

In [22]:		

```
# Part 2 Using Lasso and Ridge Regression
from scipy import stats
foldnum = 0
lr fold results = pd.DataFrame()
lasso fold results = pd.DataFrame()
ridge fold results = pd.DataFrame()
for train, test in cross validation. KFold(len(boston.data), n folds=10,
    foldnum+=1
    [boston_tr_data, boston te data,
    boston tr target, boston te target] = folds to split(boston.data,b
    lr = linear model.LinearRegression(normalize=True)
    lr.fit(boston tr data, boston tr target)
    lasso = linear model.Lasso(alpha=0.1, normalize=True, random state=
    lasso.fit(boston tr data, boston tr target)
    ridge = linear model.Ridge(alpha=0.1, normalize=True, random state=
    ridge.fit(boston tr data, boston tr target)
    # But a nicer way to store them is in a DataFrame
    lr fold results.loc[foldnum, 'R^2'] = lr.score(boston te data, bost
    lr fold results.loc[foldnum, 'MSE'] = metrics.mean squared error(bo
    lasso fold results.loc[foldnum, 'R^2'] = lasso.score(boston te data
    lasso fold results.loc[foldnum, 'MSE'] = metrics.mean squared error
    ridge_fold_results.loc[foldnum, 'R^2'] = ridge.score(boston_te_data
    ridge fold results.loc[foldnum, 'MSE'] = metrics.mean_squared_error
#Now let's look at the results:
print "TTest for LinearRegression vs RidgeRegression"
print stats.ttest rel(lr fold results['R^2'].values, ridge fold results
print "\n"
print "TTest for Lasso vs RidgeRegression"
print stats.ttest rel(lasso fold results['R^2'].values, ridge fold resu
print "\n"
print "TTest for LinearRegression vs Lasso"
print stats.ttest rel(lr fold results['R^2'].values, lasso fold results
print "\n"
#And compute the mean error across folds:
print "Linear Regression:"
print lr fold results.mean()
print "\n"
print "Lasso Regularization:"
print lasso fold results.mean()
```

```
print "\n"
print "Ridge Regression: "
print ridge fold results.mean()
TTest for LinearRegression vs RidgeRegression
Ttest relResult(statistic=1.2154248843325628, pvalue=0.2551241316570
2003)
TTest for Lasso vs RidgeRegression
Ttest relResult(statistic=-9.4005571317290268, pvalue=5.971187346761
393e-06)
TTest for LinearRegression vs Lasso
Ttest relResult(statistic=8.7799598190605046, pvalue=1.0449626130453
218e-05)
Linear Regression:
R^2
        0.713643
MSE
       23.657502
dtype: float64
Lasso Regularization:
R^2
        0.580869
MSE
       35.198977
dtype: float64
Ridge Regression:
R^2
        0.708486
MSE
       24.133678
dtype: float64
```

# **Logistic Regression**

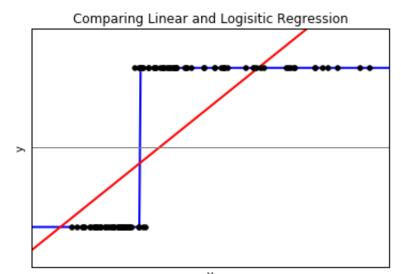
Now let's look at a different application for regression algorithms: classification. As we learned in class, regression algorithm can be adapted to a binary classification fairly simply. Instead of predicting some continuous value, we predict the class. So, for example, instead of predicting the diabetes condition of a patient we would predict whether or not a patient has diabetes by having the label be 1 when the patient has diabetes and 0 when the patient does not have diabetes.

