```
1 %% MACHINE LEARNING PROJECT
  2 % ALGORITHMS PERFORMING THE PREDICTION OF PRICES OF USED CARS (REGRESSION TASK)
  4 % The two algorithms to be performed and compared are :
  5 % 1. Random Forrest Regressor
  6 % 2. Decision Tree for Regression
  7
  8 %% About the dataset
 10 % The data set contains 17,965 rows with 9 columns. The columns and their \checkmark
descriptions are as follows:
 11
 12 %- model --> Model Name.
 13 %- year --> The year it was bought.
 14 %- price --> The price at which the used car will be sold.
 15 %- transmission --> The transmission type i.e. manual, automatic or semi-automatic.
 16 %- mileage --> The miles that used car has driven.
 17 %- fuelType --> The fuel type of the car i.e. petrol, diesel, hybrid, electric or ✓
other
 18 %- tax --> The tax that will be applied on the selling price of that used car.
 19 %- mpg --> The miles per gallon ratio telling us how many mies it can drive per ✓
gallon of fuel.
 20 %- engineSize --> The engine size of the used car.
 22 %% Uploading the ford dataset from the 100,000 UK Used Car Data set from kaggle
 24 % reading my ford csv file & converting it into a table
 25 fordcar = readtable("C:\Users\FIDVI FATEMA\OneDrive\Desktop\machine learning\10k ✓
used car dataset\ford.csv");
 26 % displaying the head of the dataset
 27 head(fordcar);
 29 %% EXPLORATORY DATA ANALYSIS
 30
 31 %% 1. Checking Missing Values in my dataset
 32 clc
 33 missing values = sum(ismissing(fordcar));
 34
 35 % Displaying the number of missing values for each variable
 36 disp('Number of missing values for each variable:');
 37 disp(missing values);
 38
 39 %there are no missing values in my dataset
 41 %% 2. Checking if there are any duplicates in my dataset
 42 clc
 43 [~, unique indices, ~] = unique(fordcar);
 44
 45 % Getting the duplicate rows and removing them
  46 duplicate rows = fordcar(setdiff(1:height(fordcar), unique indices), :);
  47 fordcar noduplicates = fordcar(unique indices, :);
```

```
48 % There were 154 duplicate rows
 49
 50 % Looking at the dimensions of the new fordcar noduplicates dataset after the {m \ell}
removal of the duplilcates
 51 dimensions = size(fordcar_noduplicates);
 52 disp(['Number of rows: ', num2str(dimensions(1))]);
 53 disp(['Number of columns: ', num2str(dimensions(2))]);
 55 % After the removal of the duplicates, the shape of fordcar final is (17811 \times 9)
 57 %% 3. Descriptive Summary of the variables in my dataset
 59 summary(fordcar noduplicates)
 60
 61 % By doing this, we can see the minimum, maximum and the median values of
  62 % each of the clumns in the dataset.
 64 %% 4. Visually analyzing the distribution for price column
 65
 66 % Extracting the 'price' column
 67 price = fordcar noduplicates.price;
 68 % Creating a histogram
 69 figure;
 70 histogram(price, 'BinWidth', 5000);
 71 title('Distribution of Price');
 72 xlabel('Price');
 73 ylabel('Frequency');
 74
 75 % from this distribution graph, it is clear that the overwhelming majority
 76 % of price falls in the 0-25,000 range.
 78 %% 5. Understanding the distribution of categorical columns
 79 clc
 80 % Extracting the 'fuelType' column
 81 fuelType = fordcar noduplicates.fuelType;
 82 % Using the tabulate function to get fuelType counts
 83 fuelType_counts = tabulate(fuelType);
 84 % Displaying the results
 85 disp(fuelType counts);
 86
 87 % Performing the same for 'model' and 'transmission' columns
 88 model = fordcar noduplicates.model;
  89 model counts = tabulate(model);
 90 disp(model counts);
 91
 92 transmission = fordcar noduplicates.transmission;
 93 transmission counts = tabulate(transmission);
 94 disp(transmission counts);
 96 %% 6. Checking if there are outliers using boxplots
 97
```

```
98 % creating a list for numerical columns
 99 num list = {'price', 'mileage', 'tax', 'mpg'};
100
101 % Creating Boxplots for them
102 figure('Position', [100, 100, 900, 600]);
103 sqtitle('Distribution of Outliers', 'FontSize', 16);
104 \text{ rows} = 2;
105 \text{ cols} = 2;
106
107 % Iterating through numerical columns
108 for k = 1:numel(num list)
        % Creating subplot
109
110
        subplot(rows, cols, k);
111
        % Plot the boxplot
112
113
       boxplot(fordcar noduplicates.(num list{k}));
114
        title(num list{k}, 'FontSize', 14);
        if k > numel(num list) - cols
115
             xlabel('Variable');
116
117
        end
118
        if mod(k, cols) == 1
119
             ylabel('Value');
120
121
         end
122 end
124 %Even though there are a lot of outliers seen in the boxplots, removing them would \checkmark
not be the best option.
125 % For example, for price column, it is very much likely to have varied set of values \checkmark
ranging from a very high to very low values.
126 % Similarly, we can conclude for tax or mileage or mpg (mileage per gallon). If i\checkmark
remove them, it may break the flow of the dataset.
127 % Even though some of the data points can be actual outliers, I would conduct my \checkmark
research without removing them.
128
129 %% 7. Looking at the correlation between year and the sale price of the used cars
130
131 figure;
132 scatter(fordcar noduplicates.year, fordcar noduplicates.price, 'filled');
133 title('Ford cars sale price by year');
134 xlabel('Year');
135 ylabel('Price');
136
137 % From this, we can see that, the more recent the car is, the more pricier it is.
138 % There is one point that we can see as an outlier which shows the price of the car \checkmark
in 2060.
139 % This seems like an obvious outlier and should be removed from my dataset.
140
141 % removing the obvious outlier
142 logical index = (fordcar noduplicates.year ~= 2060);
143 fordcar final = fordcar noduplicates(logical index, :);
```

```
144
145 % Again looking at the distribution after removal of the outlier
146 figure;
147 scatter(fordcar final.year, fordcar final.price, 'filled');
148 title('Ford used car sale price by year');
149 xlabel('Year');
150 ylabel('Price');
151 %After removing this outlier, we can see the distribution of the price with respect \checkmark
to years much more clearly.
152 % In this, we can see that, the more recent the car is i.e, the closer the year of \checkmark
buying the car, the higher is the selling price of the car.
153 % That is, the lesser the age of the car, the higher is the price.
155 %% 8. Correlation matrix between my numeric variables
 156
157 % Extracting the columns of interest
158 selected columns final = fordcar final(:, {'price', 'year', 'mileage', 'tax', 'mpg', ✓
'engineSize'});
159
160 % Converting the table to a matrix
161 selected matrix = table2array(selected columns final);
162
163 % Calculating the correlation matrix
164 correlation matrix final = corrcoef(selected matrix, 'rows', 'pairwise');
165
166 % Creating a new figure and heatmap
167 figure;
168 heatmap(selected columns final.Properties.VariableNames, selected columns final. V
Properties. VariableNames, correlation matrix final, 'Colormap', bone, 'ColorLimits', [-1

✓
1], 'FontSize', 10, 'CellLabelColor', 'k', 'CellLabelFormat', '%.2f');
169 title('Correlation Matrix (Pearson) for Numeric Values');
170 xlabel('Variables');
171 ylabel('Variables');
172
173 % This heatmap tells us a lot about our dataset and how our variables are correlated \checkmark
to each other as well as with our target column i.e
174 % the price. The highest positive correlation with price column is 0.65 and it is \checkmark
with 'year' column, while the strongest negative
175 % correlation with price column is -0.53 and it is with 'mileage' column. The \checkmark
correlations of all numeric variables with my
176 % target column are quite high which is a good thing. The heatmap obtained from this \checkmark
code only has the numerical columns and
177 % not the categorical columns of 'model', 'fuelType' and 'transmission'as they are \checkmark
not label encoded yet.
178
179 %% 9. Distribution of the average price of the cars with the year column
181 groupedData = groupsummary(fordcar final, 'year', 'mean', 'price');
182 disp(groupedData);
183
 184 %from this, we can see that the prices of the older cars i.e the cars from 1996-1998 \checkmark
```

```
are sold at a higher rates and then there is
185 %decline in the prices after those years as cars from after 1998 are not considered \checkmark
as antiques but just older and non-functional.
186 %This might justify why we see a positice correlation between 'price' and 'year' ✓
column.
187
188 %% 10. Distribution of the models of ford car in the dataset
189 clc
190 model distribution = tabulate(fordcar final.model);
191 disp(model distribution);
193 %% 11. Correlation between model of the Ford car and the price
194 clc
195 % Extracting relevant columns
196 models = fordcar final.model;
197 prices = fordcar final.price;
198
199 % Creating a table with unique models and their corresponding total prices
 200 unique models = unique (models);
 201 total prices = zeros(size(unique models));
 203 for i = 1:length(unique models)
         total prices(i) = sum(prices(strcmp(models, unique models{i})));
 205 end
 206
 207 % Creating a bar plot
 208 figure;
 209 bar(unique models, total prices);
 210 title('Sum of Model Prices for Ford Cars');
 211 xlabel('Ford Models');
212 ylabel('Total Price');
 213 xtickangle(90);
 214
 215 % from this bar graph, it is very clear that there are 3 specific models of the ford \checkmark
company car namely 'Fiesta', 'Focus'
216 % & 'Kuga' which have a huge sum of prices i.e the sale price of these models might \checkmark
be higher than the others.
 217 % the other reason might be that the numbers of these models in our datset is \checkmark
greater than the other models and so
218 % the total prices of these models exceed the others. The latter seems more accurate \checkmark
looking at the distribution of the model columns above.
 219
220 %% Feature Engineering (Label Encoding)
 222 %'Label Encoding is a technique that is used to convert categorical columns into
numerical ones so that they can be fitted by machine
 223 % learning models which only takes numerical data.
224 % for my dataset, we have 3 categorical columns namely 'transmission', 'fuelType' 🗸
and 'model' and one 'year' column.
 225
 226 %% a) One-Hot Encoding:
```

```
227 clc
 228 % code help was taken from 'https://uk.mathworks. ✓
com/help/deeplearning/ref/onehotencode.html'
 229 % The 'transmission' column has 3 unique features while 'fuelType' column has 5\checkmark
unique features which are not ordinal.
230 \% So, the decision was made to apply one-hot encoding on these two categorical \checkmark
columns.
 231
 232
 233 transmission = fordcar final.transmission;
 234 fuelType = fordcar final.fuelType;
 235
 236 % Converting categorical variables to numerical indices
 237 transmission indices = grp2idx(categorical(fordcar final.transmission));
 238 fuelType indices = grp2idx(categorical(fordcar final.fuelType));
 240 % Performing one-hot encoding
 241 transmission encoded = full(sparse(1:numel(transmission indices), ✓
transmission indices, 1));
 242 fuelType encoded = full(sparse(1:numel(fuelType indices), fuelType indices, 1));
 244 % Concatenating the encoded variables to the original table
 245 fordcar final encoded = [fordcar final, array2table(transmission encoded, ✓
'VariableNames', strcat('transmission encoded ', cellstr(unique(fordcar final. 🗸
transmission))));
 246 fordcar final encoded = [fordcar final encoded, array2table(fuelType encoded, ✓
'VariableNames', strcat('fuelType encoded ', cellstr(unique(fordcar final.fuelType))))];
247
 248 % Displaying the head of the table
249 head(fordcar final encoded);
250
 251 %% b) Integer Label Encoding
252
 253 % This type of label encoding is the basic one as you can just assign an integer to \checkmark
the specific categorical value in the column.
254 % The disadvantage of using this type of labelEncoder is that the model assumes that \checkmark
the higher integer has more value than the
255 % lower one which is not true if the categorical variables in our column are not ✓
 256 % using this labelEncoding method for my 'model' column is more practical as the \checkmark
model column has 23 unique categorical
257 % values and using one-hot encoding will just increase the dimensionality and make \checkmark
my dataset more sparsed.
259 % Using integer label encoding on my 'model' column and assigning higher integer to 2
the more frequent models in my dataset.
 260
 261
 262 % Getting the unique models and their counts
 263 [uniqueModels, ~, modelIdx] = unique(fordcar final encoded.model);
 264 % Getting the frequency of each model
```

```
265 modelCounts = histcounts(modelIdx, 1:numel(uniqueModels)+1);
266 % Getting the rank of each model based on frequency
 267 [~, rankedModels] = sort(modelCounts, 'descend');
 268 % Assigning the ranked models as 'model encoded'
 269 fordcar final encoded.model encoded = zeros(size(fordcar final encoded, 1), 1);
270
271 % Loop through uniqueModels and assigning corresponding rank
272 for i = 1:numel(uniqueModels)
        modelIndices = ismember(fordcar final encoded.model, uniqueModels{i});
 274
         fordcar final encoded.model encoded(modelIndices) = rankedModels(i);
275 end
276
277 % We can later see whether the results of the algorithms vary if we just
278 % assign the variables in 'model' columns with integers without taking into
 279 % account the frequency of the model in our dataset.
281 %% c) Ordinal label encoding for my 'year' Column
282 % code help taken from 'https://uk.mathworks.com/help/matlab/matlab prog/ordinal-✓
categorical-arrays.html'
283 % Performing Ordinal encoding for my year column as the recent the year is, the \checkmark
higher the integer value it will be assigned.
 284 fordcar final encoded.year = categorical(fordcar final encoded.year);
285 fordcar final encoded.year encoded = grp2idx(fordcar final encoded.year);
286
287 % Dropping the original categorical columns
288 categorical columns to remove = { 'transmission', 'fuelType', 'model', 'year'};
289 fordcar final encoded = removevars(fordcar final encoded, ✓
categorical columns to remove);
 290
291 % Displaying the head of the updated table
292 head(fordcar final encoded)
294 % Dimensions of the dataset after encoding
295 dim = size(fordcar final encoded);
296 disp(['Number of rows: ', num2str(dim(1))]);
297 disp(['Number of columns: ', num2str(dim(2))]);
 298
299 %% Splitting the data into training, testing and validation (70:20:10 ratio)
 300
301 % 70 is the training set ratio
302 % 20 is the testing set ratio
 303 % 10 is the validation set ratio which we will use for performing Grid
 304 % Search for Hyperparameter tuning for our models
306 % Converting table to matrix
307 data = fordcar final encoded{:,:};
309 % Setting the random seed for reproducibility i.e each time we run this code, we get \checkmark
the same results.
310 rng(43);
 311
```

```
312 % Create a partition for 70% training and 30% testing
 313 cv = cvpartition(size(data, 1), 'HoldOut', 0.3);
 314
 315 % Training data
 316 training data = data(cv.training, :);
 318 % Hold-out data (combined testing and validation i.e 20:10)
 319 Holdoutdata = data(cv.test, :);
 321 % Further splitting the hold-out data into testing (20%) and validation (10%)
 322 cv2 = cvpartition(size(Holdoutdata, 1), 'HoldOut', 0.3333);
 323
 324 % Testing data
 325 testing data = Holdoutdata(cv2.training, :);
 326
 327 % Validation data
 328 validation data = Holdoutdata(cv2.test, :);
 329
 330 % Display the sizes of the datasets
 331 disp(['Training Data Size: ' num2str(size(training data, 1))]);
                                                                        %70
 332 disp(['Testing Data Size: ' num2str(size(testing data, 1))]);
                                                                        %20
 333 disp(['Validation Data Size: ' num2str(size(validation data, 1))]); %10
 334
 335 %% Extracting the features and target variable for training, testing and validation
 336
 337 % as price is our second column
 338 % Extracting features and target variable from validation data
 339 X validation = validation data(:, (2:end));
 340 y validation = validation data(:, 1);
 341
 342 % Extracting features and target variable from training data
 343 X train = training data(:, (2:end));
 344 y train = training data(:, 1);
 345
 346 % Extracting features and target variable from testing data
 347 X test = testing data(:, (2:end));
 348 y_test = testing_data(:, 1);
 349
 350 %-----
 351 %% RANDOM FORREST REGRESSION ALGORITHM
 353 % Statistics and Machine Learning Toolbox installed
 355 %% 1. Grid Search for finding the best Hyperparameters for Random Forest model
 357 % Code help taken from 'https://uk.mathworks.com/help/stats/regression-tree-\checkmark
ensembles.html'
 358 clc
 359 % setting random seed for reproducibility
 360 rng(43);
 361
```

```
362 % Defining the range of hyperparameters to search over
 363 numTreesRange = 10:50:300;
 364 maxNumSplitsRange = 10:10:150;
 366 % Initializing variables to store results
 367 bestNumTrees = 0;
 368 bestMaxNumSplits = 0;
 369 \text{ rfbestR2} = -Inf;
 370 rfbestMSE = Inf;
 371
 372 % Performing grid search
 373 for numTrees = numTreesRange
 374
        for maxNumSplits = maxNumSplitsRange
 375
             % Creating Random Forest model for regression for grid search
 376
 377
             rfmodel gridsearch = TreeBagger(numTrees, X train, y train, 'Method', ✓
'regression', 'MaxNumSplits', maxNumSplits);
 378
 379
             % Evaluating random forest on the validation set
             y predrf validation = predict(rfmodel gridsearch, X validation);
 380
 381
 382
             % Calculating R-squared value
             rfR2 validation = corr(y predrf validation, y validation)^2;
 383
             % Calculating Mean Squared Error (MSE)
 384
 385
             rfMSE validation = mean((y predrf validation - y validation).^2);
 386
 387 % Checking if the current hyperparameters result in better R-squared value and \checkmark
lower MSE
 388
             if rfR2 validation > rfbestR2 && rfMSE validation < rfbestMSE</pre>
 389
                 bestNumTrees = numTrees;
 390
                 bestMaxNumSplits = maxNumSplits;
 391
                 rfbestR2 = rfR2 validation;
 392
                 rfbestMSE = rfMSE validation;
 393
             end
 394
         end
 395 end
 396
 397 % Displaying the best hyperparameters, R-squared and MSE value
 398 disp(['Best Number of Trees for random forest: ' num2str(bestNumTrees)]);
 399 disp(['Best MaxNumSplits for random forest: ' num2str(bestMaxNumSplits)]);
 400 disp(['Best R-squared Value for random forest after gridsearch: ' num2str⊻
(rfbestR2)]);
 401 disp(['Best MSE Value for random forest after gridsearch: ' num2str(rfbestMSE)]);
 403 % The computational time taken for this step is more than anticipated.
 404 % The result of this grid search gives us the hyperparameters to get the best \checkmark
results for our prediction.
 405
 406 %% 2. Performing Random Forest Regression on my training data using the best\checkmark
hyperparameters from GridSearch
 407 clc
```

```
408 % using 60 no. of decision trees and 150 splits from grid search to get best
 409 % results from training the model on training set.
 410
 411 % setting random seed for reproducibility
 412 rng(43);
 413
 414 % Best hyperparameters from grid search
 415 bestNumTrees = 60;
 416 bestMaxNumSplits = 150;
 417
 418 % Training my random forest model on training set
 419 randomforestmodel = TreeBagger(bestNumTrees, X train, y train, 'Method', ✓
'regression', 'MaxNumSplits', bestMaxNumSplits, 'OOBPredictorImportance', 'on');
 420
 421 % save the best model for random forest
 422 save ("bestmodel for random forest.mat", 'randomforestmodel');
 423
 424 % Making predictions on the training set
 425 y predrf train = predict(randomforestmodel, X_train);
 426
 427 % Calculating R-squared value and MSE value for my training model
 428 rfMSE train = mean((y train - y predrf train).^2);
 429 rfR2 train = 1 - sum((y train - y predrf train).^2) / sum((y train - mean(y train)). ✓
^2);
 430 disp(['MSE Value for Random Forest Model : ' num2str(rfMSE train)]);
 431 disp(['R-squared Value for Random Forest Model: ' num2str(rfR2 train)]);
 433 % Creating a table containing original and predicted prices
 434 trainingprediction rf = table(y train, y predrf train);
 435 trainingprediction rf.Properties. VariableNames = { 'Original Price', 'Predicted ✓
Price'};
 436
 437 % Displaying the top 10 rows of the table
 438 display(trainingprediction rf(1:10, :));
 439
 440 % Using the bestNumTrees as 60 and bestMaxNumSplits as 150 from gridsearch, the \checkmark
baseline model gives
 441 % a high r-squared value of 0.9215 which means that 92.15% variance in our predicted ✓
values are explained by the model.
 442 % Although MSE value of 1751473.18 is huge but it depends on the scale of our data \checkmark
which is varied.
 443
 444 %% 3. Visualizing the difference between Original price and the predicted price in \checkmark
training data
 445
 446 % Creating a line plot for predictive price and actual price for training set
 447 figure('Position', [100, 100, 800, 500]);
 448 plot(1:length(y train), y train, 'r-', 'LineWidth', 0.2);
 450 plot(1:length(y predrf train), y predrf train, 'b-', 'LineWidth', 0.2);
 451 grid on;
```

```
452 xlabel('Data Point Index');
453 ylabel('Price');
 454 legend('Actual Price', 'Predicted Price');
 455 title('Actual vs. Predicted Prices for training set for Random Forest');
 456
457 %% 4. Feature Importance selection for Random Forest
458 % code help taken from 'https://uk.mathworks.com/help/stats/select-predictors-for-✓
random-forests.html'
 459 rng(43);
460 % Extract feature importance scores
461 importance scores = randomforestmodel.OOBPermutedVarDeltaError;
 462 % Specify the number of top features to select
463 num top features = 5;
 464
 465 % Find indices of top N features
 466 [~, top feature indices] = maxk(importance scores, num top features);
 467 % Filter datasets to include only top features
 468 X train selected = X train(:, top feature indices);
 469 X validation selected = X validation(:, top feature indices);
 470 X test selected = X test(:, top feature indices);
 472 % Retrain Random Forest on the selected features
473 randomforestmodel selected = TreeBagger(bestNumTrees, X train selected, y train, ✓
'Method', 'regression', 'MaxNumSplits', bestMaxNumSplits);
 474
475 % Make predictions from the selected columns on the training set
476 y predrf train selected = predict(randomforestmodel selected, X train selected);
 477
478 % Calculate R-squared value and MSE value for my training model
479 rfMSE train selected = mean((y train - y predrf train selected).^2);
480 rfR2 train selected = 1 - sum((y train - y predrf train selected).^2) / sum((y train ✓
- mean(y train)).^2);
481
482 % Display R-squared value
483 disp(['MSE Value for Random Forest Model for selected columns : ' num2str⊻
(rfMSE train selected)]);
484 disp(['R-squared Value for Random Forest Model for selected models: ' num2str⊻
(rfR2 train selected)]);
485
486 % While Feature selection does seem like an important step for predictive
487 % modeling, its application to my model yielded minimal impact on performance.
 488 % Notably, the R-squared (r2) value remained consistent, and though there
 489 % was a modest reduction in Mean Squared Error (MSE), it was not substantial
 490 % enough to alter the overall model outcomes. This suggests that the initially
 491 % chosen features already provided relevant information for prediction,
 492 % and the removal of some features did not yield a significant improvement.
 493 % I prefer having a model with all features for interpretability, even if it
 494 % comes at the cost of a slightly more complex model.
 496 %% 5. Evaluating the baseline model on the test data (unseen data)
 497 clc
```

```
498 % evaluating the performance of random forest model on the unseen test data withoutarksim \prime
the feature selection.
 499 % setting random seed for reproducibility
 500 rng(43);
 501
 502 load ("bestmodel for random forest.mat");
 504 % Evaluating the baseline model on the testing set
 505 y predrf test = predict(randomforestmodel, X test);
 506
 507 % Calculating MSE and R-squared values on the testing set
 508 rfMSE_test = mean((y_test - y_predrf_test).^2);
 509 rfR2 test = 1 - sum((y test - y predrf test).^2) / sum((y test - mean(y test)).^2);
 510 disp(['MSE on Testing Set for Random Forest Model: ' num2str(rfMSE test)]);
 511 disp(['R-squared Value on Testing Set for Random Forest Model: ' num2str⊻
(rfR2 test)]);
 512
513 \% R2 value for my test set or unseen data is 0.91799 which is considered outstanding \checkmark
but the MSE value
 514 % of this model on my test set is 1828169.73. The lower the MSE value, the better \checkmark
the result is, but our
515 % MSE value depends on the variability of my data and I want to check whether this \checkmark
high mse value is for
516 % the entire data or just the test set via k-fold cross validation.
517
518 %% 6. Visualizing the difference between Original price and the predicted price in \checkmark
testing set
519
520 % Creating a line plot for predictive price vs. actual price on test set
 521 figure('Position', [100, 100, 800, 500]);
 522 plot(1:length(y test), y test, 'g-', 'LineWidth', 0.2);
 523 hold on;
 524 plot(1:length(y predrf test), y predrf test, 'm-', 'LineWidth', 0.2);
 525 grid on;
 526 xlabel('Data Point Index');
 527 ylabel('Price');
 528 legend('Actual Price', 'Predicted Price');
 529 title('Actual vs. Predicted Prices for test set for Random Forest');
 531 %% 7. Visualizing the predictive and actual prices for the training and testing data
 532
 533 % Creating a line plot for collectively seeing the performance of the model on \checkmark
training as well as testing set
 534 figure('Position', [100, 200, 1100, 500]);
 535 plot(y train, 'r', 'LineWidth', 0.2, 'DisplayName', 'Actual Training');
 536 hold on;
 537 plot(y_predrf_train, 'b', 'LineWidth', 0.2, 'DisplayName', 'Fitted Training');
 538 plot(length(y train) + (1:length(y test)), y test, 'g', 'LineWidth', 0.2,\checkmark
'DisplayName', 'Actual Testing');
 539 plot(length(y train) + (1:length(y test)), y predrf test, 'm', 'LineWidth', 0.2, ✓
'DisplayName', 'Fitted Testing');
```

```
540 grid on;
541
542 title('RANDOM FOREST MODEL PREDICTIONS (on training and test set)');
543 xlabel('Observation');
544 ylabel('Price');
545 legend('show');
546
547 %% 8. Performing K-fold Cross Validation on my entire dataset for Random Forest
549 % Performing K-fold Cross Validation after running my model on the training
550 % as well as testing data is crucial for robust model evaluation.
551 % It provides a more reliable estimate of the model's performance,
552 % reducing the risk of overfitting or underfitting observed with a single train-test \checkmark
split.
553
554 % setting random seed for reproducibility
555 rng(43);
556 % Combine the training, validation, and testing sets
557 combined data = [training data; validation data; testing data];
558
559 %extract the features to put in X and Y
560 X = combined data(:, (2:end));
561 y = combined data(:,1);
562
563 % Define k-fold partition
564 cvpart = cvpartition(size(X, 1), 'kfold', 10);
566 % Performing k-fold cross-validation
567 mse scores rf = [];
568 r2 scores rf = [];
569 for i = 1:cvpart.NumTestSets
 570
        % Extract training and testing data
571
        trainIdx = cvpart.training(i);
        testIdx = cvpart.test(i);
572
573
       X train = X(trainIdx, :);
574
       X \text{ test} = X(\text{testIdx, :});
 575
       y_train = y(trainIdx);
576
       y \text{ test} = y(\text{testIdx});
577
578
        % Getting my random forest model
        randomforestmodel = TreeBagger(bestNumTrees, X train, y train, 'Method', ✓
579
'regression', 'MaxNumSplits', bestMaxNumSplits);
580
        % Evaluating the model
581
       y predrf = predict(randomforestmodel, X_test);
582
583
584
        % Calculating MSE
 585
        kfMSE rf = mean((y predrf - y test).^2);
        mse_scores_rf = [mse scores rf; kfMSE rf];
 586
587
 588
        % Calculating R-squared value
```

```
589
         kfR2 rf = corr(y predrf, y test)^2;
 590
         r2 scores rf = [r2 scores rf; kfR2 rf];
 591
 592
        % Displaying results for each fold
         disp(['Fold ' num2str(i) ' - MSE: ' num2str(kfMSE_rf) ', R-squared: ' num2str⊄
 593
(kfR2 rf)]);
 594 end
 595
 596 % Analyzing overall results
 597 average mse rf = mean(mse scores rf);
 598 average r2 rf = mean(r2 scores rf);
 599 disp(['Average MSE for k-fold: ' num2str(average mse rf)]);
 600 disp(['Average R-squared for k-fold: ' num2str(average r2 rf)]);
 602 %The results obtained from k-fold cross-validation for our Random Forest model show \checkmark
promising performance.
 603 % The model effectively captures a substantial portion of the variability in the \checkmark
target variable.
604 % The consistent MSE values indicate that overfitting is not a concern; instead, the \checkmark
observed higher
605 % MSE values are suggestive of the inherent complexity or dispersion within the \checkmark
dataset.
606 % It is important to recognize that these elevated MSE values do not necessarily \checkmark
reflect poor model
607 % performance but rather highlight the nuanced characteristics of the data, \checkmark
signaling potential complexity or variability.
 608
 609 %-----
610 %% DECISION TREE REGRESSION ALGORITHM
 611
612 % For decision tree, I will be using the split data and the extracted features and \checkmark
target variables
 613 % for training, testing and validation that I used earlier for random forest \checkmark
regression model.
 614
 615 %% 1. Grid Search to find the best Hyperparameters for Decision Tree Model
 617 % code help taken from 'https://uk.mathworks.com/help/stats/decision-trees. ✓
 618 % Performing Grid Search method for getting the best Hyperparameters for Decisionarksim
Tree Model.
