Linear Regression - Ridge Regression

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Project Guide

- Project Overview
- Data Review
- Coding Ridge Regression
- Ridge Regression in sklearn

Project Overview

EXPECTED TIME 2.5 HRS

This assignment will test your ability to code your own version of ridge-regularized regression in Python. This assignment draws upon and presupposed the knowledge found in the lectures for Week 2. If ever a theoretical questions arises as to "why" we are doing something, please refer back to those lectures.

The assignment also builds upon the work performed in assignment 1 "Linear Regression - Least Squares". The data used will be the same. Though the last assignment tested your ability to read data into Pandas from a .csv . Those fundamental processes will not be directly tested here.

In coding Ridge Regression you will be asked to:

- Mean center target variable and mean center / standardize observation
- Calculate Ridge Regression w eights using linear algebra
- Create a hyperparameter tuning process

Motivation: Ridge Regression offers a way to mitigate some of the weaknesses of Least Squares Linear Regression to build more robust models.

Objectives: This assignmet will -

- Test Python competency and mathematical understandings of Ridge Regression
- Begin to introduce the concept of hyper-parameter tuning

Problem: Using housing data, we will attempt to predict house price using living area with a regression model.

Data: Data comes from Kaggle's House Prices Dataset.

See above link for Description of data

Data Exploration

Below provides a review of the "Housing" dataset.

```
### This cell imports the necessary modules and sets a few plotting parameters for display

Mmatplotlib inline
import numpy as np
import pandas as pd
import matplotlib.pyplot as plt
plt.ncParams['figure.figsize'] = (20.0, 10.0)

### Read in the data
tr_path = '../resource/asnlib/publicdata/train.csv'
test_path = '../resource/asnlib/publicdata/test.csv'
data = pd.read_csv(tr_path)

### The .head() function shows the first few lines of data for perspective
data.head()
```

Question 1

```
### GRADED
### Which column has the most "null" values? assign name as string to ans1.
### ### CAPITALIZATION/SPELLING MATTERS e.g. 'Street' != 'street'
### How many nulls are in that column? assign number as int to ans2
### YOUR ANSWER BELOW

ans1 = 'PoolQC'
ans2 = 1453
```

```
data.plot('GrLivArea', 'SalePrice', kind = 'scatter', marker = 'x');
```

```
### GRADED
### In building regressions below, a subset of our data will be used.
### Practice subsetting a DataFrame below.
### Create a DataFrame only containing the "Street" and "Alley" columns from
### the 'data' DataFrame.
```

```
### Assign to 'ans1'
### YOUR ANSWER BELOW

cols = ['Street','Alley']
ans1 = data[cols]
```

Coding Ridge Regression

Preprocessing

Before implementing Ridge Regression, it is important to mean-center our target variable and mean-center and standardize observations. We will do this according to the following:

Mean Center Target

```
\sy_{cent} = y_0 - \propty
```

Standardize Observations

```
X_{std} = \frac{X_0-\bar{X}}{s_{X}}
```

Where \bar{X} is the sample mean of X and s_{X} is the sample standard deviation of X

Question 3:

```
### GRADED
### Why are the centering / standardization transformations described above important for ridge regression?
### 'a') Regression works best when values are unitless
### 'b') The transformations makes the regression more interpretable
### 'c') Ridge penalizes large coefficients; the transformations make the coefficients of similar scales
### 'd') It isn't important
### Assign character associated with your choice as a string to ans1
### YOUR ANSWER BELOW
ans1 = 'C'
```

Question 4:

 $\SX_{std} = \frac{X_0-\bar{X}}{s_{X}}$

```
### CADED
### Code a function called "standardize"
### ACCEPT one input, a list of numbers
### To standardize, subtract the mean of the list and divide by standard deviation.
### YOUR ANSWER BELOW

def standardize (num_list):
    """

Standardize the given list of numbers

Positional arguments:
    num_list = [1,2,3,3,4,4,5,5,5,5]
    nl_std = standardize(num_list)
    print(np-round(nl_std2))
#--> np.array([-2.11, -1.36, -0.61, -0.61, -0.61, -0.14, -0.14, -0.88, 0.88, 0.88]

### Calculate standard deviation and mean
std * np.std(num_list)

## Inplement equation
toRet = [(x-mean)/std for x in num_list]

return toRet
```

Below $\ w \ e \ w \ ill \ create \ a \ function \ w \ hich \ w \ ill \ preprocess \ our \ data \ by \ performing:$

- mean subtraction from \$y\$,
- dimension standardization for \$x\$.