In this setting the regression algorithm tries to learn some combination of features to get 0 for the negative instance and 1 for the positive instances. Of course, linear regression algorithms produce continuous values, not just 1s and 0s, so we have to convert the

continuous values into 1s and 0s. Usually this is done by rounding values above 0.5 to 1 and the remaining values to 0 (using, for example, the numpy <u>round</u> (<a href="http://docs.scipy.org/doc/numpy-1.10.0/reference/generated/numpy.round">http://docs.scipy.org/doc/numpy-1.10.0/reference/generated/numpy.round</a> .html) function).

However, there's still something messy about this setting. Linear regression wants to predict continuous values and we're conscripting it to produce 1s and 0s. What if we had a regression algorithm that, by design, predicted 1s and 0s? This is exactly what logistic regression does. Let's look at an example of the difference, taken from the <a href="scikit-learn-documentation">scikit-learn-documentation</a> (http://scikit-learn.org/stable/auto examples/linear model/plot logistic.html)

```
In [23]: # Code source: Gael Varoquaux
         # License: BSD 3 clause
         # this is our test set, it's just a straight line with some
         # Gaussian noise
         xmin, xmax = -5, 5
         n \text{ samples} = 100
         np.random.seed(0)
         X = np.random.normal(size=n samples)
         y = (X > 0).astype(np.float)
         X[X > 0] *= 4
         X += .3 * np.random.normal(size=n samples)
         X = X[:, np.newaxis]
         #Learn a Linear Regression model
         lr = linear model.LinearRegression()
         lr.fit(X, y)
         # Learn a Logistic Regression model
         lgr = linear model.LogisticRegression(C=1e5)
         lgr.fit(X, y)
         # Plot the training data (as black dots)
         plt.scatter(X.ravel(), y, color='black', zorder=20)
         #Create some test data
         X \text{ test} = \text{np.linspace}(-5, 10, 300).\text{reshape}(300,1)
         # Plot the predictions of the logistic regression model (blue line)
         plt.plot(X test, lgr.predict(X test), color='blue', linewidth=2)
         # Plot the prediction of the linear regresion model (red line)
         plt.plot(X test, lr.predict(X_test), color='red', linewidth=2)
         # Make the plot prettier
         plt.axhline(.5, color='.5')
         plt.ylabel('y')
         plt.xlabel('X')
         plt.xticks(())
         plt.yticks(())
         plt.ylim(-.25, 1.25)
         plt.xlim(-4, 10)
         plt.title('Comparing Linear and Logisitic Regression')
         plt.show()
```



The black dots in the plot above are the training data. The linear regression model fits that training data with the line shown in red. The logistic regression model fits the same training data with the line shown in blue. As you can see, the logistic regression model is a much closer fit to the binary training data than the linear regression model.

Now let's explore logistic regression with a different data set. The <u>Adult dataset</u> (<a href="http://archive.ics.uci.edu/ml/datasets/Adult">http://archive.ics.uci.edu/ml/datasets/Adult</a>) consists of 14 attributes from census data and tries to predict whether an individual has an income exceeding \$50K. I've covered some of the basic steps to get the data loaded, but you can explore the original data census\_orig (which has readable, nominal attributes) and the transformed data census\_data and census\_labels to get a better understanding of the dataset.

```
census data = urllib2.urlopen("http://archive.ics.uci.edu/ml/machine-le
In [24]:
         census_orig = pd.read_csv(census_data, quotechar='"', skipinitialspace=
                                         names=['Age','WorkClass', 'FnlWgt', 'Edu
                                                'Occupation', 'Relationship', 'Ra
                                                'CapitalGain', 'CapitalLoss', 'Ho
                                                'NativeCountry', 'Label'],
                                        na values="?", index col=False)
         census_orig = census_orig.dropna()
         # Convert labels from strings to boolean
         label encoder = preprocessing.LabelEncoder()
         census labels = pd.DataFrame(label encoder.fit transform(census orig.il
         # Convert nominal attributes to encoded versions
         attr encoder = feature extraction.DictVectorizer(sparse=False)
         census data = pd.DataFrame(attr encoder.fit transform(census orig.iloc[
         census data.columns = attr encoder.get feature names()
```

