



**SRM INSTITUTE OF SCIENCE & TECHNOLOGY,
NCR CAMPUS, MODINAGAR
(FACULTY OF SCIENCE AND HUMANITIES)**

DEPARTMENT OF COMPUTER APPLICATIONS

PRACTICAL FILE

DATA ANALYSIS USING R

(PCA20S02J)

MCA-1st Year, 2nd Semester

Session: Dec 2023 - May 2024

**Submitted to:
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**Submitted by:
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BONAFIDE CERTIFICATE

Certified to be the bonafide record of the work done by **KIRTI GARG** of MCA-First year, Second Semester (Section A) for the award of **Master's** degree course in DEPARTMENT OF COMPUTER APPLICATIONS in the FACULTY OF SCIENCE & HUMANITIES in DATA ANALYSIS USING R [**PCA20S02J**] laboratory during the Academic year- 2023-24.

SUBJECT -INCHARGE

HEAD OF THE DEPARTMENT

Submitted for the university examination held on _____

INTERNAL EXAMINER 1

INTERNAL EXAMINER 2

INDEX

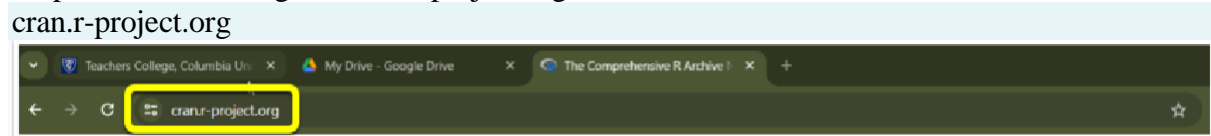
S.N	LIST OF EXPERIMENT Using R	DATE	PAGE NO.	SIGNATURE/ REMARKS
1.	Write steps on how to install R studio.			
2.	Create two vectors in R for numeric data.			
3.	Create a list in a data structure that has components of mixed datatypes.			
4.	Create a code to display Fibonacci series.			
5.	Implement decision tree on credit card issue dataset (import fromkaggale).			
6.	Implement the KNN algorithm on the Brest cancer dataset.			
7.	Implement the Naïve Bayes algorithm on the iris dataset.			
8.	Random forest algorithm implement on iris dataset.			
9.	Implement K-mean clustering inyour own dataset create database using vector.			
10	Implement linear regression on iris dataset.			
11	SVM algorithm using WDBC dataset.			
12	Implement logistics regression on the IRIS dataset.			
13	Implement Apriori algorithm.			

PRACTICAL:1

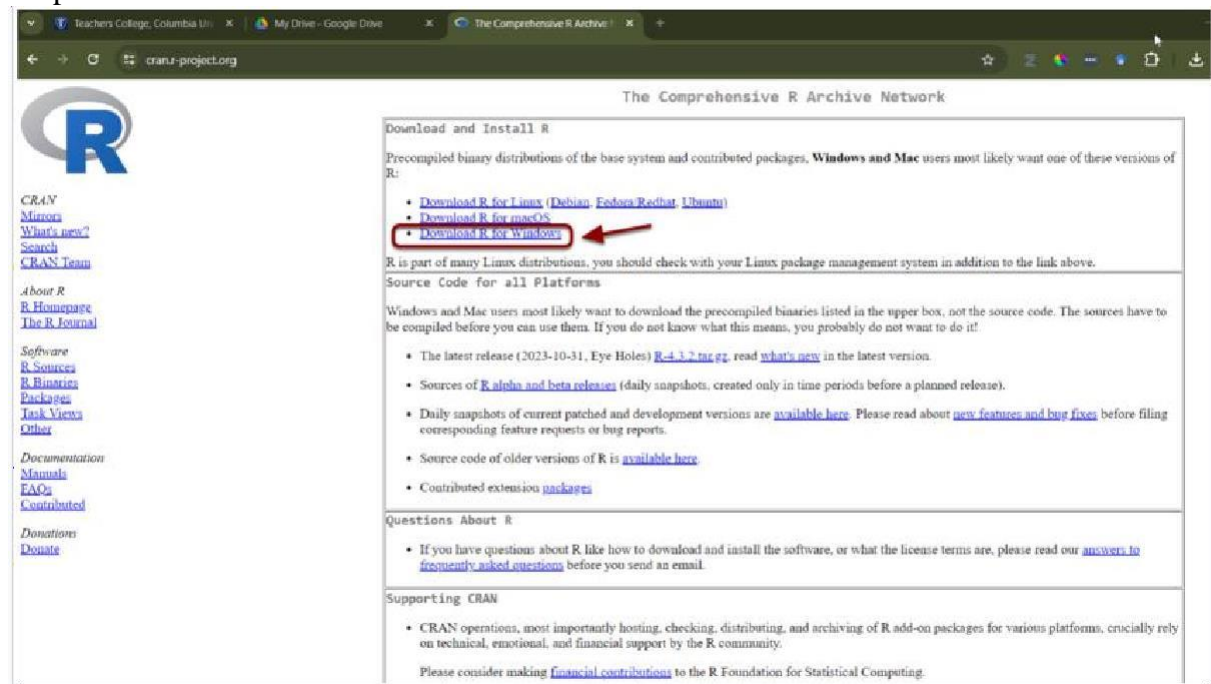
AIM: Write steps of installation of R and R Studio.

Procedure:

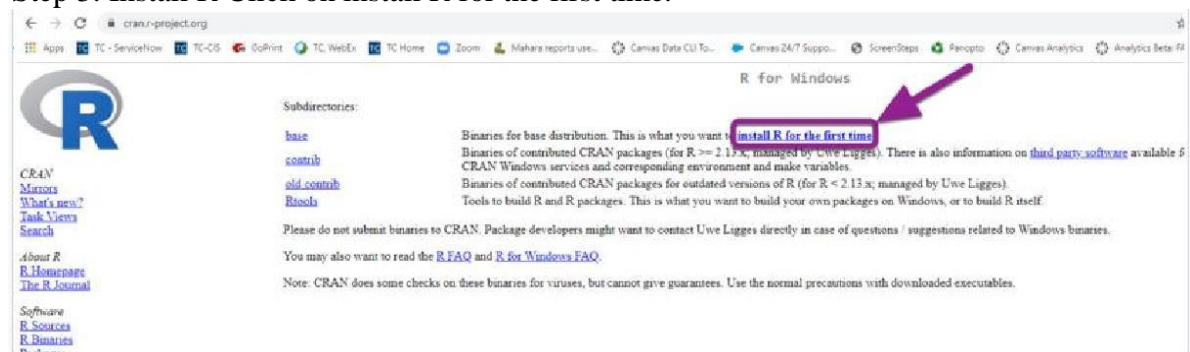
Step 1. To install R, go to cran.r-project.org



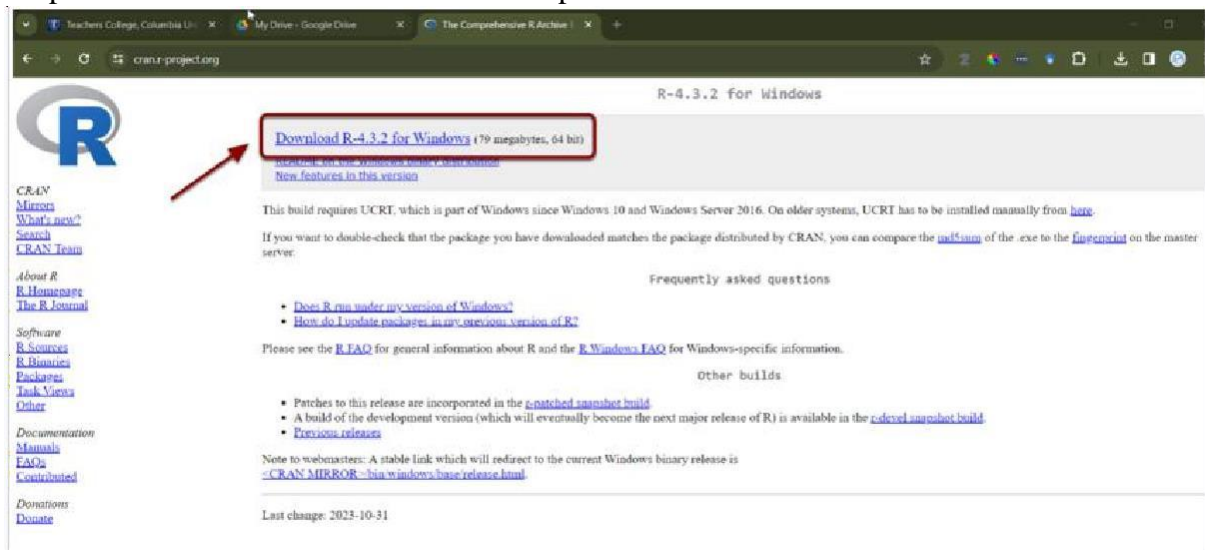
Step 2. Click Download R for Windows.



