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 over 25,000 sq.ft.
    - INDUS
               proportion of non-retail business acres per town
               Charles River dummy variable (= 1 if tract bounds river; 0 otherwis
    - CHAS
e)
               nitric oxides concentration (parts per 10 million)
    - NOX
               average number of rooms per dwelling
    - RM
               proportion of owner-occupied units built prior to 1940
    - AGE
               weighted distances to five Boston employment centres
    - DIS
    - RAD
               index of accessibility to radial highways
               full-value property-tax rate per $10,000
    - TAX
        - PTRATIO pupil-teacher ratio by town
                   1000(Bk - 0.63)^2 where Bk is the proportion of blacks by town
                   % lower status of the population

    LSTAT

        - 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.uci.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. 'Hedonic prices and the demand for clean air', J. Environ. Economics & Management, vol.5, 81-102, 1978. Used in Belsley, Kuh & Welsch, 'Regression diagnostics ...', 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 papers 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 Learning. In Proceedings on the Tenth International Conference of Machine Learning, 236-243, University of Massachusetts, Amherst. Morgan

Kaufmann.

many more! (see http://archive.ics.uci.edu/ml/datasets/Housing))

In [1]:

```
import warnings #Code to remove warnings
warnings.filterwarnings('ignore')
```

1. Loading Libraries

In [2]:

```
import numpy as np
import pandas as pd
import matplotlib.pyplot as plt
from sklearn.preprocessing import StandardScaler
from sklearn.datasets import load_boston
from sklearn.cross_validation import train_test_split
from sklearn.linear_model import SGDRegressor
from sklearn.model_selection import GridSearchCV
from sklearn.metrics import make_scorer
from sklearn.metrics import mean_squared_error
import prettytable
```

/usr/local/lib/python3.5/site-packages/sklearn/cross_validation.py:41: Depre cationWarning: This module was deprecated in version 0.18 in favor of the mo del_selection module into which all the refactored classes and functions are moved. Also note that the interface of the new CV iterators are different fr om that of this module. This module will be removed in 0.20.

"This module will be removed in 0.20.", DeprecationWarning)

2. Dataset Preprocessing

```
In [3]:
```

```
boston = load_boston() #loading boston housing dataset
```

```
In [4]:
```

```
print('Shape of dataset is:',boston.data.shape)#boston.data gives us an array. Here we are
Shape of dataset is: (506, 13)
```

In [5]:

```
bos = pd.DataFrame(boston.data,columns = boston.feature_names)#Here we are converting our a
#- and our column names are contained in 'boston.feature_names'
bos.head()
```

Out[5]:

	CRIM	ZN	INDUS	CHAS	NOX	RM	AGE	DIS	RAD	TAX	PTRATIO	В	LS
0	0.00632	18.0	2.31	0.0	0.538	6.575	65.2	4.0900	1.0	296.0	15.3	396.90	
1	0.02731	0.0	7.07	0.0	0.469	6.421	78.9	4.9671	2.0	242.0	17.8	396.90	•
2	0.02729	0.0	7.07	0.0	0.469	7.185	61.1	4.9671	2.0	242.0	17.8	392.83	4
3	0.03237	0.0	2.18	0.0	0.458	6.998	45.8	6.0622	3.0	222.0	18.7	394.63	1
4	0.06905	0.0	2.18	0.0	0.458	7.147	54.2	6.0622	3.0	222.0	18.7	396.90	ţ
4													•

In [6]:

bos['Price'] = boston.target #boston.target contains our y_i's and hence we are now putting

In [7]:

```
x = bos.drop('Price',axis = 1)#Here we are splitting our dataset 'bos' into datapoint dataf
y = bos['Price']
```

In [8]:

x_train,x_test,y_train,y_test = train_test_split(x,y)#splitting our dataset into train and

In [9]:

standardized_train = StandardScaler().fit_transform(x_train) #standardizing our training do standardized_test = StandardScaler().fit_transform(x_test) #standardizing our test dataset

In [10]:

```
#Creating training dataframe which contains both datapoints and y_i's
d_train = pd.DataFrame(standardized_train)
d_train['Price'] = np.array(y_train)

#Creating test dataframe which contains both datapoints and y_i's
d_test = pd.DataFrame(standardized_test)
d_test['Price'] = np.array(y_test)
```

In [11]:

```
print(d_train.shape)
print(d_test.shape)
```

(379, 14)