# Question 2: Logistic Regression (20 points)

- 1. Learn linear and logistic regression classifiers for the census data. Perform 10-fold cross-validation (with shuffle=True and random\_state=20160202) and report the mean accuracy of each classifier. Remember, you have to transform the linear regression result to get a classification.
- 2. One of the important aspects data mining is data normalization. Use the StandardScaler to normalize your data. Now re-run the same 10-fold experiment for both classifiers. What changed?
- 3. The Lasso and Ridge methods above are also part of logistic regression, although you must specify them slightly differently. You can specify the parameter penalty='11' or penalty='12' to specify Lasso or Ridge, respectively. The importance of regularization can be specified using the parameter C. Perform 10-fold CV with the L1 and L2 penalties and C=[0.1, 10]. Report the results of each combination (for a total of 4) averaged across folds.

Note: if this were a real experiment, you would want to do your parameter search with a validation set.

# **Answers 2:**

## 1.

Linear Regression Mean Accuracy 0.833996 Logistic Regression Accuracy 0.790465

## 2.

The accuracies changed between the two models. Now with the data normalized, the Logistic Regression performs better than the Linear Regression, whereas in the original, non-standardized data, the reverse was true. It's worthy to note that the accuracy of the Logistic Regression improved significantly ~5%

Linear Regression Mean Accuracy 0.833864 Logistic Regression Accuracy 0.848551

### 3.

Lasso Log Regression L1 0.1: 0.848087 L1 10: 0.848551

Ridge Log Regression:

L2 0.1: 0.848485 L2 10: 0.848551

```
In [25]: # Part 1 no normalization
         from sklearn.metrics import accuracy score
         #maybe do your EDA here?
         foldnum = 0
         lr fold results = pd.DataFrame()
         lgr fold results = pd.DataFrame()
         for train, test in cross validation. KFold(len(census data), n folds=10,
             foldnum+=1
             [adult tr data, adult te data,
              adult tr target, adult te target] = folds to split(census data,cen
             lr = linear model.LinearRegression()
             lr.fit(adult tr data, adult tr target)
             prediction = lr.predict(adult te data)
             classified prediction = []
             for value in range(len(prediction)):
                 classified prediction.append(round(prediction[value]))
             adult te classified data = pd.DataFrame(classified prediction)
             lgr = linear model.LogisticRegression(C=1e5)
             lgr.fit(adult tr data, np.reshape(adult tr target.values,[len(adult
             # But a nicer way to store them is in a DataFrame
             lr fold results.loc[foldnum, 'LinearReg Mean Accuracy'] = accuracy
             lgr fold results.loc[foldnum, 'LogRegr Accuracy'] = lgr.score(adult
         #Now let's look at the results:
         print lr fold results
         #And compute the mean error across folds:
         print lr fold results.mean()
         print lgr_fold_results
         print lgr fold results.mean()
```

```
LinearReg Mean Accuracy
1
                    0.840570
2
                    0.833941
3
                    0.834881
4
                    0.839523
5
                    0.830902
6
                    0.831565
7
                    0.834218
8
                    0.827586
9
                    0.832560
10
                    0.834218
LinearReg Mean Accuracy
                             0.833996
dtype: float64
    LogRegr Accuracy
1
             0.783560
2
             0.790520
3
             0.797082
4
             0.802056
5
             0.781167
6
             0.790119
7
             0.793103
8
             0.783820
9
             0.784483
             0.798740
10
LogRegr Accuracy
                     0.790465
```