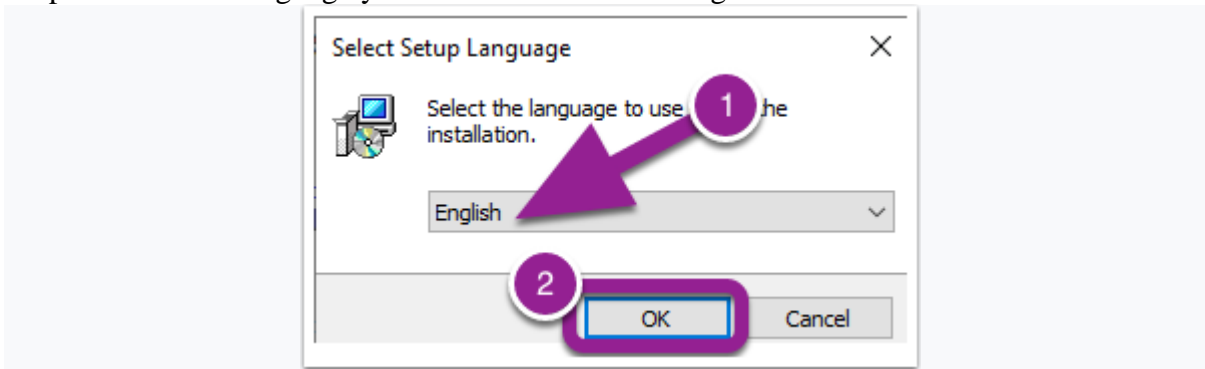
Step 3. Install R Click on install R for the first time.



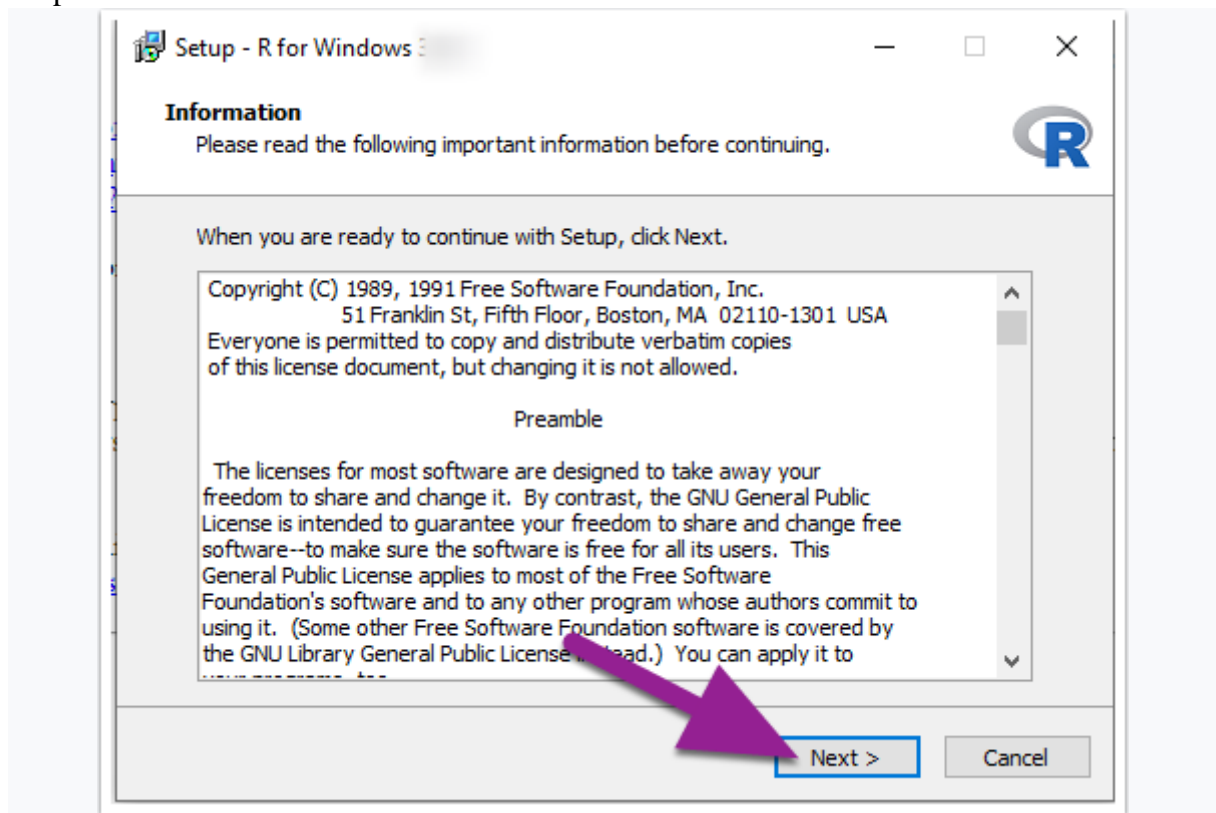
Step 4. Click Download R for Windows. Open the downloaded file.



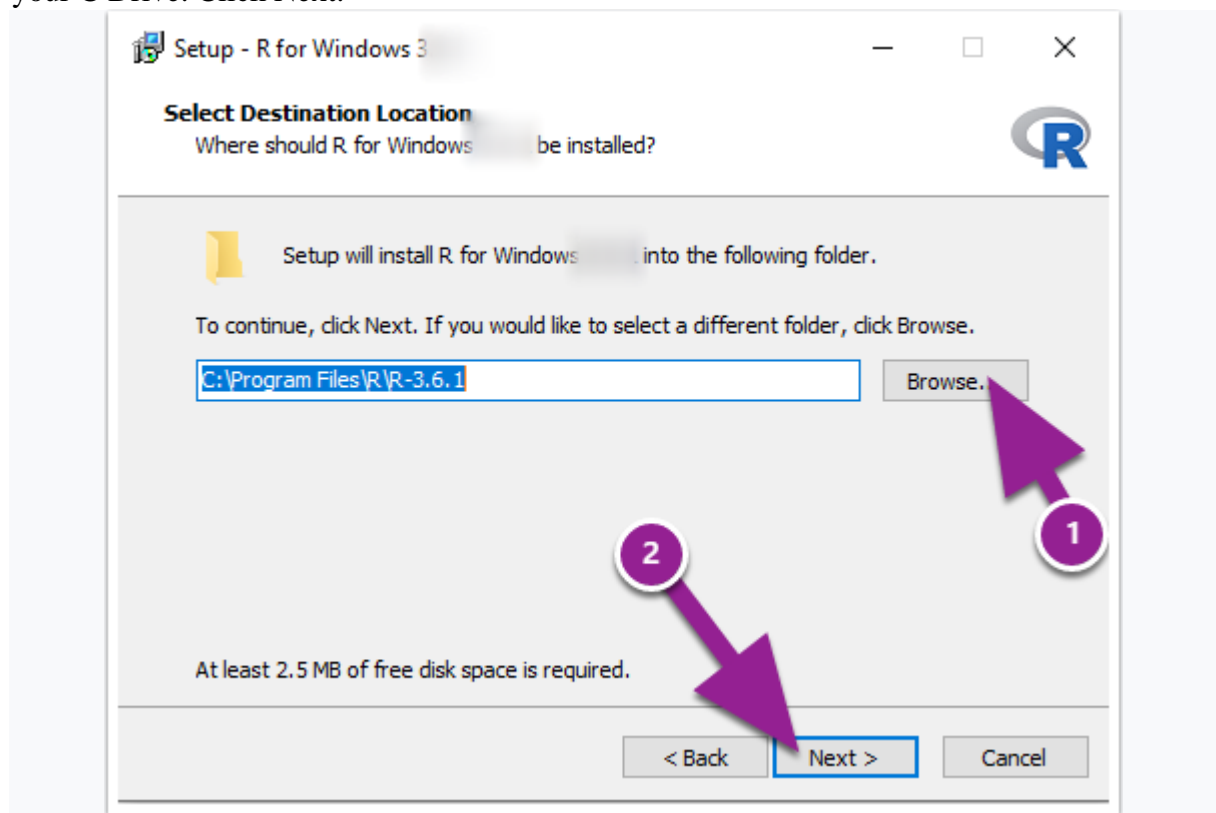
Step 5. Select the language you would like to use during the installation. Then click OK.