Both according to the equations set out below.

Mean Center Target

```
$$y_{cent} = y_0 - \bar{y}$$
```

Standardize Observations

 $X_{std} = \frac{X_0-\bar{X}}{s_{X}}$

```
def preprocess_for_regularization(data, y_column_name, x_column_names):
      Perform mean subtraction and dimension standardization on data
      Positional argument:
           data -- a pandas dataframe of the data to pre-process y_column_name -- the name (string) of the column that contains
                 the target of the training data.
olumn_names -- a *list* of the names of columns that contain the
           x_column_names -- a *list* of the na
observations to be standardized
      Returns:
            Return a DataFrame consisting only of the columns included
            in 'y_column_name' and 'x_column_names'. Where the y_column has been mean-centered, and the x_columns have been mean-centered/standardized.
            data = pd.read_csv(tr_path).head()
            prepro_data = preprocess_for_regularization(data,'SalePrice', ['GrLivArea','YearBuilt'])
           print(prepro_data) #-->
                     repro_data) #-->
Griivfrea YearBuilt SalePrice
0 -0.082772 0.716753 7800.0
1 -1.590161 -0.089594 -19200.0
2 0.172946 0.657024 22800.0
3 -0.6959219 -1.911342 -60700.0
4 1.559205 0.627159 49300.0
      # Create list of all columns
      toRetCol = x_column_names + [y_column_name]
      # subset datafram
      toRet = data[toRetCol].copy() # For "setting with copy" warning
      # calculate mean of y, then subtract from all y's
      y_mean = np.mean(data[y_column_name])
toRet[y_column_name] = data[y_column_name].apply(lambda x: x - y_mean)
      # Calc mean and std for every column in x # Then apply standardization
      for column in x_column_names:
           mean = np.mean(data[column])
std = np.std(data[column])
           {\tt toRet[column] = data[column].apply(lambda x: (x - mean)/std)}
     return toRet
```

Next, you'll implement the equation for ridge regression using the closed form equation:

```
\ \ _{RR}=(\lambda+X^TX)^{-1}X^Ty$$
```

The function will be very similar to the function you wrote for Least Squares Regression with a slightly different matrix to invert.

 $NB: \textit{Many} \ \ \textit{numpy} \ \ \textit{matrix} \ \ \textit{functions} \ \ \textit{will} \ \ \textit{be} \ \ \textit{useful}. \ \textit{e.g.} \ \ \textit{np.matmul} \ , \ \ \textit{np.linalg.inv} \ , \ \ \textit{np.ones} \ , \ \ \textit{np.transpose} \ , \ \textit{and} \ \ \textit{np.identity} \ .$

The main change from Least Squares Regression is that \$\ambda\$ is a parameter we must set. This is different from the \$w\$ parameters that we calculate from either closed form or approximation algorithms.