dtype: float64

```
In [26]: # Part 2 normalizing the data
         from sklearn import preprocessing
         sscaler = preprocessing.StandardScaler()
         census data norm = pd.DataFrame(sscaler.fit transform(census data.value
         #maybe do your EDA here?
         foldnum = 0
         lr fold results = pd.DataFrame()
         lgr fold results = pd.DataFrame()
         for train, test in cross validation. KFold(len(census data norm), n fold
             foldnum+=1
             [adult tr data, adult te data,
              adult tr target, adult te target] = folds to split(census data nor
             lr = linear model.LinearRegression()
             lr.fit(adult_tr_data, adult_tr_target)
             prediction = lr.predict(adult te data)
             classified prediction = []
             for value in range(len(prediction)):
                 classified prediction.append(round(prediction[value]))
             adult te classified data = pd.DataFrame(classified prediction)
             lgr = linear model.LogisticRegression(C=1e5)
             lgr.fit(adult tr data, np.reshape(adult tr target.values,[len(adult
             # But a nicer way to store them is in a DataFrame
             lr fold results.loc[foldnum, 'LinearReg Mean Accuracy'] = accuracy
             lgr fold results.loc[foldnum, 'LogRegr Accuracy'] = lgr.score(adult
         #Now let's look at the results:
         print lr fold results
         #And compute the mean error across folds:
         print lr fold results.mean()
         print lgr fold results
         print lgr fold results.mean()
```

```
LinearReg Mean Accuracy
1
                    0.842559
2
                    0.832615
3
                    0.835212
4
                    0.839191
5
                    0.830570
6
                    0.831233
7
                    0.832560
8
                    0.826260
9
                    0.832891
10
                    0.835544
LinearReg Mean Accuracy
                            0.833864
dtype: float64
    LogRegr Accuracy
1
            0.857474
2
            0.846536
3
            0.855106
4
             0.852122
5
            0.842838
6
            0.845159
7
            0.852122
8
            0.838196
9
            0.848806
10
            0.847149
LogRegr Accuracy
                     0.848551
dtype: float64
```

```
In [27]: # Part 3 Lasso and Ridge using Logistic Regression
         foldnum = 0
         fold 11 01 results = pd.DataFrame()
         fold 11 10 results = pd.DataFrame()
         fold 12 01 results = pd.DataFrame()
         fold 12 10 results = pd.DataFrame()
         for train, test in cross validation.KFold(len(census_data_norm), n_fold
             foldnum+=1
             [adult_tr_data, adult_te_data,
              adult tr target, adult te_target] = folds_to_split(census_data_nor
             lgr 11 01 = linear model.LogisticRegression(penalty = '11', C=0.1)
             lgr_l1_10 = linear_model.LogisticRegression(penalty = 'l1', C=10)
             lgr 12 01 = linear model.LogisticRegression(penalty = '12', C=0.1)
             lgr 12 10 = linear model.LogisticRegression(penalty = '12', C=10)
             lgr 11 01.fit(adult tr data, np.reshape(adult tr target.values,[len
             lgr 11 10.fit(adult tr data, np.reshape(adult tr target.values,[len
             lgr 12 01.fit(adult tr data, np.reshape(adult tr target.values,[len
             lgr 12 10.fit(adult tr data, np.reshape(adult tr target.values,[len
             # But a nicer way to store them is in a DataFrame
             fold 11 01 results.loc[foldnum, 'L1 0.1: '] = lgr 11 01.score(adult
             fold_l1_10_results.loc[foldnum, 'L1 10: '] = lgr_l1_10.score(adult_
             fold_12_01_results.loc[foldnum, 'L2 0.1: '] = lgr_12_01.score(adult
             fold 12 10 results.loc[foldnum, 'L2 10: '] = lgr_l2_10.score(adult_
         #And compute the mean error across folds:
         print fold 11 01 results.mean()
         print fold 11 10 results.mean()
         print fold 12 01 results.mean()
         print fold 12 10 results.mean()
         L1 0.1:
                     0.848087
         dtype: float64
         L1 10:
                    0.848551
         dtype: float64
         L2 0.1:
                     0.848485
         dtype: float64
         L2 10:
                    0.848551
```

dtype: float64