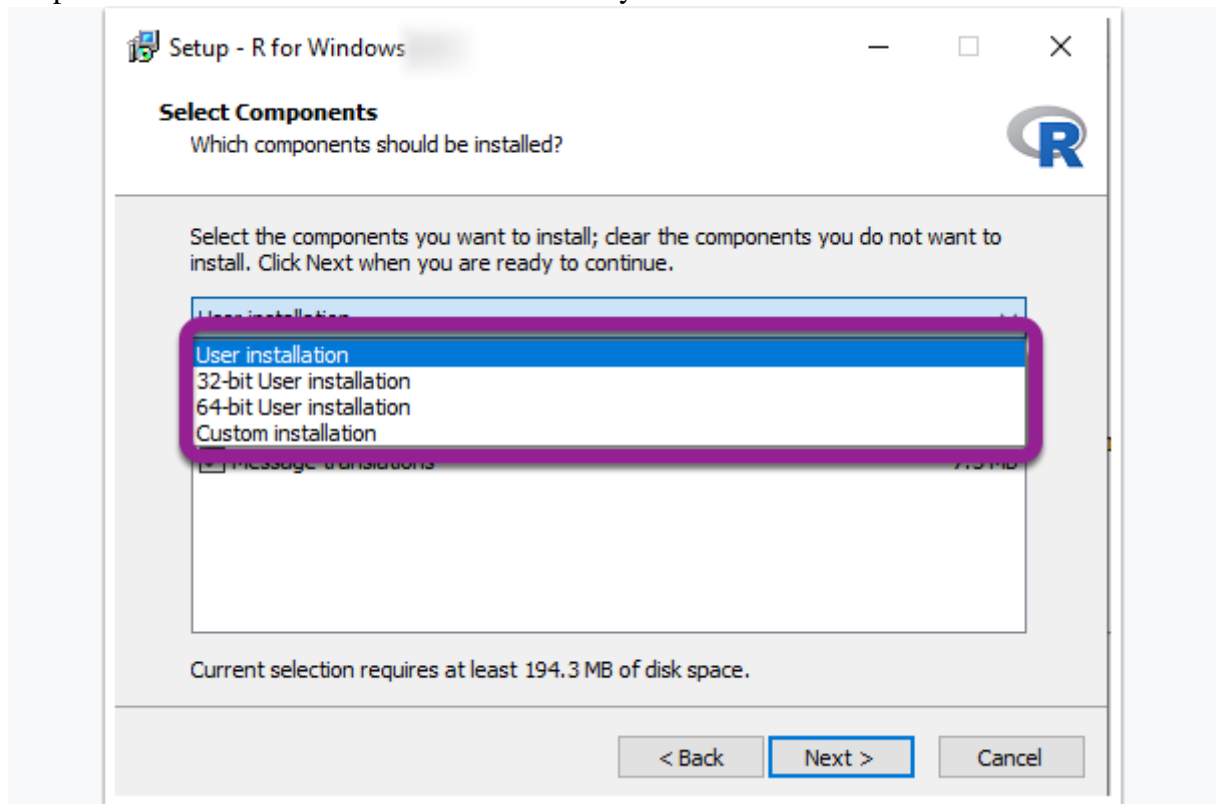
Step 6. Click Next.



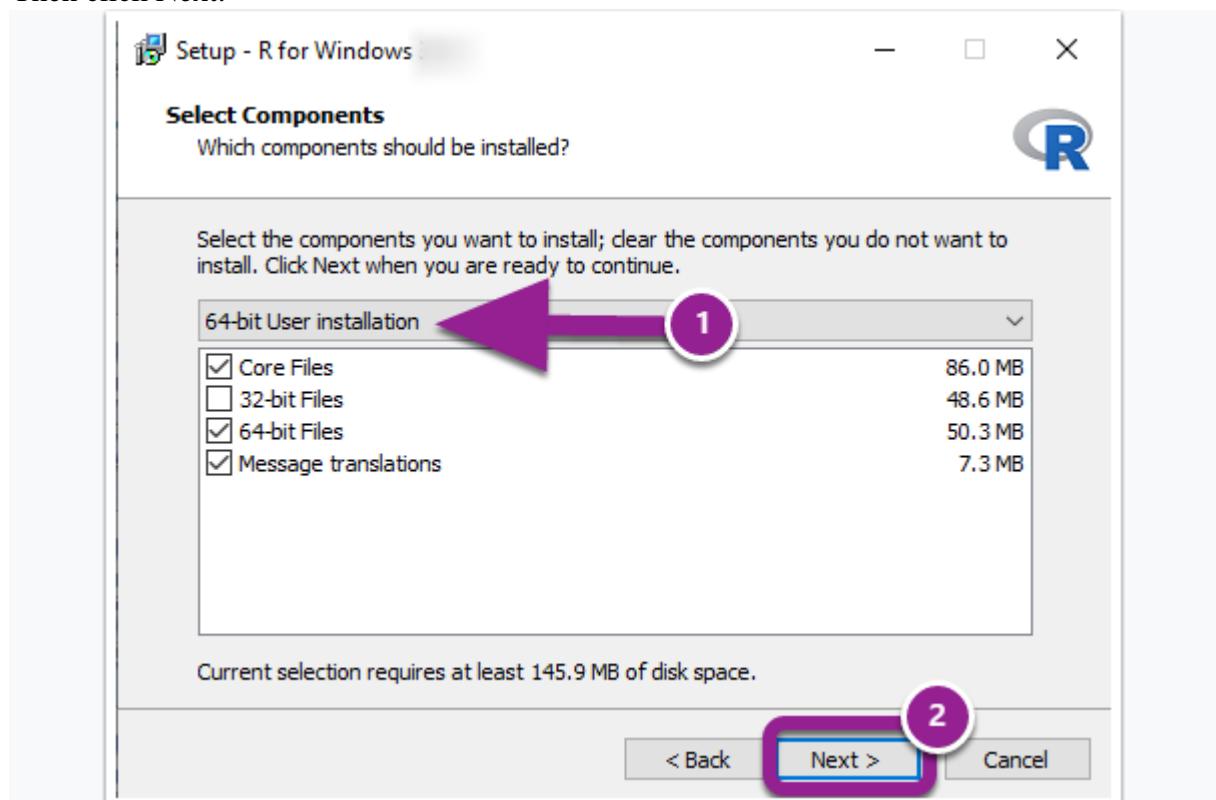
Step 7. Select where you would like R to be installed. It will default to your Program Files on your C Drive. Click Next.



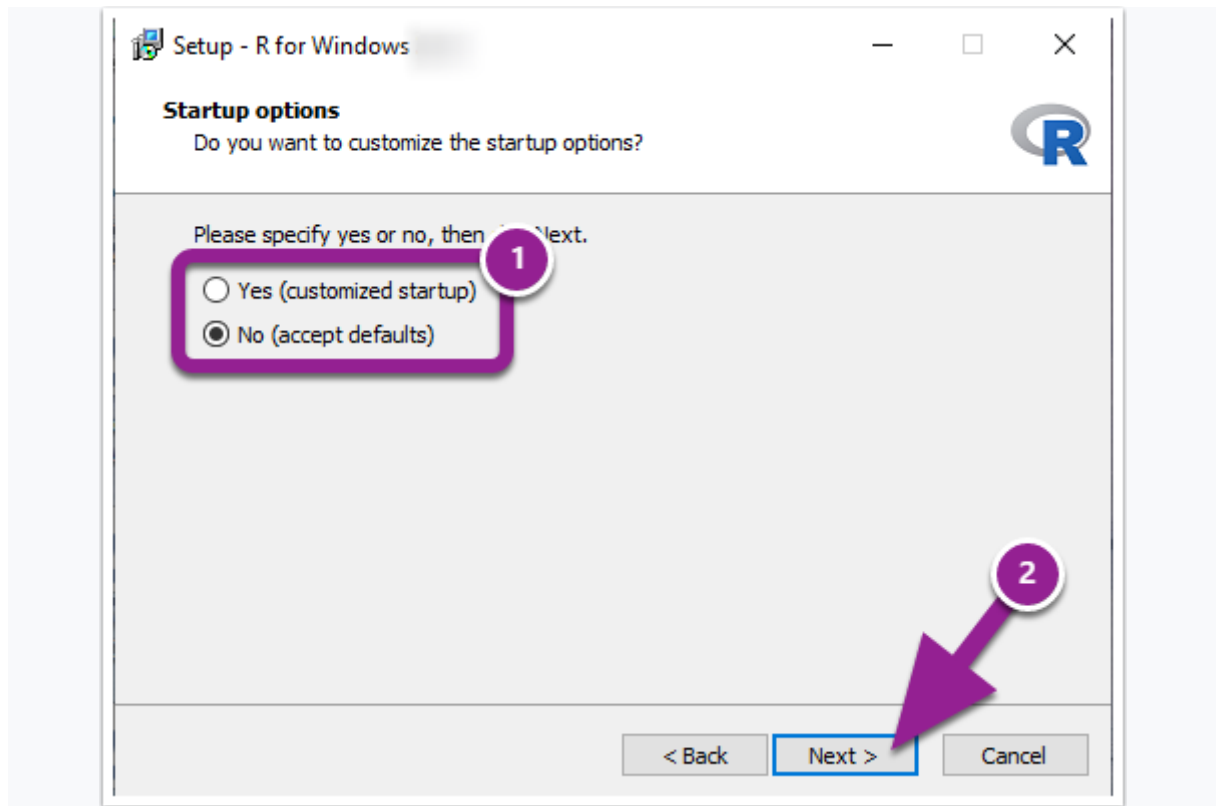
Step 8. You can then choose which installation you would like.



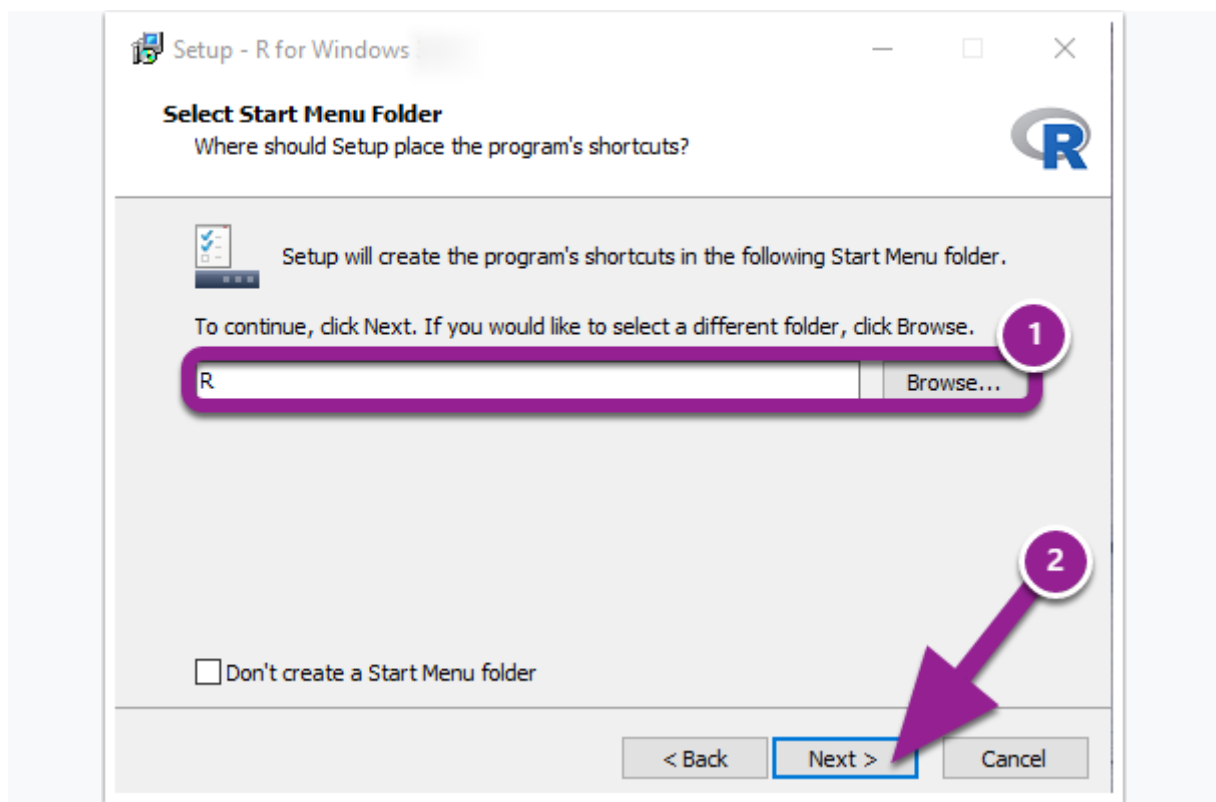
Step 9. (Optional) If your computer is a 64-bit, you can choose the 64-bit User Installation. Then click Next.



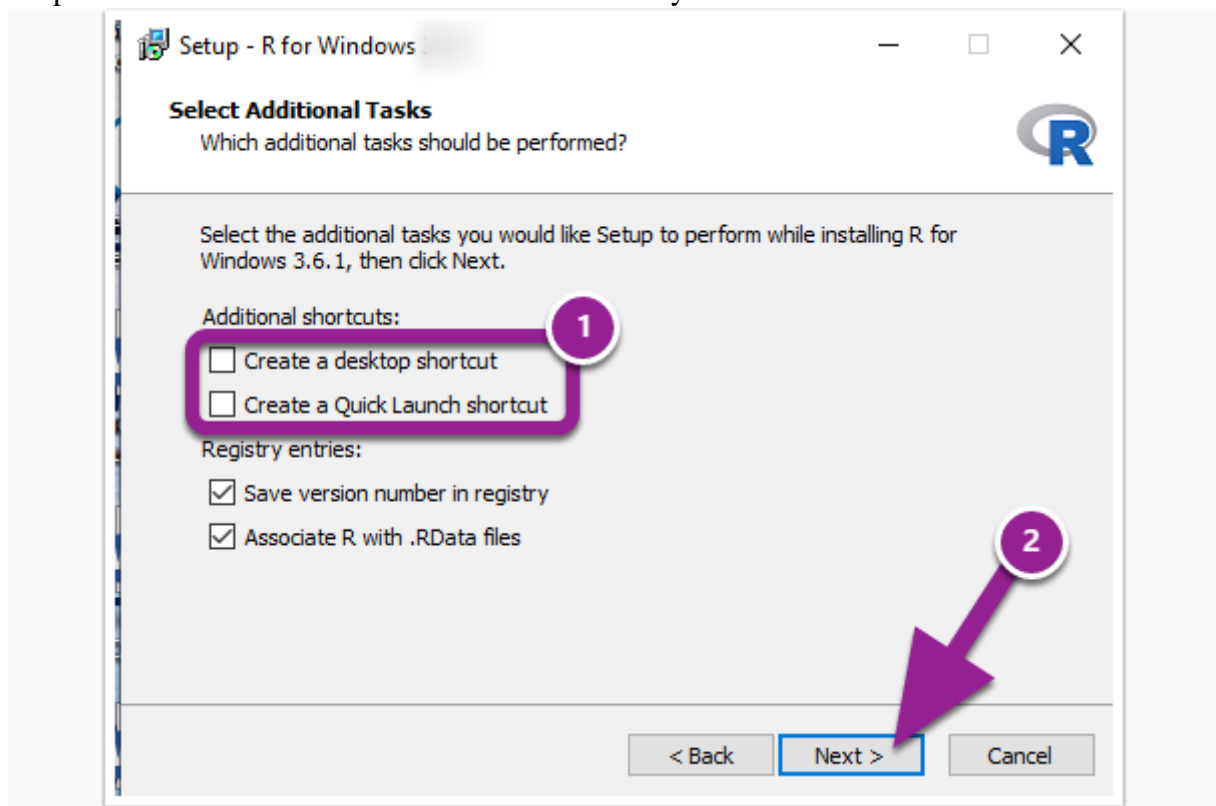
Step 10. Then specify if you want to customized your startup or just use the defaults. Then click Next.



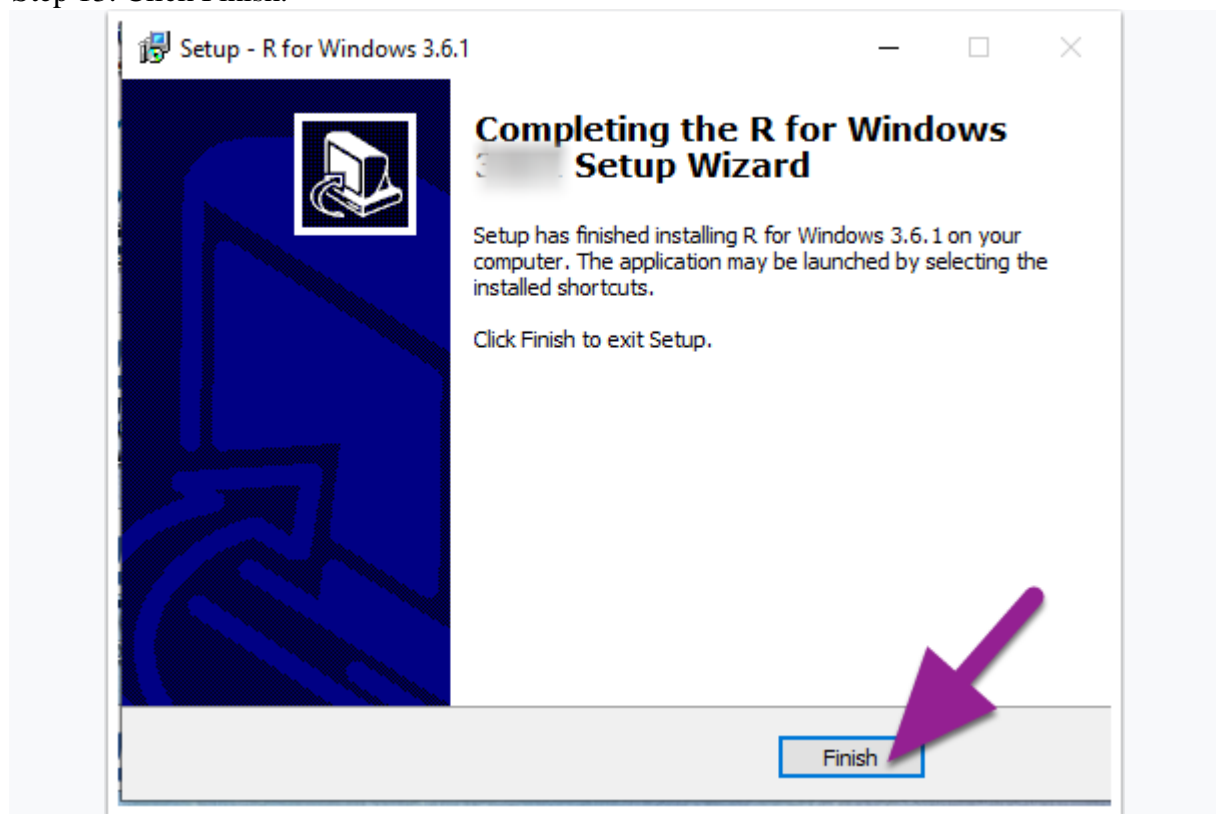
Step 11. Then you can choose the folder that you want R to be saved within or the default if the R folder that was created. Once you have finished, click Next. You can also choose if you do not want a Start Menu folder at the bottom.