We will address tuning parameters such as \$\lambda\$ in the next section.

```
### GRADED
### Code a function called "ridge_regression_weights" ### ACCEPT three inputs:
### Two matricies corresponding to the x inputs and y target ### and a number (int or float) for the lambda parameter \,
### RETURN a numpy array of regression weights
### The following must be accomplished:
### Ensure the number of rows of each the X matrix is greater than the number of columns.
### ### If not, transpose the matrix.
### Ultimately, the y input will have length n.
### Thus the x input should be in the shape n-by-p
### *Prepend* an n-by-1 column of ones to the input x matrix
### Use the above equation to calculate the least squares weights.
### This will involve creating the lambda matrix---
### ### a p+1-by-p+1 matrix with the "lambda_param" on the diagonal
### ### p+1-by-p+1 because of the prepended "ones".
### NB: Pay close attention to the expected format of the returned ### weights. It is different / simplified from Assignment 1.
### YOUR ANSWER BELOW
def ridge_regression_weights(input_x, output_y, lambda_param):
          Calculate ridge regression least squares weights
     Positional arguments:
           input x - 2-d matrix of input data
output_y -- 1-d numpy array of target values
lambda_param -- lambda parameter that controls how heavily
                to penalize large weight values
     Example:
           training_y = np.array([208500, 181500, 223500,
```

```
140000, 250000, 143000,
                                   307000, 200000, 129900
                                   118000])
     training x = np.array([[1710, 1262, 1786,
                                  1717, 2198, 1362,
1694, 2090, 1774,
                                   1077],
                                 [2003, 1976, 2001,
1915, 2000, 1993,
2004, 1973, 1931,
     lambda_param = 10
     rrw = ridge_regression_weights(training_x, training_y, lambda_param)
     print(rrw) #--> np.array([-576.67947107, 77.45913349, 31.50189177])
     print(rrw[2]) #--> 31.50189177
Assumptions:
     -- output_y is a vector whose length is the same as the number of observations in input_x
-- lambda_param has a value greater than 0
# Check to ensure dataframe is long not wide
if input_x.shape[0] < input_x.shape[1]:
   input_x = np.transpose(input_x)</pre>
# Create column of ones
       = np.ones((len(output_y), 1), dtype=int)
# Add column of ones to X
  ugmented_x = np.concatenate((ones, input_x), axis=1)
# Create square lambda_matrix, with size equal to number of columns in X
lambda_matrix = lambda_param * np.identity(min(augmented_x.shape))
\# Invert lambda + dot-prod of x and transposed x
inv = np.linalg.inv(lambda_matrix + np.matmul(np.transpose(augmented_x), augmented_x))
# dot-prod of inverted matrix and transposed X
left_multiplier = np.matmul(inv , np.transpose(augmented_x))
# final dot-prod with the ys
weights = np.matmul(left_multiplier, output_y)
return weights
```

Selecting the \$\lambda\$ parameter

For our final function before looking at the sklearn implementation of ridge regression, we will create a hyper-parameter tuning algorithm.

In ridge regression, we must pick a value for \$\lambda\$. We have some intuition about \$\lambda\$ from the equations that define it: small values tend to emulate the results from Least Squares, while large values will reduce the dimensionality of the problem. But the choice of \$\lambda\$ can motivated with a more precise quantitative treatment.

Eventually, we will look to choose the value of \$\ambda\$ that minimizes validation error, which we will determine using \$k\$-fold cross-validation.

For this example here, we will solve a simpler but more problem: Find a value that minimizes the of the list returned by a function.

```
### Example of hiden function below:
### 'hidden' takes a single number as a parameter (int or float) and returns a list of 1000 numbers
### the input must be between 0 and 50 exclusive

def hidden(hp):
    if (hp<=0) or (hp >= 50):
        print("input out of bounds")

nums = np.logspace(0,5,num = 1000)
    vals = nums** 43.123985172351235134687934

user_vals = nums** hp
return vals-user_vals
```