(127, 14)

```
In [12]:
```

```
d_train.head()
```

Out[12]:

	0	1	2	3	4	5	6	7	
0	0.694276	-0.480911	0.998510	-0.248243	0.245557	-1.100563	0.071415	-0.828240	1.6489
1	-0.390557	-0.480911	-0.186256	-0.248243	-0.073535	-0.321752	0.956263	-0.592255	-0.418′
2	-0.048540	-0.480911	1.215253	-0.248243	0.426662	-0.308451	0.870633	-0.721134	-0.5329
3	-0.414197	0.981168	-0.760333	-0.248243	-1.056680	0.414203	-1.006101	0.841536	-0.3033
4	-0.401406	-0.480911	-0.101316	-0.248243	-0.573731	-0.618793	-1.651898	0.088378	-0.6478
4									•

3. Custom implementation of SGDRegressor

```
In [13]:
```

```
def sgdreg fit(x,alpha,batch size=100, n iter = 1000,eta0 = 0.01):
    #Here our learning rate == 'invscaling' and we have used L2 regularization
    """This function is custom implementation of Stochastic Gradient Descent Regressor.
    Input:
          x: This is our dataset on which we want to calculate our weights so that we can f
          alpha: This is our coefficient of L2 Regularizer.
          batch_size: This is the size of our batch of our datapoints which we randomly sam
          n iter: No. of iterations we want to perform.
          eta0: This is our learning rate.
    Output:
           This function return a dictionary containing weight and intercept corresponding
    dict_1 = dict()#dictionary this function returns
    t = 1 #invscaling parameter and is given by t = n iter*batch size
    power_t = 0.25 #invscaling parameter which will be used as power of 't'
    eta = eta0 #learning rate
    w_old = np.zeros(shape = (1,13))#zero intitialization of our weights.
    b_old = 0 #zero initialization of our intercept term
    for i in range(n_iter):
        gradient_w = np.zeros(shape = (1,13)) #zero intitialization of gradient of loss ter
        gradient_b = 0 #zero intitialization of gradient of loss term w.r.t b
        batch_data=x.sample(n = batch_size) #randomly sampling datapoints from dataset
        batch_x= np.array(batch_data.drop('Price',axis=1)) # separating our x_i's from batd
        batch_y = np.array(batch_data['Price']) # separating our y_i's from dataset from bd
        #Here we are calculating gradient of loss term w.r.t w and w.r.t b
        for j in range(batch size):
            y = np.dot(w old,batch x[j])+b old
            gradient_w += (batch_x[j]*(batch_y[j]-y)) + 2*alpha*w_old
            gradient_b += (batch_y[j]-y)
        gradient_w *= (-2/batch_size)
        gradient b *= (-2/batch size)
        #updating our weights
        w_new = w_old - eta*gradient_w
        b_new = b_old - eta*gradient_b
        #updating our old weights
        w old = w new
        b old = b new
        eta = eta/pow(t,power_t) #Here we are updating our learning rate at each iteration
        t=(i*batch_size)+(alpha*eta0) #updating our invscaling parameter at each iteration
        #Here we are assigning values to various keys in the dictionary
        if i == (n iter-1)
```

```
dict_1['alpha'] = alpha
            dict_1['weight'] = w_old
            dict 1['intercept'] = b old
            break
    return dict 1
def sgdreg_predict(w,b,x):
    """This function will predict the y_i values corresponding to our query points.
       Input: This function takes 'w' which is the best weight on which our dataset has bee
              This function also takes 'b' which is the best intercept term on which our da
              'x' is our test dataset.
       Output: This function returns an array of predicted values.
   y_pred = []
    for i in range(x.shape[0]):
        y_pred.append(np.dot(w,x[i])+b)
    return np.array(y_pred)
def error(y_true,y_pred):
    """This function calculates mean squared error"""
    error = mean_squared_error(y_true,y_pred)
    return error
def plot(y_true,y_pred):
    """This function plots y_pred as a function of y_true"""
    plt.figure()
    plt.scatter(y = y_pred,x = y_true,label = 'Actual values Vs Pred values')
    plt.xlabel('Actual values')
    plt.ylabel('Predicted values')
    plt.title('Actual values Vs Pred values')
    plt.legend()
    plt.grid()
    plt.show()
```

