Regression Models Report

Improving machine learning skills by developing regression models

Student:

Anurag Banger

Student Id:

Introduction

Regression models are developed to analysis the data. We train our model so that it gives better output for our test data. In this report we will be developing 3 regression models: linear regression model, support vector regression model and k-nearest neighbour regression model. There are two types of machine learning algorithms: Regression which predicts continuous value outputs and Classification which predicts discrete outputs.

```
#Import the necessary libraries
from pandas.plotting import scatter_matrix
from pandas import read_csv
import pandas as pd
import numpy as np
from matplotlib import pyplot
import seaborn as seabornInstance

#Reading the data
filename = 'housing.csv'
names = ['CRIM', 'ZN', 'INDUS', 'CHAS', 'NOX', 'RM', 'AGE', 'DIS', 'RAD',
'TAX', 'PTRATIO', 'B', 'LSTAT', 'MEDV']
dataset = read_csv(filename, delim_whitespace=True, names=names)
```

By "print(dataset.shape)" we can get the shape of the dataset as (506, 14).

The info of the dataset can be checked by "dataset,info()"

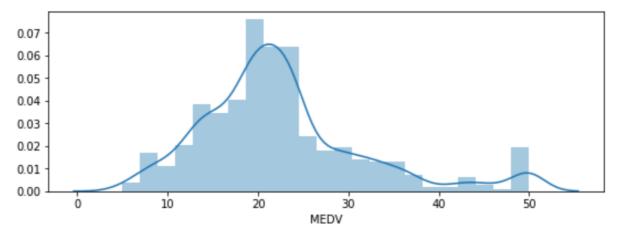
Below is the data description of the "housing.csv" data that we are using for this report.

	CRIM	ZN	INDUS	CHAS	NOX	RM	\
count	506.000000	506.000000	506.000000	506.000000	506.000000	506.000000	
mean	3.613524	11.363636	11.136779	0.069170	0.554695	6.284634	
std	8.601545	23.322453	6.860353	0.253994	0.115878	0.702617	
min	0.006320	0.000000	0.460000	0.000000	0.385000	3.561000	
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%	3.677082	12.500000	18.100000	0.000000	0.624000	6.623500	
max	88.976200	100.000000	27.740000	1.000000	0.871000	8.780000	
	AGE	DIS	RAD	TAX	PTRATIO	В	\
count	506.000000	506.000000	506.000000	506.000000	506.000000	506.000000	
mean	68.574901	3.795043	9.549407	408.237154	18.455534	356.674032	
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	375.377500	
50%	77.500000	3.207450	5.000000	330.000000	19.050000	391.440000	
75%	94.075000	5.188425	24.000000	666.000000	20.200000	396.225000	
max	100.000000	12.126500	24.000000	711.000000	22.000000	396.900000	
	LSTAT	MEDV					
count	506.000000	506.000000					
mean	12.653063	22.532806					
std	7.141062	9.197104					
min	1.730000	5.000000					
25%	6.950000	17.025000					
50%	11.360000	21.200000					
75%	16.955000	25.000000					
max	37.970000	50.000000					