Step 12. You can then select additional shortcuts if you would like. Click Next.

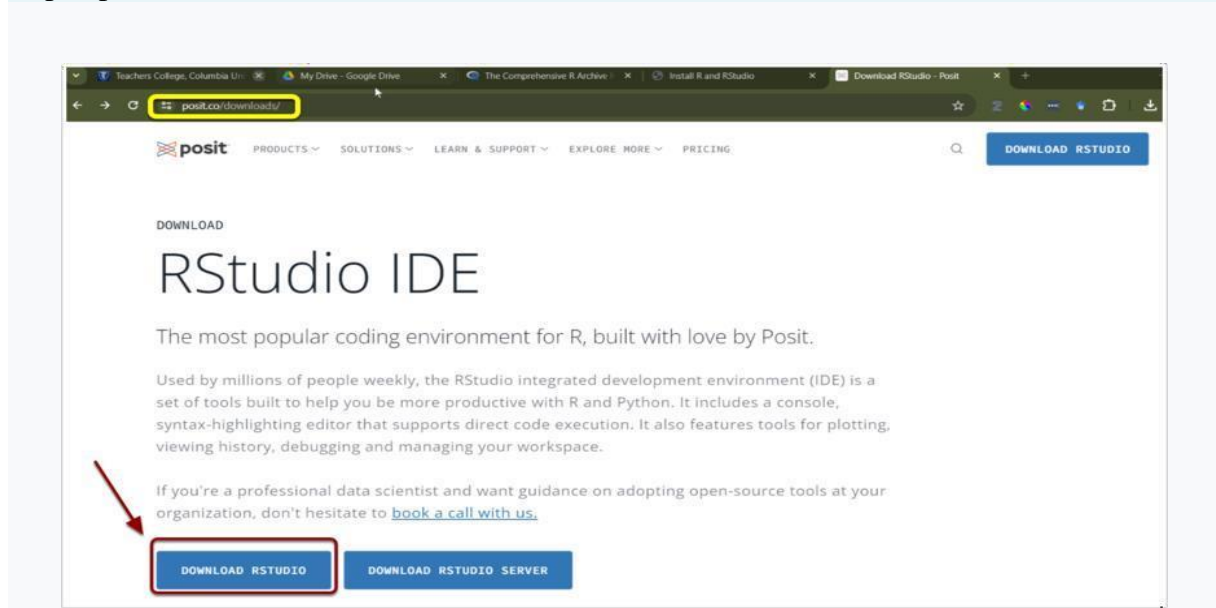


Step 13. Click Finish.

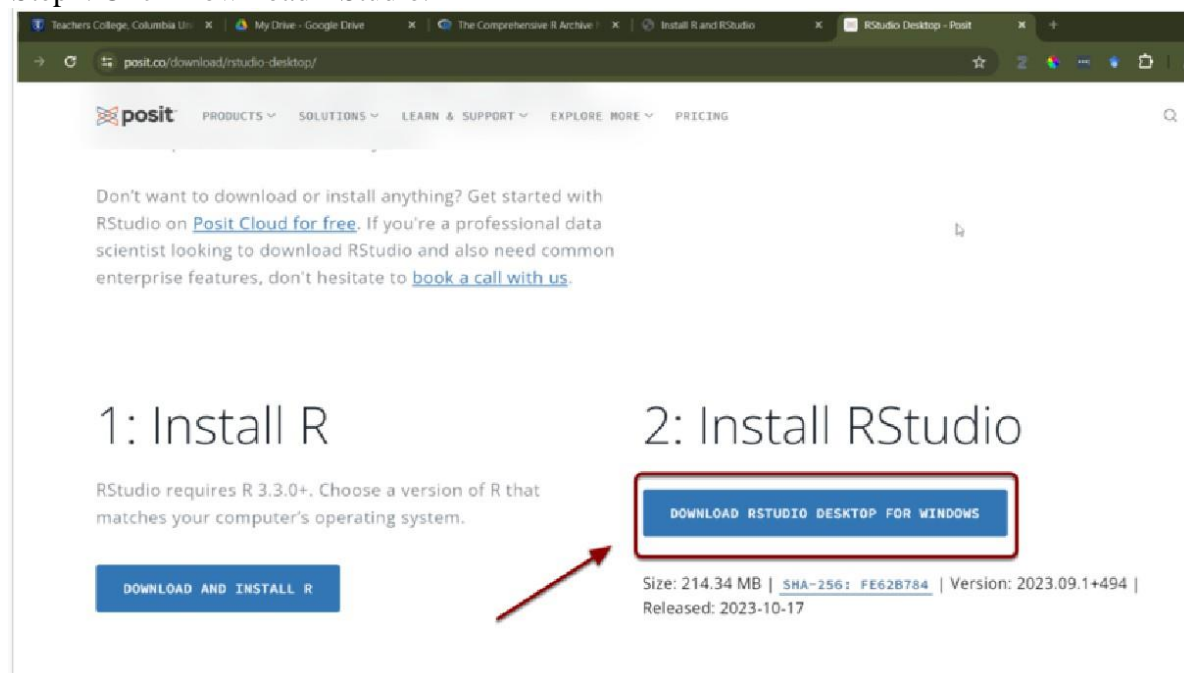


STEPS OF INSTALLATION OF R STUDIO

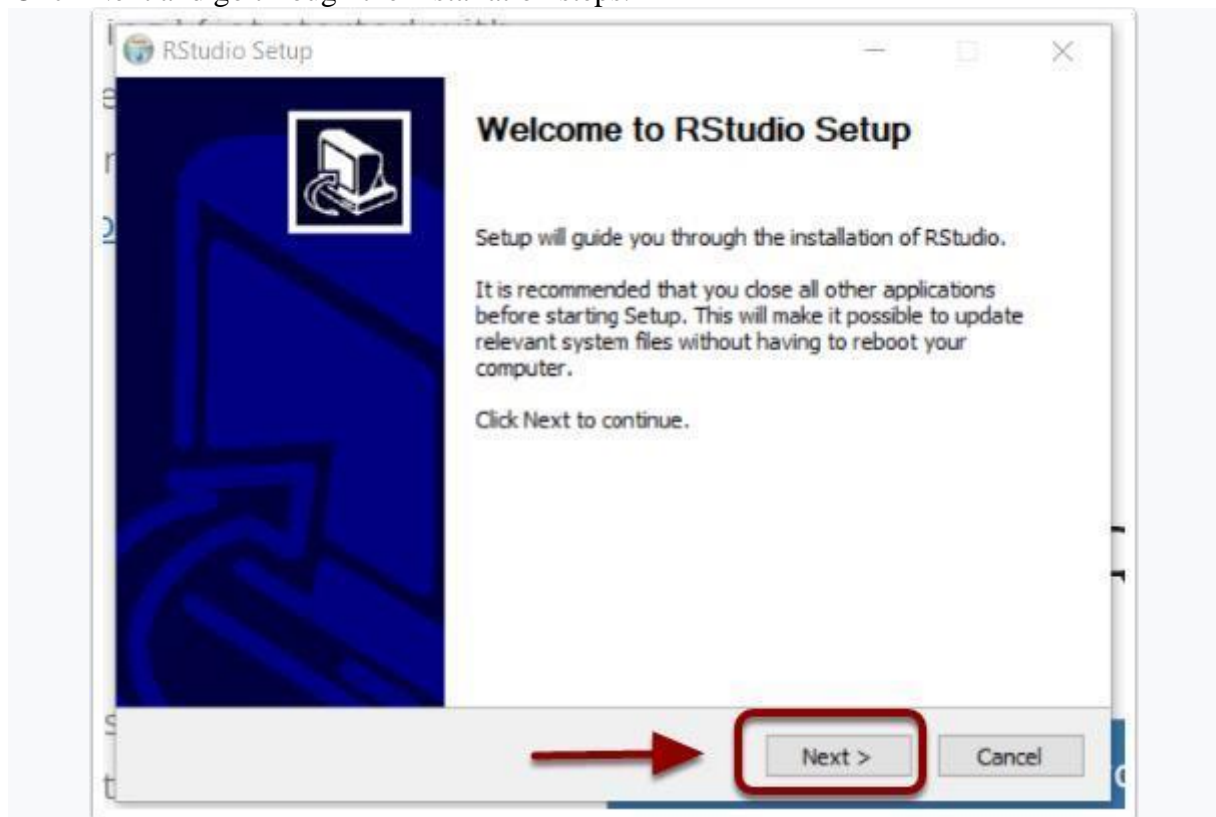
Step1:. Next, download RStudio. Go to <https://posit.co/downloads/>



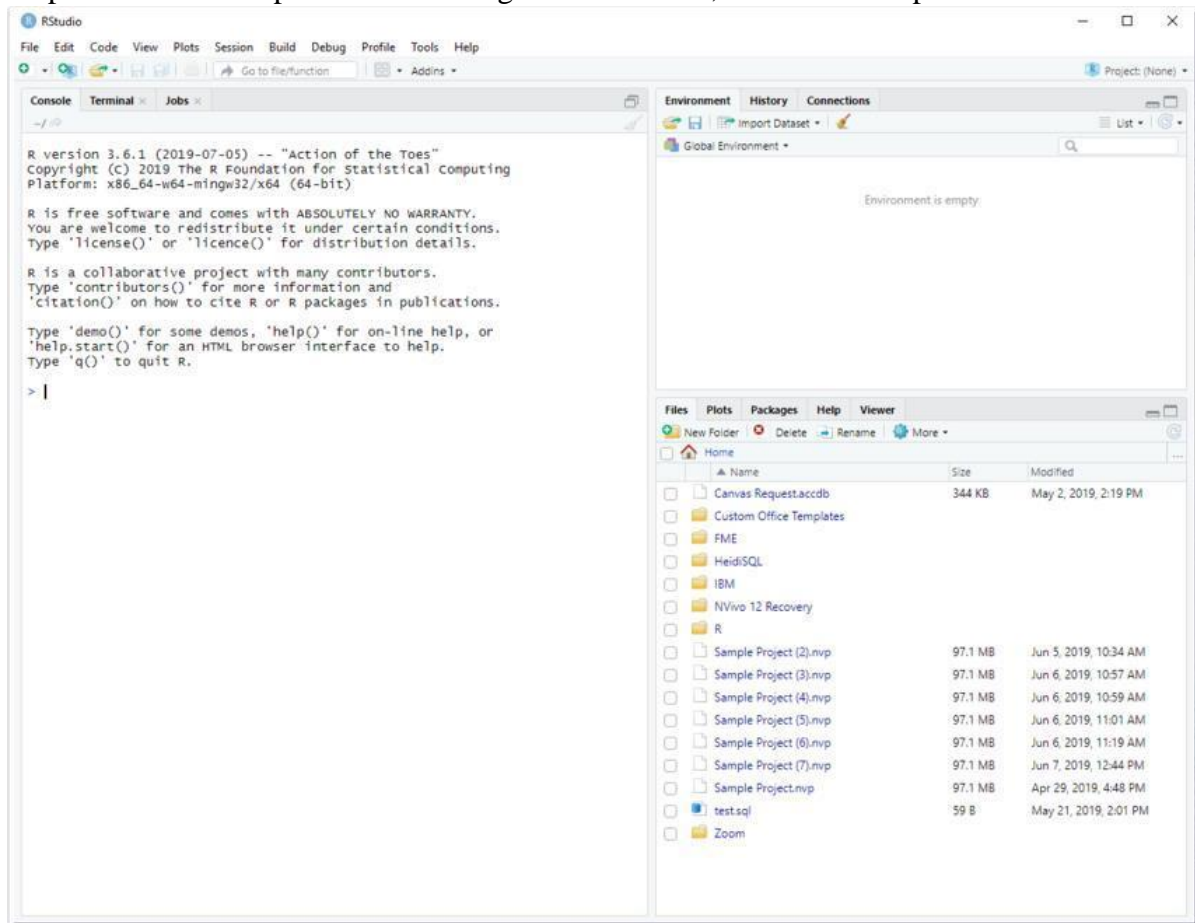
Step2. Click Download RStudio.



Step3. Once the packet has downloaded, the Welcome to RStudio Setup Wizard will open. Click Next and go through the installation steps.




Step4. After the Setup Wizard finishing the installation, RStudio will open.

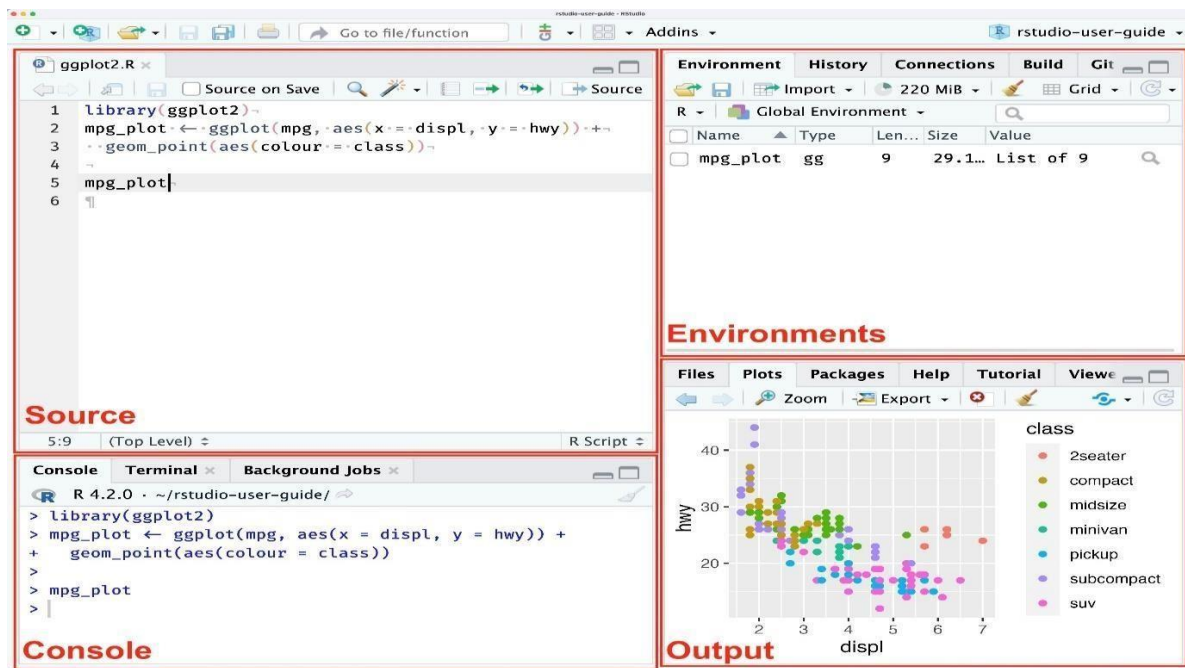


Pane Layout

The RStudio user interface has 4 primary panes:

- **Source pane**
- **Console pane**
- **Environment pane**, containing the **Environment**, **History**, **Connections**, **Build**, **VCS**, and **Tutorial** tabs
- **Output pane**, containing the **Files**, **Plots**, **Packages**, **Help**, **Viewer**, and **Presentation** tabs

Each pane can be minimized or maximized within the column by clicking the minimize/maximize buttons .



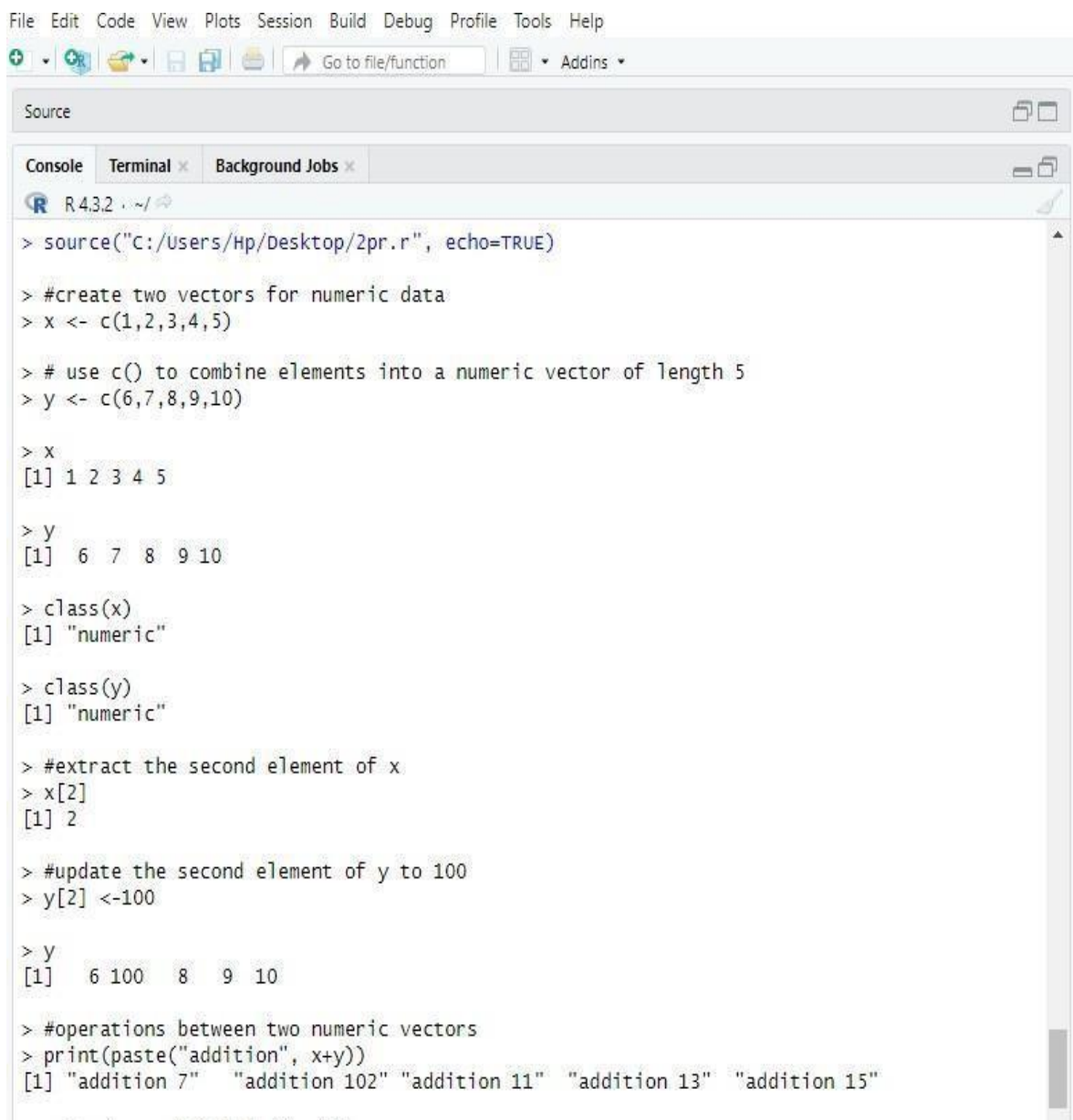
PRACTICAL-2

AIM: Create two vectors in R for Numeric Data.

PROCEDURE:

```
#create two vectors for numeric data
x <- c(1,2,3,4,5)
# use c() to combine elements into a numeric vector of length 5
y <- c(6,7,8,9,10)
x
y
class(x)
class(y)
#extract the second element of x
x[2]
#update the second element of y to 100
y[2] <-100
y
#operations between two numeric vectors
print(paste("addition", x+y))
print(paste("division",y/5))
print(paste("subtraction",y-x)) #create two vectors for numeric data
x <- c(1,2,3,4,5)
# use c() to combine elements into a numeric vector of length 5
y <- c(6,7,8,9,10)
x
y
class(x)
class(y)
#extract the second element of x
x[2]
#update the second element of y to 100
y[2] <-100
y
#operations between two numeric vectors
print(paste("addition", x+y))
```

INPUT/OUTPUT:



The screenshot shows the RStudio interface with the following components:

- Menu Bar:** File, Edit, Code, View, Plots, Session, Build, Debug, Profile, Tools, Help.
- Toolbar:** Includes icons for saving, running, and other standard IDE functions, along with a search bar labeled "Go to file/function" and a dropdown menu labeled "Addins".
- Source Panel:** Displays the R script being executed. The script content is as follows:

```
> source("C:/Users/Hp/Desktop/2pr.r", echo=TRUE)

> #create two vectors for numeric data
> x <- c(1,2,3,4,5)

> # use c() to combine elements into a numeric vector of length 5
> y <- c(6,7,8,9,10)

> x
[1] 1 2 3 4 5

> y
[1] 6 7 8 9 10

> class(x)
[1] "numeric"

> class(y)
[1] "numeric"

> #extract the second element of x
> x[2]
[1] 2

> #update the second element of y to 100
> y[2] <-100

> y
[1] 6 100 8 9 10

> #operations between two numeric vectors
> print(paste("addition", x+y))
[1] "addition 7" "addition 102" "addition 11" "addition 13" "addition 15"
```
- Console Panel:** Shows the output of the script execution, which matches the printed output in the source panel.
- Terminal/Background Jobs Panel:** Currently empty.

PRACTICAL-3

AIM: Create a list in a data structure that has components of mixed data types.

PROCEDURE:

#a vector having elements of different type is called list.

#Use the list() function to create a list.

```
x <- list(1, "first no.", 22, 2, "true")
```

```
print(x)
```

```
print(paste("type of x", class(x)))
```

```
y<-list("z"=1:10)
```

```
print(y)
```

```
print(paste("length of list y", length(y)))
```

#Multiple lists can be merged

```
list_combined <- c(x, y)
```

#create an empty list of a prespecified length with the vector() function

```
z<- vector("list", length=4 )
```

```
print(z)
```

INPUT/OUTPUT:


```
Source
Console Terminal x Background Jobs x
R 4.3.2 ~ /
> source("~/active-rstudio-document", echo=TRUE)

> #Create a list in a data structure that has components of mixed data types.
> #a vector having elements of different type is called list.
> #Use the .... [TRUNCATED]

> print(x)
[[1]]
[1] 1

[[2]]
[1] "first no."

[[3]]
[1] 22

[[4]]
[1] 2

[[5]]
[1] "true"

> print(paste("type of x", class(x)))
[1] "type of x list"

> y<-list("z"=1:10)

> print(y)
$z
[1] 1 2 3 4 5 6 7 8 9 10

> print(paste("length of list y", length(y)))
[1] "length of list y 1"
```

```
Console Terminal x Background Jobs x
R 4.3.2 ~ /
> #Multiple lists can be merged
> list_combined <- c(x, y)

> #create an empty list of a prespecified length with the vector() function
> z<- vector("list", length=4 )

> print(z)
[[1]]
NULL

[[2]]
NULL

[[3]]
NULL

[[4]]
NULL

> |
```

PRACTICAL-4

AIM: Create a code to display a Fibonacci series

PROCEDURE:

```
x=readline("enter a no")
```

```
x=as.integer(x)
```

```
a=0
```

```
b=1
```

```
c=0
```

```
while(x>0){
```

```
  print(c)
```

```
  c=a+b
```

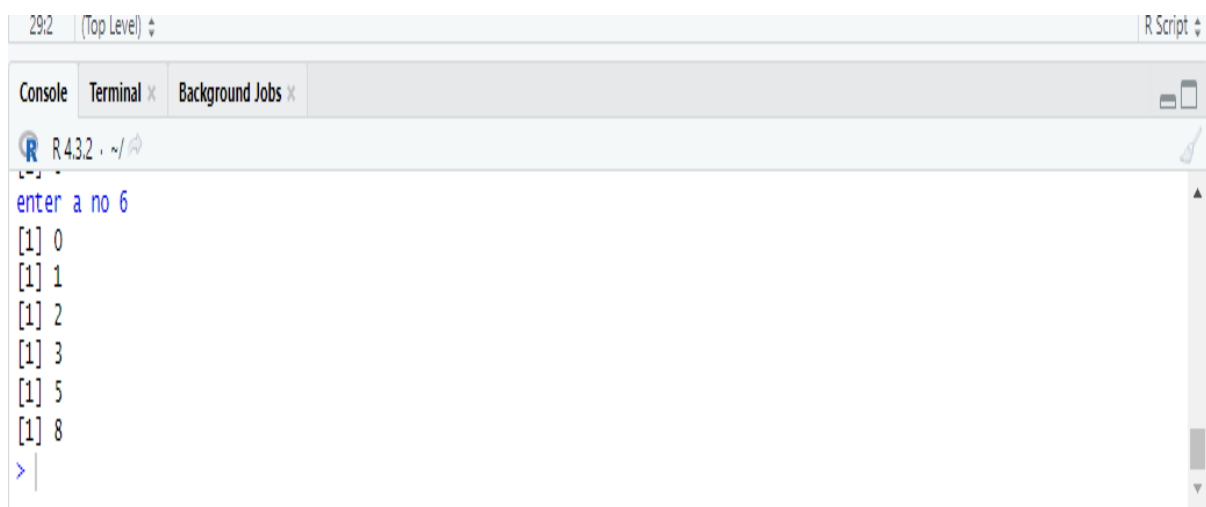
```
  a=b
```

```
  b=c
```

```
  x=x-1
```

```
}
```

OUTPUT



```
29/2 (Top Level) R Script
Console Terminal Background Jobs
R 4.3.2 ~
enter a no 6
[1] 0
[1] 1
[1] 2
[1] 3
[1] 5
[1] 8
> |
```

PRACTICAL-5

AIM: Implement decision tree on credit card issue dataset(import from kaggle).