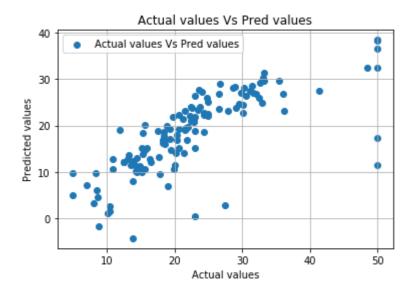
```
In [14]:
```

```
#Here we are performing cross validation w.r.t our training dataset using custom built SGD
#- alpha and eta0.
error i = []
err 1 = dict()
alpha = [10**-6,10**-5,10**-4,10**-3,10**-2,10**-1,1,10**1,10**2,10**3,10**4,10**5,10**6]
eta0 = [10**-6,10**-5,10**-4,10**-3,10**-2,10**-1,1,10**1,10**2,10**3,10**4,10**5,10**6]
for i in alpha:
    for j in eta0:
        sgd = sgdreg_fit(d_train,alpha = i,eta0 = j)#Fitting our training dataset
        y pred = sgdreg predict(sgd['weight'],sgd['intercept'],standardized train) #predict
        # the weights that our model has learned.
        err = error(y_train,y_pred) #calculating our mean squared error
        error_i.append(err) #Appending our errors
        err_1[err] = [i,j,err] #storing the mean squared error corresponding to a particula
        print('alpha = \{0\}, eta = \{1\}, mse = \{2\}'.format(i,j,err))
alpha = 1e-06, eta = 1e-06, mse = 585.5912974903646
alpha = 1e-06, eta = 1e-05, mse = 568.4120011329558
alpha = 1e-06, eta = 0.0001, mse = 477.42740964084373
alpha = 1e-06, eta = 0.001, mse = 185.4550531700009
alpha = 1e-06, eta = 0.01, mse = 537.4241268396811
alpha = 1e-06, eta = 0.1, mse = 9102493.203844715
alpha = 1e-06, eta = 1, mse = 8.009229853387747e+17
alpha = 1e-06, eta = 10, mse = 1.4330462550748914e+31
alpha = 1e-06, eta = 100, mse = 2.418153055558553e+44
alpha = 1e-06, eta = 1000, mse = 2.202044048965857e+60
alpha = 1e-06, eta = 10000, mse = 7.517358960021271e+76
alpha = 1e-06, eta = 100000, mse = 5.42896297609114e+95
alpha = 1e-06, eta = 1000000, mse = 6.164461811029461e+115
alpha = 1e-05, eta = 1e-06, mse = 588.0098637370008
alpha = 1e-05, eta = 1e-05, mse = 577.1803377850713
alpha = 1e-05, eta = 0.0001, mse = 522.5942352294663
alpha = 1e-05, eta = 0.001, mse = 304.980486739748
alpha = 1e-05, eta = 0.01, mse = 115.92005296807415
alpha = 1e-05, eta = 0.1, mse = 112224.76569128312
In [19]:
a = min(error_i)#Finding the alpha and eta0 values corresponding to minimum mean squared er
err 1[a]
Out[19]:
[0.1, 0.1, 37.84139064712439]
In [20]:
#Here we are trying to predict the y_i's corresponding to x_i's of test dataset using our b
sgd = sgdreg fit(d train,alpha = 0.1,eta0 = 0.1)
y pred = sgdreg predict(sgd['weight'],sgd['intercept'],standardized test)
```

In [22]:

```
print(error(y_test,y_pred))
plot(y_test,y_pred)
```

58.59034191193969



Observation:

From above plot we can see that the actual values and predicted values are linearly aligned with each other which indicates that our custom model is predicting our actual values with very less error and is thus somewhat perfectly performing Stoch astic Gradient Descent Regression.

In [23]:

```
w_1 = sgd['weight'] #weights corresponding to our best fit model
w_1
```

Out[23]:

```
array([[ 0.38355789, -0.65461713, -0.54872229, 1.18067096, -0.61003668, 3.54662039, -0.34283599, -0.79518735, -1.05936229, -0.67868092, -1.81454486, 0.96494767, -2.69118232]])
```

4. Sklearn SGDRegressor

In [24]:

mse = make_scorer(mean_squared_error) #As mean_squared_error is not a scoring parameter in #- using make_score to make mean_squared_error as one of the scoring parameters.