Run the above cell and test out the functionality of hidden . Remember it takes a single number, between 0 and 50, as an argument

```
### GRADED
### Code a function called "minimize"
### That function will be similar to 'hidden' created above and available for your exploration.
### That function will be similar to 'hidden' created above and available for your exploration.
### Like 'hidden', the passed function will take a single argument, a number between 0 and 50 exclusive
### and then, the function will return a numpy array of 1000 numbers.

### RETURN the value that makes the mean of the array returned by 'passed_func' as close to 0 as possible

### Note, you will almost certainly NOT be able to find the number that makes the mean exactly 0

### YOUR ANSWER BELOW

def minimize( passed_func):
    """
    Find the numeric value that minimizes the output of 'passed_func'

Positional Argument:
    passed_func -- a function that takes a single number (between 0 and 50 exclusive)
    as input, and returns a list of 1000 floats.

Example:
```

```
passed_func = hidden
min_hidden = minimize(passed_func)
print(round(min_hidden,4))
#--> 43.1204 (answers will vary, must be close to 43.123985172351)

"""

# Create values to test
test_vals = list(np.linspace(.1,49.9, 1000))

# Find mean of returned array from function
ret_vals = [abs(np.mean(passed_func(x))) for x in test_vals]

# Find smallest mean
min_mean = min(ret_vals)

# Return the test value that creates the smallest mean
return test_vals[ret_vals.index(min_mean)]
```

The above simulates hyper parameter tuning.

In the case of ridge regression, you would be searching lambda parameters to minimize validation error.

The hidden function would be analogous to the model building; the returned list analogous to residuals; and the mean of that list analogous to validation error.

See below for an example of using the functions built above that automatically performs hyper-parameter tuning using mean-absolute-deviation.

```
def lambda_search_func(lambda_param):

# Define X and y
# with preprocessing
df = preprocess_for_regularization(data.head(50), 'SalePrice', ['GrLivArea', 'YearBuilt'])

y_true = df['SalePrice'].values
X = df[['GrLivArea', 'YearBuilt']].values

# Calculate Weights then use for predictions
weights = ridge_regression_weights(X, y_true, lambda_param )
y_pred = weights[0] + np.matmul(X,weights[1:])

# Calculate Residuals
resid = y_true - y_pred

# take absolute value to tune on mean-absolute-deviation
# Alternatively, could use:
# return resid **2-5
# for tuning on mean-squared-error
return abs(resid)

minimize(lambda_search_func) # --> about 1.4957957957957957957
```

Implementing a k-folds cross-validation strategy will come in later assignments.

Question 8

```
### GRADED
### Why is cross-validation useful?
### 'a') to minimize the liklihood of overfitting
### 'b') Cross-validation allows us to fit on all our data
### 'b') Cross-validation standardizes outputs
### 'd') cross-validation is not useful
### assing the character associated with your choice as a string to ans1
### YOUR ANSWER BELOW
ans1 = 'a'
```

Ridge Regression in sklearn

Below gives the syntax for implementing ridge regression in sklearn.

```
from sklearn.linear_model import Ridge, LinearRegression
### Note, the "alpha" parameter defines regularization strength.
### Lambda is a reserved word in 'Python' -- Thus "alpha" instead

### An alpha of 0 is equivalent to least-squares regression
Ir = LinearRegression()
reg = Ridge(alpha = 100000)
reg0 = Ridge(alpha = 0)

# Notice how the consistent sklearn syntax may be used to easily fit many kinds of models
for m, name in zip([Ir, reg, reg0], ['LeastSquares", "Ridge alpha = 100000", "Ridge, alpha = 0"]):

m.fit(data[['GrLivArea', 'YearBuilt']], data['SalePrice'])
print(name, "Intercept:", m.intercept_, "Coefs:",m.coef_,"\n")
```

Note, in the above example, an alpha of 100,000 is set for the ridge regularization. The reason an alpha value this high is required is because standardization / mean centering of our inputs did not occur, and instead of w orking with inputs on the order of [-4,4] we are on the interval of [0,2000].

Question 9:

```
### GRADED
### Above, the coefficent around 95/96 corresponds with:
### 'a') Living Area
```

```
### 'b') Year Built
### 'c') Sale Price
### Assign character associated with your choice as string to ans1
### YOUR ANSWER BELOW
ans1 = 'a'
```

Queston 10:

```
### GRADED
### True or False:
### A larger "alpha" corresponds to a greater amount of regularization
### assign boolean choice to ans1
ans1 = True
```