Correlation data for the dataset is

```
CRIM
                                INDUS
                         ZN
                                           CHAS
                                                      NOX
                                                                 RM
                                                                          AGE
CRTM
        1.000000 - 0.200469 0.406583 - 0.055892 0.420972 - 0.219247 0.352734
ZN
        -0.200469 \quad 1.000000 \quad -0.533828 \quad -0.042697 \quad -0.516604 \quad 0.311991 \quad -0.569537
        0.406583 -0.533828 1.000000 0.062938 0.763651 -0.391676 0.644779
INDUS
        -0.055892 -0.042697 0.062938 1.000000 0.091203 0.091251 0.086518
NOX
        0.420972 -0.516604 0.763651 0.091203 1.000000 -0.302188 0.731470
       -0.219247 0.311991 -0.391676 0.091251 -0.302188 1.000000 -0.240265
RM
        0.352734 -0.569537 0.644779 0.086518 0.731470 -0.240265 1.000000
AGE
       -0.379670 0.664408 -0.708027 -0.099176 -0.769230 0.205246 -0.747881
DIS
        0.625505 -0.311948 0.595129 -0.007368 0.611441 -0.209847 0.456022
RAD
        0.582764 \ -0.314563 \quad 0.720760 \ -0.035587 \quad 0.668023 \ -0.292048 \quad 0.506456
PTRATIO 0.289946 -0.391679 0.383248 -0.121515 0.188933 -0.355501 0.261515
       -0.385064 0.175520 -0.356977 0.048788 -0.380051 0.128069 -0.273534
В
        LSTAT
MEDV
       -0.388305 \quad 0.360445 \ -0.483725 \quad 0.175260 \ -0.427321 \quad 0.695360 \ -0.376955
                        RAD
                                  TAX PTRATIO
                                                        В
        -0.379670 0.625505 0.582764 0.289946 -0.385064 0.455621 -0.388305
CRIM
        0.664408 -0.311948 -0.314563 -0.391679 0.175520 -0.412995 0.360445
ZN
INDUS
       -0.708027 0.595129 0.720760 0.383248 -0.356977 0.603800 -0.483725
CHAS
       -0.099176 \ -0.007368 \ -0.035587 \ -0.121515 \ \ 0.048788 \ -0.053929 \ \ 0.175260
NOX
        -0.769230 \quad 0.611441 \quad 0.668023 \quad 0.188933 \ -0.380051 \quad 0.590879 \ -0.427321
        0.205246 -0.209847 -0.292048 -0.355501 0.128069 -0.613808 0.695360
RM
       -0.747881 0.456022 0.506456 0.261515 -0.273534 0.602339 -0.376955
AGE
DIS
        1.000000 -0.494588 -0.534432 -0.232471 0.291512 -0.496996 0.249929
       -0.494588 1.000000 0.910228 0.464741 -0.444413 0.488676 -0.381626
RAD
TAX
       -0.534432 \quad 0.910228 \quad 1.000000 \quad 0.460853 \quad -0.441808 \quad 0.543993 \quad -0.468536
PTRATIO -0.232471 0.464741 0.460853 1.000000 -0.177383 0.374044 -0.507787
       0.291512 -0.444413 -0.441808 -0.177383 1.000000 -0.366087 0.333461
LSTAT
        -0.496996   0.488676   0.543993   0.374044   -0.366087   1.000000   -0.737663
MEDV
        0.249929 -0.381626 -0.468536 -0.507787 0.333461 -0.737663 1.000000
```

```
pyplot.figure(figsize=(9,3))
pyplot.tight_layout()
seabornInstance.distplot(dataset['MEDV'])
```



Our regression model should predict output similar to the above graph.

```
#Train Test Split
from sklearn.model_selection import train_test_split
array = dataset.values
X = array[:,0:13]
Y = array[:,13]
validation_size = 0.20
seed = 85
X_train, X_test, Y_train, Y_test = train_test_split(X, Y, test size=validation size, random state=seed)
```

Using the above code, we are splitting the dataset into train and test at the random state of 85. In the "train_test_split" function initially I used the random state as 50, I was getting a high error. So, I tried increased the random state to 70, the error reduced and the r2_score increased. However, when I increase it to 90 the error increase and the r2_score decreased. Finally, I Settled for 85 as the best random state for all the below regression models.

The final step is to evaluate the performance of the algorithm. This step is particularly important to compare how well different algorithms perform on a particular dataset. For regression algorithms, three evaluation metrics are commonly used:

1. **Mean Absolute Error** (MAE) is the mean of the absolute value of the errors.

$$\mathsf{MAE} = \frac{1}{n} \sum_{j=1}^{n} |y_j^{} - y_j^{}|$$

2. **Mean Squared Error** (MSE) is the mean of the squared errors and is calculated as:

$$MSE = \frac{1}{N} \sum_{i}^{n} (Y_i - y_i)^2$$

3. **Root Mean Squared Error** (RMSE) is the square root of the mean of the squared errors:

$$RMSE = \sqrt{\frac{1}{n} \sum_{j=1}^{n} (y_j - \hat{y}_j)^2}$$

Linear Regression Model

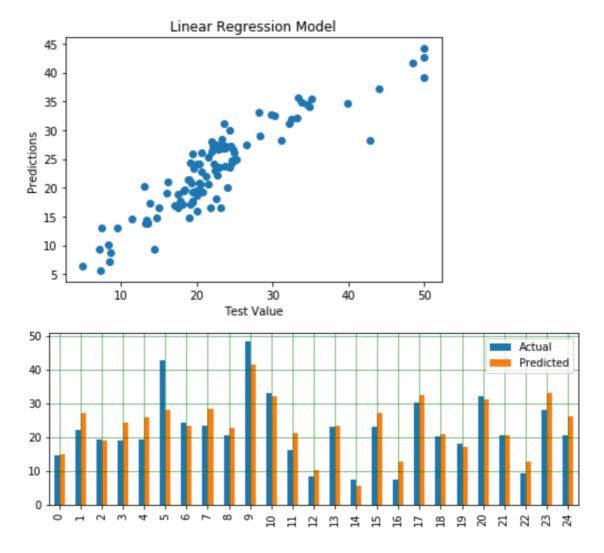
Linear regression performs the tasks to predict a dependent variable (Y) using the given variable (X). We will be using multiple linear regression for this problem. For multivariable linear regression, the linear the regression model has to find the most optimal coefficients for all the attributes. From the table 1, we can that for a unit increase in "nitric oxides concentration (NOX)", there is a decrease of 20.38 units in the median value and a unit decrease in "average number of rooms per dwelling (RM)" there is an increase of 3.14 units in the median value. We can check the difference between the actual and predicted value from the table 2 and the graph of actual and predicted.

Table 1

	Coefficient
CRIM	-0.110900
ZN	0.048214
INDUS	0.002390
CHAS	2.669476
NOX	-20.384569
RM	3.146873
AGE	0.012821
DIS	-1.594610
RAD	0.319752
TAX	-0.012424
PTRATIO	-0.996783
В	0.008374
LSTAT	-0.571749

Table 2

	Actual	Predicted
0	14.8	14.868280
1	22.1	27.281833
2	19.5	19.213757
3	19.2	24.342587
4	19.4	25.964260
5	42.8	28.179739
6	24.4	23.538190
7	23.3	28.464132
8	20.6	22.702733
9	48.5	41.669042
10	33.2	32.183417
11	16.2	21.107817
12	8.3	10.196391
13	23.0	23.472528
14	7.4	5.663803
15	23.2	27.166357
16	7.5	12.942476
17	30.3	32.473298
18	20.4	20.817842
19	18.1	17.062482
20	32.2	31.093483
21	20.5	20.646620
22	9.5	12.973219
23	28.2	33.143395
24	20.7	26.111568



regressor = LinearRegression(normalize=True)

Output for linear regression model:

Mean Absolute Error: 2.7770165268624316 Mean Squared Error: 14.053146682822419 Root Mean Squared Error: 3.74875268360323

r2_score: 0.8300388580559296

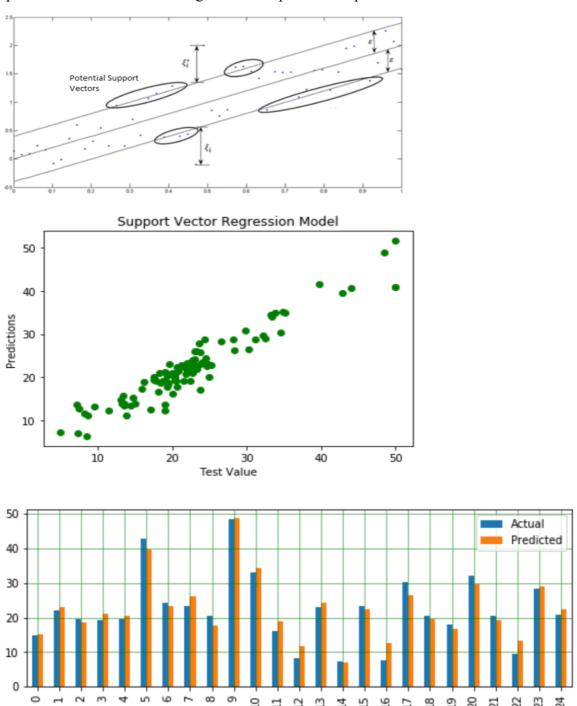
explained variance Score: 0.8328922853022158

For this model I tuned the 'random state' to 85 to get such low error.

SVM Model

Support Vector regression is a type of Support vector machine that supports linear and non-linear regression. As it seems in the below graph, the mission is to fit as many instances as possible between the lines while limiting the margin violations. The violation concept in this example represents as ϵ (epsilon). I have used the "StandardScaler" function to normalize the data so that it can be easier for the model to read the data and make better predictions.

To build a SVR model we have to first choose a kernel and parameters. I trained my model using 'rbf', 'linear' and 'poly' kernel. I found out the kernel = 'rbf' gives the best output. Changing the epsilon and C value will further give a more optimised output.



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	Actual	Predicted	
0	14.8	15.282607	
1	22.1	23.068043	
2	19.5	18.727181	
3	19.2	21.064892	
4	19.4	20.386563	
5	42.8	39.541974	
6	24.4	23.299517	
7	23.3	26.069507	
8	20.6	17.798168	
9	48.5	48.899721	
10	33.2	34.391573	
11	16.2	18.880577	
12	8.3	11.528066	
13	23.0	24.158698	
14	7.4	6.932648	
15	23.2	22.471696	
16	7.5	12.670828	
17	30.3	26.386681	
18	20.4	19.485129	
19	18.1	16.649052	
20	32.2	29.586668	
21	20.5	19.164047	
22	9.5	13.286589	
23	28.2	28.842379	
24	20.7	22.460207	

model = SVR(kernel='rbf', C=100)

Output for SVR model:

Mean Absolute Error: 2.045683998294405 Mean Squared Error: 7.432523737676976

Root Mean Squared Error: 2.7262655295618172

r2_score: 0.910109796012726

explained variance Score: 0.9107069269827175

K-Nearest Neighbour Model

"KNeighborsRegressor" implements learning based on the k nearest neighbours of each query point, where k is an integer value specified by the user. KNN used in the variety of applications such as finance, healthcare, political science, handwriting detection, image recognition and video recognition. Normalization of the data in this model has been done using "StandardScaler". Cross-validation is when the dataset is randomly split up into 'k' groups. One of the groups is used as the test set and the rest are used as the training set. The model is trained on the training set and scored on the test set. Then the process is repeated until each unique group has been used as the test set. Below are the cross validation scores for the KNN model. In our model we are using a 6 fold cross validation to get the best output. The train-test-split method we used in earlier is called 'holdout'. Cross-validation is better than using the holdout method because the holdout method score is dependent on how the data is split into train and test sets. Cross-validation gives the model an opportunity to test on multiple splits so we can get a better idea on how the model will perform on unseen data.

```
[0.82105336 0.6195712 0.74470466 0.78805673 0.78792551 0.80417367] cv_scores mean:{} 0.7609141886383952
```

Hypertuning parameters is when you go through a process to find the optimal parameters for your model to improve accuracy. In our case, we will use GridSearchCV to find the optimal value for 'n_neighbors'. GridSearchCV works by training our model multiple times on a range of parameters that we specify. That way, we can test our model with each parameter and figure out the optimal values to get the best accuracy results. The best params we get it for the n_neighbour of 3. KNeighborsRegressor has many parameters out of which I tuned weights and algorithm. I encountered 2 types of weights: uniform and distance, distance weight had lesser error and higher r2_score compared to uniform weight for nearest neighbor = 3. The algorithm 'brute' is very efficient for a small data sample. However, as the number of samples increases the efficiency decreases. Algorithm 'kd_tree' is very fast because partitioning is performed only along the data axes. The algorithm 'ball_tree' is a single distance calculation between a test point and the centroid is sufficient to determine a lower and upper bound on the distance to all points within the node. Because of the spherical geometry of the ball tree nodes, it can out-

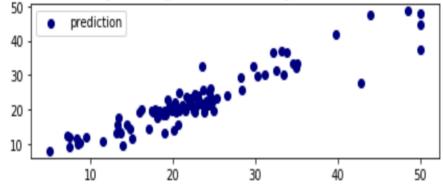
perform a *KD-tree* in high dimensions, though the actual performance is highly dependent on the structure of the training data.

Grid search using cross validation:

```
from sklearn.model_selection import GridSearchCV
knn2 = KNeighborsRegressor()
param_grid = {'n_neighbors': np.arange(1, 25)}
#use gridsearch to test all values for n_neighbors
knn_gscv = GridSearchCV(knn2, param_grid, cv=6)
#fit model to data
knn gscv.fit(X train std, Y train)
```

```
Actual Predicted
0
      14.8 14.429577
      22.1 23.296695
1
2
      19.5 19.948124
3
      19.2 19.655173
           22.990306
4
      19.4
5
      42.8 27.649959
6
      24.4
            24.879748
      23.3 22.058364
7
8
      20.6 15.507271
      48.5 48.910176
10
      33.2 37.118727
11
      16.2 20.022242
12
       8.3
           11.638765
13
      23.0 23.889208
14
       7.4
             9.239357
      23.2 21.569757
15
16
       7.5 12.004898
      30.3 29.872230
17
18
      20.4 19.371515
      18.1 17.668061
19
      32.2 36.660045
20.5 19.620360
20
21
      9.5 12.036435
22
      28.2 29.301961
23
24
      20.7 25.033085
```

KNeighborsRegressor (k = 3, weights = 'distance')



knn = KNeighborsRegressor(3, weights='distance')

Output for KNN:

```
Mean Absolute Error: 2.277490832961668
Mean Squared Error: 10.503144584772654
Root Mean Squared Error: 3.240855532845093
```

r2_score: 0.8729731861538417

explained variance Score: 0.8736631499908325

Conclusion

In the above 3 models general 6 Steps are being performed: importing libraries, loading the dat asets, dividing the train and test datasets, feature scaling, fitting the support vector regression m odel, predicting the result and visualising the support vector results by comparing it with the test t results. By observing the above regression models we can conclude that SVR model with kern el as 'rbf' and C as 100 has the best predictions. The output of SVR shows that mean absolute e rror, mean squared error and root mean squared error are lowest and r2_score and explained va riance score are the highest among all the 3 models. We can see that the value of root mean squared error is 2.726 which is greater than 10% of the mean value which is 22.533. This means t hat our algorithm is not very accurate but can still make reasonably good predictions.

Reference

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