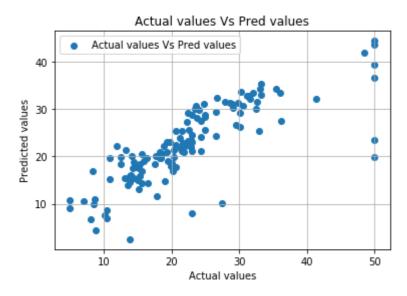
```
In [25]:
```

```
#Here we are performing gridsearch cross validation on our dataset for our sqd regressor
sgd = SGDRegressor(max_iter = 1000)
param_grid = \{ alpha': [10**-6,10**-5,10**-4,10**-3,10**-2,10**-1,1,10**1,10**2,10**3,10**4,10**-2,10**-1,1,10**1,10**1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**-1,10**
                                 'eta0':[10**-6,10**-5,10**-4,10**-3,10**-2,10**-1,1,10**1,10**2,10**3,10**4,1
gscv = GridSearchCV(sgd,param_grid,cv = 10,scoring = mse,verbose = 1)
gscv.fit(standardized_train,y_train)
print('*'*20)
Fitting 10 folds for each of 169 candidates, totalling 1690 fits
************
[Parallel(n jobs=1)]: Done 1690 out of 1690 | elapsed: 1.5min finished
In [26]:
df = pd.DataFrame(gscv.cv_results_)#Converting our grid search cross validation results int
print(df['mean_test_score'].min()) #Here we are trying to get the minimum mean squared erro
best_df = df[df['mean_test_score']==[df['mean_test_score'].min()]] #Here we are tring to ge
#- to the min mean_squared_error.
best_df
20.687995052417175
Out[26]:
         mean_fit_time mean_score_time mean_test_score mean_train_score param_alpha param_{ { } }
                  0.049928
                                                     0.000303
                                                                                    20.687995
                                                                                                                       18.543505
  56
                                                                                                                                                           0.01
1 rows × 32 columns
In [27]:
index = best df['param alpha'].index #Getting index corresponding to best values of hyperpd
index[0]
Out[27]:
56
In [28]:
print('Best hyperparameters are: alpha = {0} and eta0 = {1}'.format(best df['param alpha'][
Best hyperparameters are: alpha = 0.01 and eta0 = 0.01
In [29]:
#Here we are testing our model that we learnt above on our test dataset
sgd = SGDRegressor(alpha = 0.01,eta0 = 0.01,max iter = 1000)
sgd.fit(standardized_train,y_train)
y_pred = sgd.predict(standardized_test)
```

In [30]:

```
print(error(y_test,y_pred))
plot(y_test,y_pred)
```

33.651453623469465



Observation:

The above graph that we have got using sklearns Stochastic Gradient Descent Regres sor model is very much similar to our plot that we got using custom implementatio n.

In [31]:

```
w_2 = sgd.coef_ #weights corresponding to our best fit model
w_2
```

Out[31]:

```
array([-0.75595421, 0.78835386, 0.08226338, 0.62230719, -1.85857673, 3.6593889, -0.44289959, -2.70376005, 1.70737974, -1.63573206, -1.92452995, 0.82438522, -2.63974891])
```

In [32]:

np.linalg.norm(w_1-w_2) #Calculating euclidean distance between the weight vectors we got u #- own custom implementation.

Out[32]:

4.233214598939649

From above we can see that the weight vectors that we got using Custom implementat ion and sklearns implementation are very close to each other in 13 dimensional space as the distance between them is only 4.23 units.

5. Conclusion

In [33]:

Out[33]:

	Regressor	MSE	Weight	Distance	alpha	learning_rate
0	Custom SGDRegressor	58.59	[0.38355789, -0.65461713, -0.54872229, 1.18067	4.23	0.10	0.10
1	Sklearn SGDRegressor	33.65	[-0.75595421, 0.78835386, 0.08226338, 0.622307	4.23	0.01	0.01

From above table we can see that our weight vectors that we got using custom sgd r egressor implementation and sklearns sgd regressor implementation are very close t o each other as the distance between them is only 4.23 units which means that both the weights vectors are very similar to each other. This fact is also reflected in the mean squared error values which we got for both of these weight vectors. With Custom SGD Regressor we got best fit model for alpha = 0.1 and eta (or learning_r ate) = 0.1. Where as for Sklearns sgd we got best fit model for alpha = 0.01 and e ta = 0.01.

In []: