# PROJECT IN R



As a part of this Project, install 'swirl' package. Then type, *library(swirl)* to initiate the package. Complete this interactive R tutorial. If you have followed this course thoroughly, this assignment should be an easy task for you!



## **Problem Statement**

The data scientists at BigMart have collected 2013 sales data for 1559 products across 10 stores in different cities. Also, certain attributes of each product and store have been defined. The aim is to build a predictive model and find out the sales of each product at a particular store.

Using this model, BigMart will try to understand the properties of products and stores which play a key role in increasing sales.

Please note that the data may have missing values as some stores might not report all the data due to technical glitches. Hence, it will be required to treat them accordingly.

#### **Evaluation Metric:**

Your model performance will be evaluated on the basis of your prediction of the sales for the test data (test.csv), which contains similar data-points as train except for the sales to be predicted. We at our end, have the actual sales for the test dataset, against which your predictions will be evaluated. We will use the Root Mean Square Error value to judge your response.

Where,

N: total number of observations

Predicted: the response entered by user

Actual: actual values of sales

$$RMSE = \sqrt{\frac{\sum_{i=1}^{N} (Predicted_{i} - Actual_{i})^{2}}{N}}$$



## Exploratory Data Analysis in R

From this section onwards, we'll dive deep into various stages of predictive modeling. Hence, make sure you understand every aspect of this section. In case you find anything difficult to understand, ask me through mail.

Data Exploration is a crucial stage of predictive model. You can't build great and practical models unless you learn to explore the data from begin to end. This stage forms a concrete foundation for data manipulation (the very next stage). Let's understand it in R.

I've taken the data set from Big Mart Sales Prediction, find the data set attached with the mail. Before we start, you must get familiar with these terms:

**Response Variable (a.k.a Dependent Variable)**: In a data set, the response variable (y) is one on which we make predictions. In this case, we'll predict 'Item\_Outlet\_Sales'. (Refer to image shown below)

**Predictor Variable (a.k.a Independent Variable)**: In a data set, predictor variables (Xi) are those using which the prediction is made on response variable. (Image in next slide).

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1	Item Identifier	Item Weight	Item Fat Conten	t Item Visibility	Item Type	tem MRP	Outlet Identifi	er Outlet Estab	olishment Year	Outlet Size	Outlet Location Typ	e Outlet Type	Item Outlet Sales	
2	FDA15	9.3	Low Fat	0.01604730	1 Dairy	249.8092	OUT049	1 2 2	199	Medium	Tier 1	Supermarket	3735.138	
3	DRC01	5.92	Regular	0.01927821	6 Soft Drinks	48.2692	OUT018		200	9 Medium	Tier 3	Supermarket	443.4228	
4	FDN15	17.5	Low Fat	0.01676007	5 Meat	141.618	OUT049		199	9 Medium	Tier 1	Supermarket	2097.27	
5	FDX07	19.2	Regular		Fruits and V	182.095	OUT010		199	8	Tier 3	Grocery Store	732.38	
5	NCD19	8.93	Low Fat		0 Household	53.8614	OUT013		198	7 High	Tier 3	Supermarket	994.7052	
,	FDP36	10.395	Regular		0 Baking Goo	51.4008	OUT018		2009	9 Medium	Tier 3	Supermarket	556.6088	
3	FDO10	13.65	Regular	0.01274108	9 Snack Food	57.6588	OUT013		198	7 High	Tier 3	Supermarket	343.5528	
9	FDP10		Low Fat	0.12746985	7 Snack Food	107.7622	OUT027		198	Medium	Tier 3	Supermarket	4022.7636	
0	FDH17	16.2	Regular	0.01668711	4 Frozen Fooi	96.9726	OUT045		200	2	Tier 2	Supermarket	1076.5986	
1	FDU28	19.2	Regular	0.0944495	9 Frozen Fooi	187.8214	OUT017		200	7	Tier 2	Supermarket	4710.535	
2	FDY07	11.8	Low Fat		Fruits and V	45.5402	OUT049		199	9 Medium	Tier 1	Supermarket	1516.0266	
3	FDA03	18.5	Regular	0.04546377	3 Dairy	144.1102	OUT046		199	7 Small	Tier 1	Supermarket	2187.153	
4	FDX32	15.1	Regular	0.100013	5 Fruits and V	145.4786	OUT049		1999	9 Medium	Tier 1	Supermarket	1589.2646	
5	FDS46	17.6	Regular	0.04725732	8 Snack Food	119.6782	OUT046		199	7 Small	Tier 1	Supermarket	2145.2076	
6	FDF32	16.35	Low Fat	0.068024	3 Fruits and V	196.4426	OUT013		198	7 High	Tier 3	Supermarket	1977.426	
7	FDP49	9	Regular	0.06908896	1 Breakfast	56.3614	OUT046		199	7 Small	Tier 1	Supermarket	1547.3192	
8	NCB42	11.8	Low Fat	0.00859605	1 Health and	115.3492	OUT018		2009	9 Medium	Tier 3	Supermarket	1621.8888	
9	FDP49	9	Regular	0.06919637	6 Breakfast	54.3614	OUT049		199	9 Medium	Tier 1	Supermarket	718.3982	
0	DRI11		Low Fat	0.03423768	2 Hard Drinks	113.2834	OUT027		198	5 Medium	Tier 3	Supermarket	2303.668	
1	FDU02	13.35	Low Fat	0.1024921	2 Dairy	230.5352	OUT035		200-	4 Small	Tier 2	Supermarket	2748.4224	
2	FDN22	18.85	Regular	0.13819027	7 Snack Food	250.8724	OUT013		198	7 High	Tier 3	Supermarket	3775.086	
13	FDW12		Regular	0.03539992	3 Baking Goo	144.5444	OUT027		198	Medium	Tier 3	Supermarket	4064.0432	



#### Data

We have train (8523) and test (5681) data set, train data set has both input and output variable(s). You need to predict the sales for test data set.

Variable	Description				
Item_Identifier	Unique product ID				
Item_Weight	Weight of product				
Item_Fat_Content	Whether the product is low fat or not				
Item_Visibility	The % of total display area of all products in a store allocated to the particular product				
Item_Type	The category to which the product belongs				
Item_MRP	Maximum Retail Price (list price) of the product				
Outlet_Identifier	Unique store ID				



Outlet_Establishment_Year	The year in which store was established
Outlet_Size	The size of the store in terms of ground area covered
Outlet_Location_Type	The type of city in which the store is located
Outlet_Type	Whether the outlet is just a grocery store or some sort of supermarket
Item_Outlet_Sales	Sales of the product in the particulat store. This is the outcome variable to be predicted.



**Train Data**: The predictive model is always built on train data set. An intuitive way to identify the train data is, that it always has the 'response variable' included.

**Test Data**: Once the model is built, it's accuracy is 'tested' on test data. This data always contains less number of observations than train data set. Also, it does not include 'response variable'.

Right now, you should download the data set. Take a good look at train and test data. Cross check the information shared above and then proceed.

Let's now begin with **importing and exploring data**.



You can import the data directly using the import button or follow the steps below to do manually.

#working directory

path <- ".../Data/BigMartSales"</pre>

#set working directory

setwd(path)

As a beginner, I'll advise you to keep the train and test files in your working directly to avoid unnecessary directory troubles.

Once the directory is set, we can easily import the .csv files using commands below.

#Load Datasets

train <- read.csv("Train\_UWu5bXk.csv")</pre>

test <- read.csv("Test\_u94Q5KV.csv")</pre>



In fact, even prior to loading data in R, it's a good practice to look at the data in Excel. This helps in strategizing the complete prediction modeling process. To check if the data set has been loaded successfully, look at R environment. The data can be seen there. Let's explore the data quickly.

```
#check dimesions ( number of row & columns) in data set
```

> dim(train)

[1] 8523 12

> dim(test)

[1] 5681 11

We have 8523 rows and 12 columns in train data set and 5681 rows and 11 columns in data set. This makes sense. Test data should always have one column less. Let's get deeper in train data set now.

```
#check the variables and their types in train
> str(train)
'data.frame': 8523 obs. of 12 variables:
$ Item Identifier : Factor w/ 1559 levels "DRA12", "DRA24", ..: 157 9 663 1122 1298 759
697 739 441 991 ...
$ Item Weight : num 9.3 5.92 17.5 19.2 8.93 ...
$ Item Fat Content : Factor w/ 5 levels "LF","low fat",..: 3 5 3 5 5 5 5 5 5 ...
$ Item Visibility : num 0.016 0.0193 0.0168 0 0 ...
$ Item_Type : Factor w/ 16 levels "Baking Goods",..: 5 15 11 7 10 1 14 14 6 6 ...
$ Item MRP : num 249.8 48.3 141.6 182.1 53.9 ...
$ Outlet Identifier : Factor w/ 10 levels "OUT010", "OUT013",..: 10 4 10 1 2 4 2 6 8 3
$ Outlet Establishment Year: int 1999 2009 1999 1998 1987 2009 1987 1985 2002 2007 ...
$ Outlet_Size : Factor w/ 4 levels "","High","Medium",..: 3 3 3 1 2 3 2 3 1 1 ...
$ Outlet Location Type : Factor w/ 3 levels "Tier 1", "Tier 2",..: 1 3 1 3 3 3 3 2 2
$ Outlet_Type : Factor w/ 4 levels "Grocery Store",..: 2 3 2 1 2 3 2 4 2 2 ...
```

\$ Item Outlet Sales : num 3735 443 2097 732 995 ...



To begin with, I'll first check if this data has missing values. This can be done by using:

```
> table(is.na(train))
```

#### FALSE TRUE

#### 100813 1463

In train data set, we have 1463 missing values. Let's check the variables in which these values are missing. It's important to find and locate these missing values. Many data scientists have repeatedly advised beginners to pay close attention to missing value in data exploration stages.

## > colSums(is.na(train))

Hence, we see that column Item\_Weight has 1463 missing values. Let's get more inferences from this data.

## > summary(train)

Here are some quick inferences drawn from variables in train data set:

- 1. Item\_Fat\_Content has mismatched factor levels.
- 2. Minimum value of item\_visibility is o. Practically, this is not possible. If an item occupies shelf space in a grocery store, it ought to have some visibility. We'll treat all o's as missing values.
- 3. Item\_Weight has 1463 missing values (already explained above).
- 4. Outlet Size has a unmatched factor levels.

These inference will help us in treating these variable more accurately.



## **Graphical Representation of Variables**

I'm sure you would understand these variables better when explained visually. Using graphs, we can analyze the data in 2 ways: Univariate Analysis and Bivariate Analysis.

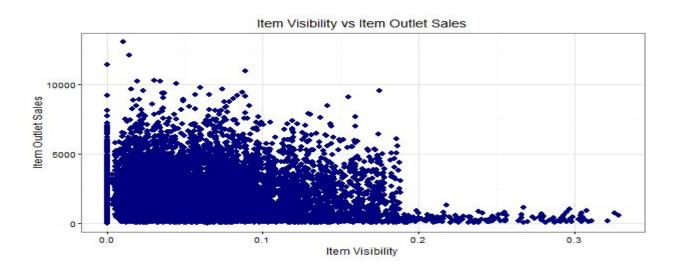
Univariate analysis is done with one variable. Bivariate analysis is done with two variables. Univariate analysis is a lot easy to do. Hence, I'll skip that part here. I'd recommend you to try it at your end. Let's now experiment doing bivariate analysis and carve out hidden insights.

For visualization, I'll use ggplot2 package.



These graphs would help us understand the distribution and frequency of variables in the data set.

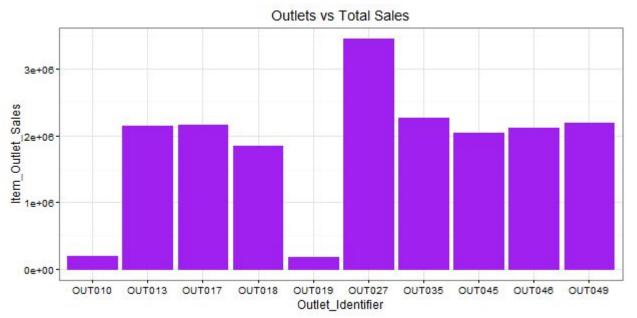
```
> ggplot(train, aes(x= Item_Visibility, y = Item_Outlet_Sales)) + geom_point(size = 2.5, color="navy") +
xlab("Item Visibility") + ylab("Item Outlet Sales") + ggtitle("Item Visibility vs Item Outlet Sales")
```



We can see that majority of sales has been obtained from products having visibility less than 0.2. This suggests that item\_visibility < 2 must be an important factor in determining sales.

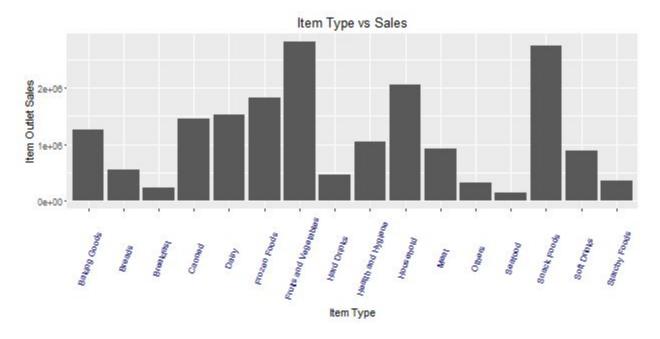
Let's plot few more interesting graphs and explore such hidden stories.

```
> ggplot(train, aes(Outlet_Identifier, Item_Outlet_Sales)) + geom_bar(stat = "identity", color = "purple")
+theme(axis.text.x = element_text(angle = 70, vjust = 0.5, color = "black")) + ggtitle("Outlets vs Total Sales")
+ theme_bw()
```



Here, we infer that OUT027 has contributed to majority of sales followed by OUT35. OUT10 and OUT19 have probably the least footfall, thereby contributing to the least outlet sales.

```
> ggplot(train, aes(Item_Type, Item_Outlet_Sales)) + geom_bar( stat = "identity") +theme(axis.text.x =
element_text(angle = 70, vjust = 0.5, color = "navy")) + xlab("Item Type") + ylab("Item Outlet
Sales")+ggtitle("Item Type vs Sales")
```

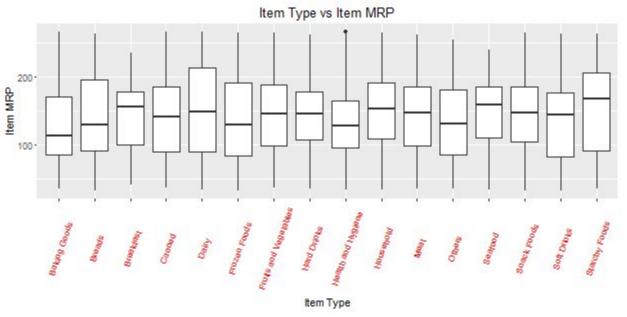


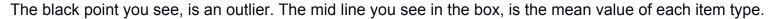
From this graph, we can infer that Fruits and Vegetables contribute to the highest amount of outlet sales followed by snack foods and household products. This information can also be represented using a box plot chart.



The benefit of using a box plot is, you get to see the outlier and mean deviation of corresponding levels of a variable (shown below).

```
> ggplot(train, aes(Item_Type, Item_MRP)) +geom_boxplot() +ggtitle("Box Plot") + theme(axis.text.x =
element_text(angle = 70, vjust = 0.5, color = "red")) + xlab("Item Type") + ylab("Item MRP") + ggtitle("Item Type
vs Item MRP")
```







Now, we have an idea of the variables and their importance on response variable. Let's now move back to where we started. Missing values. Now we'll impute the missing values.

We saw variable Item\_Weight has missing values. Item\_Weight is an continuous variable. Hence, in this case we can impute missing values with mean / median of item\_weight. These are the most commonly used methods of imputing missing value

Let's first combine the data sets. This will save our time as we don't need to write separate codes for train and test data sets. To combine the two data frames, we must make sure that they have equal columns, which is not the case.

```
> dim(train)
```

[1] 8523 12

> dim(test)

[1] 5681 11



Test data set has one less column (response variable). Let's first add the column. We can give this column any value. An intuitive approach would be to extract the mean value of sales from train data set and use it as placeholder for test variable Item \_Outlet\_ Sales. Anyways, let's make it simple for now. I've taken a value 1. Now, we'll combine the data sets.

```
> test$Item_Outlet_Sales <- 1</pre>
```

```
> combi <- rbind(train, test)</pre>
```

Impute missing value by median. I'm using median because it is known to be highly robust to outliers. Moreover, for this problem, our evaluation metric is RMSE which is also highly affected by outliers. Hence, median is better in this case.

```
> combi$Item_Weight[is.na(combi$Item_Weight)] <- median(combi$Item_Weight, na.rm = TRUE)</pre>
```

```
> table(is.na(combi$Item_Weight))
```

**FALSE** 

14204



## **Trouble with Continuous Variables & Categorical Variables**

It's important to learn to deal with continuous and categorical variables separately in a data set. In other words, they need special attention. In this data set, we have only 3 continuous variables and rest are categorical in nature. If you are still confused, I'll suggest you to once again look at the data set using *str()* and proceed.

Let's take up *Item\_Visibility*. In the graph above, we saw item visibility has zero value also, which is practically not feasible. Hence, we'll consider it as a missing value and once again make the imputation using median.

```
> combi$Item_Visibility <- ifelse(combi$Item_Visibility == 0,</pre>
```

median(combi\$Item\_Visibility), combi\$Item\_Visibility)



Let's proceed to categorical variables now. During exploration, we saw there are mis-matched levels in variables which needs to be corrected.

```
> levels(combi$Outlet_Size)[1] <- "Other"</pre>
> library(plyr)
> combi$Item Fat Content <- revalue(combi$Item Fat Content,</pre>
c("LF" = "Low Fat", "reg" = "Regular"))
> combi$Item Fat Content <- revalue(combi$Item Fat Content, c("low fat" = "Low Fat"))</pre>
> table(combi$Item Fat Content)
  Low Fat Regular
```

9185

5019

Using the commands above, I've assigned the name 'Other' to unnamed level in *Outlet\_Size* variable. Rest, I've simply renamed the various levels of Item\_Fat\_Content.



## 4. Data Manipulation in R

Let's call it as, the advanced level of data exploration. In this section we'll practically learn about feature engineering and other useful aspects.

**Feature Engineering:** This component separates an intelligent data scientist from a technically enabled data scientist. You might have access to large machines to run heavy computations and algorithms, but the power delivered by new features, just can't be matched. We create new variables to extract and provide as much 'new' information to the model, to help it make accurate predictions.

If you have been thinking all this time, great. But now is the time to think deeper. Look at the data set and ask yourself, what else (factor) could influence Item\_Outlet\_Sales? Anyhow, the answer is below. But, I want you to try it out first, before scrolling down.



1. Count of Outlet Identifiers – There are 10 unique outlets in this data. This variable will give us information on count of outlets in the data set. More the number of counts of an outlet, chances are more will be the sales contributed by it.

```
> library(dplyr)
> a <- combi%>%
            group_by(Outlet_Identifier)%>%
            tally()
> head(a)
Source: local data frame [6 x 2] Outlet Identifier n
(fctr)
                 (int)
1 OUT010
                 925
2 OUT013
                 1553
3 OUT017
                 1543
4 OUT018
                 1546
                 880
5 OUT019
```

1559

6 OUT027

```
> combi <- full_join(a, combi, by = "Outlet_Identifier")</pre>
```

As you can see, dplyr package makes data manipulation quite effortless. You no longer need to write long function. In the code above, I've simply stored the new data frame in a variable *a.* Later, the new column *Outlet\_Count* is added in our original 'combi' data set. To know more about dplyr, follow this <u>tutorial</u>.

2. **Count of Item Identifiers** – Similarly, we can compute count of item identifiers too. It's a good practice to fetch more information from unique ID variables using their count. This will help us to understand, which outlet has maximum frequency.

> head (b)

> names(a)[2] <- "Outlet Count"</pre>



```
Item_Identifier
                    Item_Count
(fctr)
                    (int)
1 DRA12
                      9
2 DRA24
                      10
3 DRA59
                      10
4 DRB01
                      8
5 DRB13
6 DRB24
                     8
> combi <- merge(b, combi, by = "Item_Identifier")</pre>
```



3. **Outlet Years** – This variable represent the information of existence of a particular outlet since year 2013. Why just 2013? You'll find the answer in problem statement. My hypothesis is, older the outlet, more footfall, large base of loyal customers and larger the outlet sales.

This suggests that outlets established in 1999 were 14 years old in 2013 and so on.



4. **Item Type New** – Now, pay attention to *Item\_Identifiers*. We are about to discover a new trend. Look carefully, there is a pattern in the identifiers starting with "FD", "DR", "NC". Now, check the corresponding *Item\_Types* to these identifiers in the data set. You'll discover, items corresponding to "DR", are mostly eatables. Items corresponding to "FD", are drinks. And, item corresponding to "NC", are products which can't be consumed, let's call them non-consumable. Let's extract these variables into a new variable representing their counts.

Here I'll use *substr()*, *gsub()* function to extract and rename the variables respectively.

```
> q <- substr(combi$Item_Identifier,1,2)
> q <- gsub("FD","Food",q)
> q <- gsub("DR","Drinks",q)
> q <- gsub("NC","Non-Consumable",q)
> table(q)
    Drinks Food Non-Consumable
    1317 10201 2686
```

Let's now add this information in our data set with a variable name 'Item Type New.

```
> combi$Item_Type_New <- q</pre>
```

I'll leave the rest of feature engineering intuition to you. You can think of more variables which could add more information to the model. But make sure, the variable aren't correlated. Since, they are emanating from a same set of variable, there is a high chance for them to be correlated. You can check the same in R using *cor()* function.

## **Label Encoding and One Hot Encoding**

Just, one last aspect of feature engineering left. Label Encoding and One Hot Encoding.

Label Encoding, in simple words, is the practice of numerically encoding (replacing) different levels of a categorical variables. For example: In our data set, the variable *Item\_Fat\_Content* has 2 levels: Low Fat and Regular. So, we'll encode Low Fat as o and Regular as 1. This will help us convert a factor variable in numeric variable. This can be simply done using if else statement in R.

```
> combi$Item_Fat_Content <- ifelse(combi$Item_Fat_Content == "Regular",1,0)</pre>
```

One Hot Encoding, in simple words, is the splitting a categorical variable into its unique levels, and eventually removing the original variable from data set. Confused? Here's an example: Let's take any categorical variable, say, *Outlet\_Location\_Type*. It has 3 levels. One hot encoding of this variable, will create 3 different variables consisting of 1s and os. 1s will represent the existence of variable and os will represent non-existence of variable. Let look at a sample on next slide



*model.matrix* creates a matrix of encoded variables. ~. -1 tells R, to encode all variables in the data frame, but suppress the intercept. So, what will happen if you don't write -1? model.matrix will skip the first level of the factor, thereby resulting in just 2 out of 3 factor levels (loss of information).

This was the demonstration of one hot encoding. Hope you have understood the concept now.



```
>library(dummies)
>combi <- dummy.data.frame(combi, names = c('Outlet_Size','Outlet_Location_Type','Outlet_Type', 'Item_Type_New'),</pre>
```

With this, I have shared 2 different methods of performing one hot encoding in R. Let's check if encoding has been done.

Let's now apply this technique to all categorical variables in our data set (excluding ID variable).

```
> str (combi)
```

sep='\_')



```
$ Outlet Size Other : int 0 1 1 0 1 0 0 0 0 0 ...
$ Outlet_Size_High : int 0 0 0 1 0 0 0 0 0 0 ...
$ Outlet_Size_Medium : int 1 0 0 0 0 0 1 1 0 1 ...
$ Outlet Size Small : int 0 0 0 0 0 1 0 0 1 0 ...
$ Outlet_Location_Type_Tier 1 : int 1 0 0 0 0 0 0 0 1 0 ...
$ Outlet_Location_Type_Tier 2 : int 0 1 0 0 1 1 0 0 0 0 ...
$ Outlet_Location_Type_Tier 3 : int 0 0 1 1 0 0 1 1 0 1 ...
$ Outlet_Type_Grocery Store : int 0 0 1 0 0 0 0 0 0 0 ...
$ Outlet Type Supermarket Type1: int 1 1 0 1 1 1 0 0 1 0 ...
$ Outlet Type Supermarket Type2: int 0 0 0 0 0 0 1 0 0 ...
$ Outlet Type Supermarket Type3: int 0 0 0 0 0 0 1 0 0 1 ...
$ Item Outlet Sales : num 1 3829 284 2553 2553 ...
$ Year : num 14 11 15 26 6 9 28 4 16 28 ...
$ Item Type New Drinks : int 1 1 1 1 1 1 1 1 1 1 ...
$ Item Type New Food : int 0 0 0 0 0 0 0 0 0 0 ...
$ Item Type New Non-Consumable : int 0 0 0 0 0 0 0 0 0 0 ...
```

As you can see, after one hot encoding, the original variables are removed automatically from the data set.



# 5. Predictive Modeling using Machine Learning

Finally, we'll drop the columns which have either been converted using other variables or are identifier variables. This can be accomplished using *select* from dplyr package.

```
> combi <- select(combi, -c(Item_Identifier, Outlet_Identifier, Item_Fat_Content,
Outlet_Establishment_Year,Item_Type))
> str(combi)
```

In this section, I'll cover Regression, Decision Trees and Random Forest. A detailed explanation of these algorithms is provided in links attached. These algorithms have been satisfactorily explained in the links for useful resources.

As you can see, we have encoded all our categorical variables. Now, this data set is good to take forward to modeling. Since, we started from Train and Test, let's now divide the data sets.

```
> new_train <- combi[1:nrow(train),]
> new test <- combi[-(1:nrow(train)),]</pre>
```



## **Linear (Multiple) Regression**

Multiple Regression is used when response variable is continuous in nature and predictors are many. Had it been categorical, we would have used Logistic Regression. Before you proceed, sharpen your basics of Regression <u>here</u>.

## Linear Regression takes following assumptions:

- 1. There exists a linear relationship between response and predictor variables
- 2. The predictor (independent) variables are not correlated with each other. Presence of collinearity leads to a phenomenon known as <u>multicollinearity</u>.
- 3. The error terms are uncorrelated. Otherwise, it will lead to <u>autocorrelation</u>.
- 4. Error terms must have constant variance. Non-constant variance leads to <u>heteroskedasticity</u>.

Let's now build out first regression model on this data set. R uses *lm()* function for regression.

```
> linear_model <- lm(Item_Outlet_Sales ~ ., data = new_train)</pre>
```

## > summary(linear\_model)

Adjusted  $R^2$  measures the goodness of fit of a regression model. Higher the  $R^2$ , better is the model. Our  $R^2 = 0.2085$ . It means we really did something drastically wrong. Let's figure it out.

In our case, I could find our new variables aren't helping much i.e. Item count, Outlet Count and Item\_Type\_New. Neither of these variables are significant. Significant variables are denoted by '\*' sign.

As we know, correlated predictor variables brings down the model accuracy. Let's find out the amount of correlation present in our predictor variables. This can be simply calculated using:

```
> cor(new_train)
```

Alternatively, you can also use corrplot package for some fancy correlation plots. Scrolling through the long list of correlation coefficients, I could find a deadly correlation coefficient:

```
cor(new train$Outlet Count, new train$`Outlet Type Grocery Store`)
```

## [1] -0.9991203

Outlet\_Count is highly correlated (negatively) with Outlet Type Grocery Store. Here are some problems I could find in this model:

- 1. We have correlated predictor variables.
- 2. We did one hot encoding and label encoding. That's not necessary since linear regression handle categorical variables by creating dummy variables intrinsically.

3. The new variables (item count, outlet count, item type new) created in feature engineering are not significant.

Let's try to create a more robust regression model. This time, I'll be using a building a simple model without encoding and new features. Below is the entire code:

```
#load directory
> path <- "C:/Users/manish/desktop/Data/February 2016"</pre>
> setwd(path)
#load data
> train <- read.csv("train Big.csv")</pre>
> test <- read.csv("test Big.csv")</pre>
#create a new variable in test file
> test$Item Outlet Sales <- 1</pre>
#combine train and test data
> combi <- rbind(train, test)</pre>
impute missing value in Item Weight
> combi$Item Weight[is.na(combi$Item Weight)] <- median(combi$Item Weight, na.rm = TRUE)</pre>
#impute 0 in item visibility
> combi$Item Visibility <- ifelse(combi$Item Visibility == 0, median(combi$Item Visibility),</pre>
combi$Item Visibility)
```

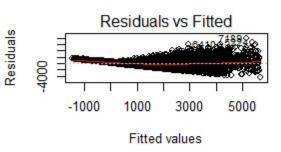
```
##rename level in Outlet Size
> levels(combi$Outlet Size)[1] <- "Other"</pre>
#rename levels of Item Fat Content
> library(plyr)
> combi$Item Fat Content <- revalue(combi$Item Fat Content,c("LF" = "Low Fat", "reg" =</pre>
"Regular"))
> combi$Item Fat Content <- revalue(combi$Item Fat Content, c("low fat" = "Low Fat"))</pre>
#create a new column 2013 - Year
> combi$Year <- 2013 - combi$Outlet Establishment Year</pre>
#drop variables not required in modeling
> library(dplyr)
> combi <- select(combi, -c(Item Identifier, Outlet Identifier, Outlet Establishment Year))</pre>
#divide data set
> new train <- combi[1:nrow(train),]</pre>
> new test <- combi[-(1:nrow(train)),]</pre>
#linear regression
> linear model <- lm(Item Outlet Sales ~ ., data = new train)</pre>
> summary(linear model)
```

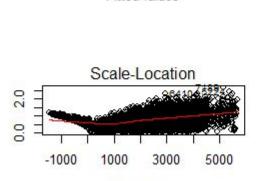
Now we have got  $R^2 = 0.5623$ . This teaches us that, sometimes all you need is simple thought process to get high accuracy. Quite a good improvement from previous model. Next, time when you work on any model, always remember to start with a simple model.

Let's check out regression plot to find out more ways to improve this model.

> par(mfrow=c(2,2))

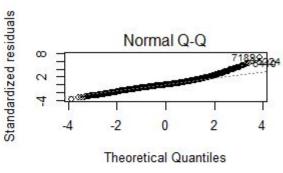
> plot(linear\_model)

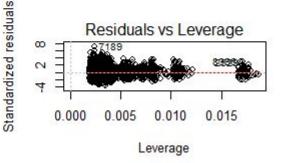




Fitted values

V|Standardized residuals







You can zoom these graphs in R Studio at your end. All these plots have a different story to tell. But the most important story is being portrayed by Residuals vs Fitted graph.

Residual values are the difference between actual and predicted outcome values. Fitted values are the predicted values. If you see carefully, you'll discover it as a funnel shape graph (from right to left). The shape of this graph suggests that our model is suffering from heteroskedasticity (unequal variance in error terms). Had there been constant variance, there would be no pattern visible in this graph.

A common practice to tackle heteroskedasticity is by taking the log of response variable. Let's do it and check if we can get further improvement.

```
> linear_model <- lm(log(Item_Outlet_Sales) ~ ., data = new_train)</pre>
```

> summary(linear\_model)

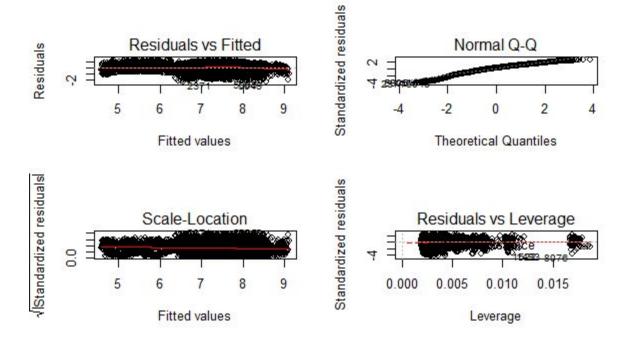


And, here's a snapshot of my model output. Congrats! We have got an improved model with  $R^2 = 0.72$ . Now, we are on the right path.

```
Item_TypeSnack Foods
Item_TypeSoft Drinks
Item_TypeStarchy Foods
Item MRP
                               ***
Outlet_SizeHigh
Outlet SizeMedium
Outlet SizeSmall
                               **
Outlet_Location_TypeTier 2
Outlet_Location_TypeTier 3
Outlet_TypeSupermarket Type1
                               ***
Outlet_TypeSupermarket Type2
                               ***
Outlet_TypeSupermarket Type3
                               ***
                               **
Year
Signif. codes: 0 '***' 0.001 '**' 0.01 '*' 0.05 '.' 0.1 ' '1
Residual standard error: 0.5378 on 8494 degrees of freedom
Multiple R-squared: 0.7214, Adjusted R-squared: 0.7205
F-statistic: 785.4 on 28 and 8494 DF, p-value: < 2.2e-16
```



Once again you can check the residual plots (you might zoom it). You'll find there is no longer a trend in residual vs fitted value plot.





This model can be further improved by detecting outliers and high leverage points. For now, I leave that part to you! I shall write a separate post on mysteries of regression soon. For now, let's check our RMSE so that we can compare it with other algorithms demonstrated below.

To calculate RMSE, we can load a package named *Metrics*.

```
> install.packages("Metrics")
```

```
> library(Metrics)
```

```
> rmse(new_train$Item_Outlet_Sales, exp(linear_model$fitted.values))
```

```
[1] 1140.004
```

Let's proceed to decision tree algorithm and try to improve our RMSE score.



## **Decision Trees**

Before you start, I'd recommend you to glance through the basics of decision tree algorithms. To understand what makes it superior than linear regression, check this tutorial <u>Part 1</u> and <u>Part 2</u>.

In R, decision tree algorithm can be implemented using *rpart package*. In addition, we'll use *caret package* for doing cross validation. Cross validation is a technique to build robust models which are not prone to overfitting. Read more about <u>Cross Validation</u>.

In R, decision tree uses a complexity parameter (*cp*). It measures the tradeoff between model complexity and accuracy on training set. A smaller cp will lead to a bigger tree, which might overfit the model. Conversely, a large cp value might underfit the model. Underfitting occurs when the model does not capture underlying trends properly. Let's find out the optimum cp value for our model with 5 fold cross validation.

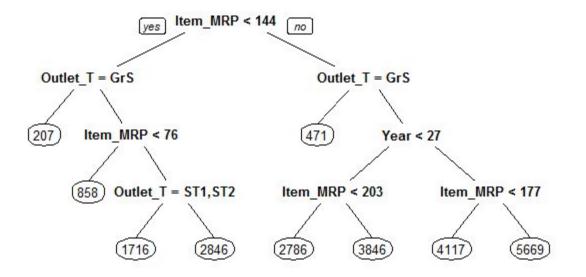


```
#loading required libraries
> library(rpart)
> library(e1071)
> library(rpart.plot)
> library(caret)
#setting the tree control parameters
> fitControl <- trainControl(method = "cv", number = 5)</pre>
> cartGrid <- expand.grid(.cp=(1:50)*0.01)</pre>
#decision tree
> tree_model <- train(Item_Outlet_Sales ~ ., data = new_train, method = "rpart", trControl = fitControl, tuneGrid
= cartGrid)
> print(tree_model)
```

The final value for cp = 0.01. You can also check the table populated in console for more information. The model with cp = 0.01 has the least RMSE. Let's now build a decision tree with 0.01 as complexity parameter.

```
> main_tree <- rpart(Item_Outlet_Sales ~ ., data = new_train, control = rpart.control(cp=0.01))</pre>
```

> prp(main\_tree)





Here is the tree structure of our model. If you have gone through the basics, you would now understand that this algorithm has marked Item\_MRP as the most important variable (being the root node). Let's check the RMSE of this model and see if this is any better than regression.

```
> pre_score <- predict(main_tree, type = "vector")
> rmse(new_train$Item_Outlet_Sales, pre_score)
[1] 1102.774
```

As you can see, our RMSE has further improved from 1140 to 1102.77 with decision tree. To improve this score further, you can further tune the parameters for greater accuracy.



## **Random Forest**

Random Forest is a powerful algorithm which holistically takes care of missing values, outliers and other non-linearities in the data set. It's simply a collection of classification trees, hence the name 'forest'. I'd suggest you to quickly refresh your basics of random forest with this <u>tutorial</u>.

In R, random forest algorithm can be implement using *randomForest* package. Again, we'll use train package for cross validation and finding optimum value of model parameters.

For this problem, I'll focus on two parameters of random forest. *mtry* and *ntree*. ntree is the number of trees to be grown in the forest. *mtry* is the number of variables taken at each node to build a tree. And, we'll do a 5 fold cross validation.



```
#load randomForest library
> library(randomForest)
#set tuning parameters
> control <- trainControl(method = "cv", number = 5)</pre>
#random forest model
> rf_model <- train(Item_Outlet_Sales ~ ., data = new_train, method = "parRF", trControl =</pre>
control, prox = TRUE, allowParallel = TRUE)
#check optimal parameters
> print(rf_model)
```



If you notice, you'll see I've used method = "parRF". This is parallel random forest. This is parallel implementation of random forest. This package causes your local machine to take less time in random forest computation. Alternatively, you can also use method = "rf" as a standard random forest function.

```
Parallel Random Forest
8523 samples
  9 predictor
No pre-processing
Resampling: Cross-Validated (5 fold)
Summary of sample sizes: 6819, 6818, 6818, 6819, 6818
Resampling results across tuning parameters:
       RMSE Rsquared RMSE SD Rsquared SD
 mtry
       1293.032 0.5157854 7.845358 <u>0.01852388</u>
 15 1122.530 0.5697400 16.109357 0.01307624
 28
       1135.780 0.5613481 16.825174 0.01294525
RMSE was used to select the optimal model using the smallest value.
The final value used for the model was mtry = 15.
```



Now we've got the optimal value of mtry = 15. Let's use 1000 trees for computation.

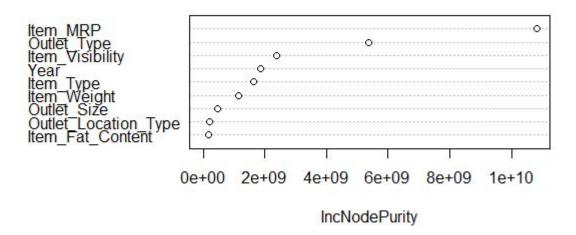
#random forest model

- > forest\_model <- randomForest(Item\_Outlet\_Sales ~ ., data = new\_train, mtry = 15, ntree = 1000)</pre>
- > print(forest\_model)
- > varImpPlot(forest\_model)

This model throws RMSE = 1132.04 which is not an improvement over decision tree model. Random forest has a feature of presenting the important variables. We see that the most important variable is Item\_MRP (also shown by decision tree algorithm).



## forest\_model



This model can be further improved by tuning parameters. But this can also be your output for submission if you want.



```
> main_predict <- predict(main_tree, newdata = new_test, type = "vector")
> sub_file <- data.frame(Item_Identifier = test$Item_Identifier, Outlet_Identifier = test$Outlet_Identifier,
Item_Outlet_Sales = main_predict)
> write.csv(sub_file, 'Decision_tree_sales.csv')
```

When predicted on out of sample data, our RMSE has come out to be 1174.33. Here are some things you can do to improve this model further:

- 1. Since we did not use encoding, I encourage you to use one hot encoding and label encoding for random forest model.
- 2. Parameters Tuning will help.
- 3. Use Gradient Boosting.
- 4. Build an ensemble of these models. Read more about **Ensemble Modeling**.



## Thank You