 619 clc
 620 % Set the random seed for reproducibility
 621 rng(43);
 622
 623 % Defining the range of hyperparameters to search over
 624 maxsplitsrange = 5:5:100;
 625 minLeafSizeRange = 5:5:50;
 627 % Initializing variables to store results
 628 bestMaxSplits = 0;
```

```
629 bestMinLeafSize = 0;
 630 dtbestR2 = -Inf;
 631 dtbestMSE = Inf;
 633 % Performing grid search on validation set
 634 for maxnumsplits = maxsplitsrange
        for minLeafSize = minLeafSizeRange
 636
             % Creating decision tree model for regression for gridsearch
 637
             dtreemodel gridsearch = fitrtree(X train, y train, 'MinLeafSize', ✓
minLeafSize, 'MaxNumSplits', maxnumsplits);
 639
 640
             % Evaluating the decision tree on the validation set
             y preddt validation = predict(dtreemodel gridsearch, X validation);
 641
 642
 643
             % Calculating R-squared value
 644
             dtR2 validation = corr(y preddt validation, y validation)^2;
 645
             % Calculating Mean Squared Error (MSE)
 646
             dtMSE validation = mean((y preddt validation - y validation).^2);
 647
             % Checking if the current hyperparameters result in better R-squared value oldsymbol{arepsilon}
and lower MSE
 649
             if dtR2 validation > dtbestR2 && dtMSE validation < dtbestMSE</pre>
                 bestMaxSplits = maxnumsplits;
 650
                 bestMinLeafSize = minLeafSize;
 651
                 dtbestR2 = dtR2 validation;
 652
 653
                 dtbestMSE = dtMSE validation;
 654
              end
 655
         end
 656 end
 657
 658 % Displaying the best hyperparameters and performance metrics
 659 disp(['Best MaxNumber of splits: ' num2str(bestMaxSplits)]);
 660 disp(['Best MinLeafSize: ' num2str(bestMinLeafSize)]);
 661 disp(['Best R-squared Value for Decision Tree Model: ' num2str(dtbestR2)]);
 662 disp(['Best MSE Value for Decision Tree Model: ' num2str(dtbestMSE)]);
 663
 664 % Computational time taken for gridsearch for decision tree is way lesser than that \checkmark
taken for random forest.
 665 % The result of this hyperparameter tuning gives us the best number of splits and \checkmark
the minimum leaf size so as
 666 % to obtain the best predictive model by using these hyperparameters.
667
 668 %% 2. Performing Decision Tree Model on my training set using the best ✓
hyperparameters from GridSearch
669 clc
 670 % to get the best results for our decision tree's model on our dataset, we'll be \checkmark
using the hyperparameters
 671 % we got from the grid search for our Decision Trees model
 673 % setting random seed for reproducibility
```

```
674 rng(43);
 675
 676 % Best hyperparameters from grid search
 677 bestMaxSplits = 100;
 678 bestMinLeafSize = 20;
 679
 680 \ % Training decision tree model for regression on training set
 681 dtreemodel = fitrtree(X train, y train, 'MinLeafSize', bestMinLeafSize, ✓
'MaxNumSplits', bestMaxSplits);
 682
 683 %save the best model for decision tree
 684 save ("bestmodel for decision tree.mat", "dtreemodel");
 685
 686 % Make predictions on the training set
 687 y preddt train = predict(dtreemodel, X train);
 689 % Calculate MSE and R2 values for the decision tree model
 690 dtMSE train = mean((y train - y preddt train).^2);
 691 dtR2 train = 1 - sum((y train - y preddt train).^2) / sum((y train - mean(y train)). ✓
^2);
692 disp(['R-squared Value for Decision Tree Model on training set: ' num2str⊻
(dtR2 train));
693 disp(['MSE Value for Decision Tree Model on training set: ' num2str(dtMSE train)]);
 695 % Creating a table containing original and predicted prices for the decision tree \checkmark
 696 trainingprediction dt = table(y train, y preddt train);
697 trainingprediction dt. Properties. VariableNames = { 'Original Price', 'Predicted ✓
Price'};
 698
 699 % Displaying the top 10 rows of the table for the decision tree model
700 display(trainingprediction dt(1:10, :));
701
 702 %% 3. Visualizing the difference between Original price and the predicted price in \checkmark
training data for Decision Tree
703
704 % Creating a line plot for predictive price and actual price for training set for \mathbf{k}'
the decision tree model
 705 figure('Position', [100, 100, 800, 500]);
706 plot(1:length(y train), y train, 'r-', 'LineWidth', 0.2);
 707 hold on;
 708 plot(1:length(y preddt train), y preddt train, 'b-', 'LineWidth', 0.2);
 709 grid on;
 710 xlabel('Data Point Index');
 711 ylabel('Price');
 712 legend('Actual Price', 'Predicted Price');
 713 title('Actual vs. Predicted Prices for training set for Decision Tree');
 714
 715 %% 4. Evaluating the baseline model on the test data (unseen data) for Decision Tree
 717 % setting random seed for reproducibility
```

```
718 rng(43);
 719
 720 load ("bestmodel for decision tree.mat");
 722 % Evaluating the decision tree model on the testing set
 723 y preddt test = predict(dtreemodel, X test);
 724
 725 % Calculating MSE and R-squared values on the testing set
 726 dtMSE test = mean((y test - y preddt test).^2);
 727 dtR2 test = 1 - sum((y test - y preddt test))^2) / sum((y test - mean(y test))^2);
728
729 % Displaying the MSE and R-squared values on the testing set for the decision tree \checkmark
model
 730 disp(['MSE on Testing Set (Decision Tree Model): ' num2str(dtMSE test)]);
731 disp(['R-squared Value on Testing Set (Decision Tree Model): ' num2str(dtR2 test)]);
733 % It is performing as good on the test set as it did on the training set.
734
735 %% 5. Visualizing the difference between Original price and the predicted price in \checkmark
testing set for Decision Tree
 737 % Creating a line plot for predictive price vs. actual price on test set for \checkmark
decision tree
738 figure('Position', [100, 100, 800, 500]);
739 plot(1:length(y test), y test, 'g-', 'LineWidth', 0.2);
740 hold on;
 741 plot(1:length(y preddt test), y preddt test, 'm-', 'LineWidth', 0.2);
742 grid on;
743 xlabel('Data Point Index');
744 ylabel('Price');
745 legend('Actual Price', 'Predicted Price (Decision Tree)');
746 title('Actual vs predicted price for test set for Decision Tree Model');
 748 %% 6. Visualizing the predictive and actual prices for the training and testing data \checkmark
for decision tree model
749
750 % Creating a line plot for collectively seeing the performance of the model on {m \varkappa}'
training as well as testing set
751 figure('Position', [100, 200, 1100, 500]);
752 plot(y train, 'r', 'LineWidth', 0.2, 'DisplayName', 'Actual Training');
753 hold on;
754 plot(y preddt train, 'b', 'LineWidth', 0.2, 'DisplayName', 'Fitted Training');
755 plot(length(y_train) + (1:length(y_test)), y_test, 'g', 'LineWidth', 0.2, \checkmark
'DisplayName', 'Actual Testing');
756 plot(length(y_train) + (1:length(y test)), y preddt test, 'm', 'LineWidth', 0.2, ✓
'DisplayName', 'Fitted Testing');
757 grid on;
758
 759 title('DECISION TREE MODEL PREDICTIONS (on training and testing set)');
 760 xlabel('Observation');
 761 ylabel('Price');
```

```
762 legend('show');
763
      %% 7. Performing K-fold Cross Validation on my entire dataset
764
765
766 % K-fold cross validation is done to make sure that my model is not overfitting the \checkmark
training data.
767 % It has the same motive of providing a more reliable estimate of the model's \checkmark
generalization performance,
768 % helping identify potential overfitting or underfitting issues and ensuring the \checkmark
model's consistency across diverse data subset.
770 % setting random seed for reproducibility
771 rng(43);
772
773 % Using the combined data and X and y features that we created above when performing \checkmark
the K-fold cross validation on random forest model.
774
775 % Defining k-fold partition
776 cvpart = cvpartition(size(X, 1), 'kfold', 10);
777
778 % Perform k-fold cross-validation on decision tree model (following the same \checkmark
procedure as for random forest)
779 mse scores dt = [];
780 r2 scores dt = [];
781 for i = 1:cvpart.NumTestSets
        % Extracting training and testing data
782
783
        trainIdx = cvpart.training(i);
        testIdx = cvpart.test(i);
784
        X train = X(trainIdx, :);
785
786
        X test = X(testIdx, :);
787
        y train = y(trainIdx);
788
         y test = y(testIdx);
789
790
       % Getting my decision tree model
       dtreemodel = fitrtree(X train, y train, 'MinLeafSize', bestMinLeafSize, ✓
791
'MaxNumSplits', bestMaxSplits);
792
793
         % Evaluating the model
794
         y preddt = predict(dtreemodel, X test);
795
796
         % Calculating MSE
797
        kfMSE dt = mean((y preddt - y test).^2);
798
        mse_scores_dt = [mse_scores_dt; kfMSE_dt];
799
800
         % Calculating R-squared value
801
         kfR2 dt = corr(y preddt, y test)^2;
802
         r2 scores dt = [r2 scores dt; kfR2 dt];
803
804
         % Displaying results for each fold
805
         disp(['Fold ' num2str(i) ' - MSE: ' num2str(kfMSE dt) ', R-squared: ' num2str

✓
(kfR2 dt)]);
```

```
806 end
807
808 % Analyzing overall results
 809 average mse dt = mean(mse scores dt);
810 average r2 dt = mean(r2 scores dt);
811 disp(['Average MSE for k-fold for Decision Tree model: ' num2str(average mse dt)]);
812 disp(['Average R-squared for k-fold for Decision Tree model: ' num2str⊻
(average r2 dt)]);
813
814 %-----
815 \% Q. Does scaling the data before applying Random Forest and Decision Tree models \checkmark
yield comparable results to unscaled data, or does it significantly impact model \checkmark
performance?
816 %(extra analytical question)
818 % - Scaling of my variables (Through minmax scaling)
820 % Before scaling
821 figure;
822 subplot(2, 1, 1);
823 histogram(X train(:, 1), 'BinEdges', linspace(min(X train(:, 1)), max(X train(:, ✓
824 title('Before Scaling - Feature 1');
826 % Calculate min and max values for scaling
 827 min values train = min(X train, [], 1);
 828 max values train = max(X train, [], 1);
829
830 % Scale training data
831 for feature index = 1:size(X train, 2)
       X train(:, feature index) = (X train(:, feature index) - min values train ✓
(feature index)) / (max values train(feature index) - min values train(feature index));
833 end
834
835 % Scale validation data using the same min and max values
836 for feature index = 1:size(X validation, 2)
        X_validation(:, feature_index) = (X_validation(:, feature_index) - 
min values train(feature index)) / (max values train(feature index) - min values trainarksim
(feature index));
838 end
839
840 % Scale test data using the same min and max values
841 for feature_index = 1:size(X_test, 2)
        X test(:, feature index) = (X test(:, feature index) - min values train ✓
(feature index)) / (max values train(feature index) - min values train(feature index));
843 end
844
845 % After scaling
846 subplot(2, 1, 2);
847 histogram(X train(:, 1), 'BinEdges', linspace(0, 1, 20));
848 title('After Scaling - Feature 1');
```

```
849
 850 % The subplots are just to check whether my scaling worked or not.
 852 %% Random Forest with scaled data
853 % 1.Grid Search
854 % setting random seed for reproducibility
855 rng(43);
856
857 % Defining the range of hyperparameters to search over
 858 numTreesRange = 10:50:300;
859 maxNumSplitsRange = 10:10:150;
860
861 % Initializing variables to store results
862 bestNumTrees = 0;
 863 bestMaxNumSplits = 0;
864 rfbestR2 = -Inf;
865 rfbestMSE = Inf;
866
867 % Performing grid search
 868 for numTrees = numTreesRange
         for maxNumSplits = maxNumSplitsRange
870
871
             % Creating Random Forest model for regression
             randomforestmodel = TreeBagger(numTrees, X_train, y train, 'Method', ✓
872
'regression', 'MaxNumSplits', maxNumSplits);
873
874
             % Evaluating random forest on the validation set
             y predrf = predict(randomforestmodel, X validation);
875
876
877
            % Calculating R-squared value
878
            rfR2 = corr(y predrf, y validation)^2;
 879
             % Calculating Mean Squared Error (MSE)
880
             rfMSE = mean((y predrf - y validation).^2);
881
882 % Checking if the current hyperparameters result in better R-squared value and ✓
lower MSE
             if rfR2 > rfbestR2 && rfMSE < rfbestMSE</pre>
 883
                bestNumTrees = numTrees;
884
885
                bestMaxNumSplits = maxNumSplits;
                rfbestR2 = rfR2;
886
887
                 rfbestMSE = rfMSE;
888
             end
 889
         end
890 end
891
 892 % Display the best hyperparameters and R-squared value
893 disp(['Best Number of Trees: ' num2str(bestNumTrees)]);
 894 disp(['Best MaxNumSplits: ' num2str(bestMaxNumSplits)]);
 895 disp(['Best R-squared Value for random forest: ' num2str(rfbestR2)]);
 896 disp(['Best MSE Value for random forest: ' num2str(rfbestMSE)]);
 897
```

```
898 % 2. Running the model on training set for Random Forest (Scaled Data)
899
 900 % setting random seed for reproducibility
 901 rng(43);
 902
903 % Best hyperparameters from grid search
904 bestNumTrees = 60;
905 bestMaxNumSplits = 140;
 906
 907 % Training my random forest model on training set
908 randomforestmodel = TreeBagger(bestNumTrees, X train, y train, 'Method', ✓
'regression', 'MaxNumSplits', bestMaxNumSplits);
909
910 % Make predictions on the training set
911 y predrf train = predict(randomforestmodel, X train);
913 % Calculate R-squared value and MSE value for my training model
914 rfMSE train = mean((y train - y predrf train).^2);
 915 rfR2 train = 1 - sum((y train - y predrf train).^2) / sum((y train - mean(y train)).\checkmark
^2);
 916
917 % Display R-squared value
918 disp(['MSE Value for Random Forest Model : ' num2str(rfMSE train)]);
919 disp(['R-squared Value for Random Forest Model: ' num2str(rfR2 train)]);
 920
 921 % There isn't much of a difference between these results and the ones
 922 % obtained without scaling
 923
924 %% Decision Tree with scaled data
 925 % 1. Grid Search
 926 % Set the random seed for reproducibility
 927 rng(43)
928
929 % Define the range of hyperparameters to search over
 930 maxsplitsrange = 5:5:250;
931 minLeafSizeRange = 5:5:50;
 933 % Initialize variables to store results
934 bestMaxSplits = 0;
935 bestMinLeafSize = 0;
936 dtbestR2 = -Inf;
937 dtbestMSE = Inf;
 938
 939 % Performing grid search on validation set
 940 for maxnumsplits = maxsplitsrange
 941
        for minLeafSize = minLeafSizeRange
 942
 943
             % Creating decision tree model for regression
             dtreemodel = fitrtree(X train, y train, 'MinLeafSize', minLeafSize, ✓
'MaxNumSplits', maxnumsplits);
945
```

```
946
             % Evaluating the decision tree on the validation set
 947
             y preddt = predict(dtreemodel, X validation);
 948
 949
             % Calculating R-squared value
 950
             dtR2 = corr(y preddt, y validation)^2;
951
952
             % Calculating Mean Squared Error (MSE)
953
             dtMSE = mean((y preddt - y validation).^2);
954
 955
             % Checking if the current hyperparameters result in better R-squared value ✓
and lower MSE
 956
             if dtR2 > dtbestR2 && dtMSE < dtbestMSE</pre>
957
                 bestMaxSplits = maxnumsplits;
                 bestMinLeafSize = minLeafSize;
958
 959
                 dtbestR2 = dtR2;
960
                 dtbestMSE = dtMSE;
961
              end
962
         end
 963 end
 964
 965 % Display the best hyperparameters and performance metrics
 966 disp(['Best MaxNumber of splits: ' num2str(bestMaxSplits)]);
 967 disp(['Best MinLeafSize: ' num2str(bestMinLeafSize)]);
 968 disp(['Best R-squared Value for Decision Tree Model: ' num2str(dtbestR2)]);
 969 disp(['Best MSE Value for Decision Tree Model: ' num2str(dtbestMSE)]);
 971 % 2. Running the model on training set for Decision Tree (Scaled Data)
 972
973 % setting random seed for reproducibility
 974 rng(43);
975
 976 % Best hyperparameters from grid search
 977 bestMaxSplits = 100;
978 bestMinLeafSize = 20;
 979
980 % Creating decision tree model for regression
981 dtreemodel = fitrtree(X_train, y_train, 'MinLeafSize', bestMinLeafSize, <a href="mailto:Y">Y</a>
'MaxNumSplits', bestMaxSplits);
983 % Make predictions on the training set
984 y preddt train = predict(dtreemodel, X train);
 986 % Calculate MSE value for the decision tree model
 987 dtMSE train = mean((y train - y preddt train).^2);
988
989 % Calculate R-squared value for the decision tree model
 990 dtR2 train = 1 - sum((y train - y preddt train).^2) / sum((y train - mean(y train)).\checkmark
^2);
 991
992 % Display the calculated values
993 disp(['R-squared Value for Decision Tree Model: ' num2str(dtR2 train)]);
```

```
994 disp(['MSE Value for Decision Tree Model: 'num2str(dtMSE_train)]);
995
996 % I am getting the same values as I got before scaling for grid research for 
decision tree model.
997
998 % Scaling the data had minimal impact on the results of Random Forest, while having 
no impact on
999 % Decision Tree Model's performance . Consequently, it can be inferred that both 
Random Forest and Decision Tree algorithms
1000 % exhibit robustness to non-scaled data, performing admirably without the need for 
normalization or scaling.
```