PROCEDURE:

Algorithm :

1. Import the dataset: Import the dataset which contains the input variables (or the independent variables) and the output variable (or the dependent variable).
2. Data Pre-processing: Check for missing values and handle them, handle outliers if any, and remove or transform variables that do not contribute to the model.
3. Split the dataset: Split the dataset into a training set and a testing set. The training set is used to train the model, while the testing set is used to evaluate the performance of the model.
4. Feature Scaling: If the input variables (or the independent variables) have different ranges, scale them to have the same range to ensure that no variable dominates the others in the modeling process.
5. Model Training: Train the decision tree model using the training set. The decision tree model recursively partitions the input space into subsets, each defined by a set of conditions on the input variables. The partitioning is done in a way that the subsets are as homogeneous as possible with respect to the output variable.
6. Model Evaluation: Evaluate the performance of the model using the testing set. There are several metrics to evaluate the performance of the model, such as accuracy, precision, recall, F1 score, and area under the ROC curve.
7. Model Improvement: If the model does not perform well, try to improve the model by changing the hyperparameters, selecting the relevant features, or using other modeling techniques.
8. Model Deployment: Once the model is finalized and performs well on the testing set, deploy it to the production environment for making predictions on new data.

```

install.packages("party")
install.packages("Rtools")
library(party)
#suffling iris dataset
iris_1=iris[sample(150),]
#spliting dataset into training and testing
train=iris_1[1:100,]
test=iris_1[101:150,]
tree=ctree(Species~Petal.Length+Petal.Width,data =train)
plot(tree)
p=predict(tree,test)
p
table(p,test$ Species)
accuracy=(21+16+10)/50*100
accuracy

```

OUTPUT :

```

> library(party)
> #suffling iris dataset
> iris_1=iris[sample(150),]

> #spliting dataset into training and testing
> train=iris_1[1:100,]
> test=iris_1[101:150,]

> tree=ctree(Species~Petal.Length+Petal.width,data =train)
> plot(tree)
> p=predict(tree,test)

> p
 [1] versicolor setosa    versicolor setosa
 [5] setosa      setosa    virginica  setosa
 [9] versicolor virginica versicolor virginica
[13] virginica  setosa    virginica versicolor
[17] virginica  setosa    virginica setosa
[21] versicolor virginica versicolor setosa
[25] setosa     setosa    setosa    virginica
[29] versicolor virginica setosa    versicolor
[33] setosa     virginica versicolor virginica
[37] versicolor setosa    setosa    virginica
[41] setosa     virginica versicolor versicolor
[45] setosa     virginica virginica  versicolor
[49] versicolor virginica

Levels: setosa versicolor virginica

> table(p,test$ Species)

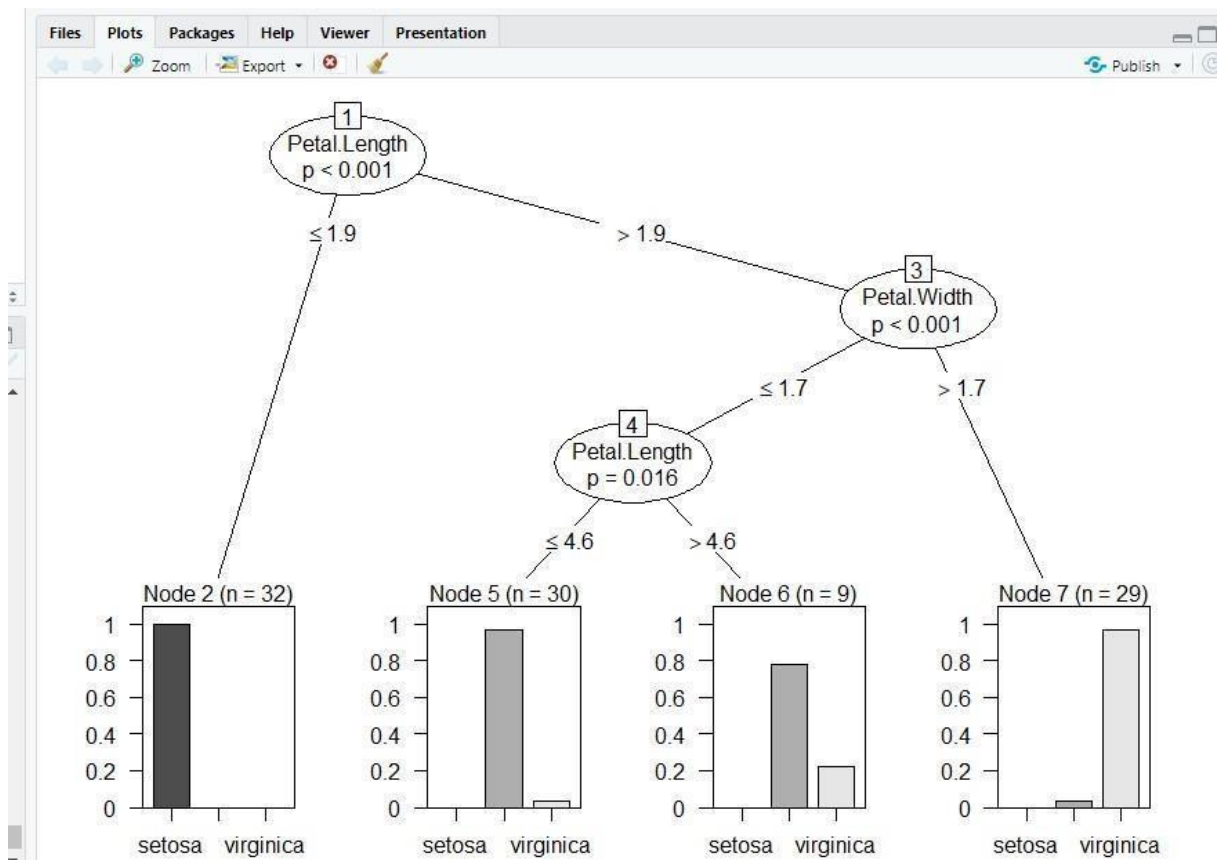
p          setosa versicolor virginica
setosa      18          0          0
versicolor   0         13          2
virginica    0          0         17

> accuracy=(21+16+10)/50*100

> accuracy
[1] 94
>

```

Data	
iris_1	150 obs. of 5 variables
test	50 obs. of 5 variables
train	100 obs. of 5 variables
tree	Formal class BinaryTree



RESULT:

The Decision tree have been implemented successfully.

PRACTICA-6

AIM:- Implement the KNN algorithm on the Breast cancer dataset.

PROCEDURE:-

To implement the k-Nearest Neighbors (KNN) algorithm on the Breast Cancer dataset, you can use the caret and class packages in R. The Breast Cancer dataset is often available through the datasets package in R. Here's an example.

```
# Install and load necessary libraries
install.packages("caret") install.packages("class")
library(caret)
library(class)

# Load the Breast Cancer dataset data("BreastCancer")

# Explore the structure of the dataset str(BreastCancer)

# Split the data into training and testing sets set.seed(123) # Set
seed for reproducibility
sample_index <- createDataPartition(BreastCancer$Class, p = 0.8, list = FALSE) train_data <-
BreastCancer[sample_index, ] test_data <- BreastCancer[-sample_index, ]

# Preprocess the data
# In this example, we'll scale the features
preprocess_params <- preProcess(train_data[, -1], method = c("center", "scale"))

train_data_scaled <- predict(preprocess_params, train_data[, -1]) test_data_scaled <-
predict(preprocess_params, test_data[, -1])

# Train the KNN model
knn_model <- knn(train_data_scaled, test_data_scaled, train_data$Class, k = 5)

# Evaluate the model
conf_matrix <- table(knn_model, test_data$Class)
print("Confusion Matrix:")
print(conf_matrix)

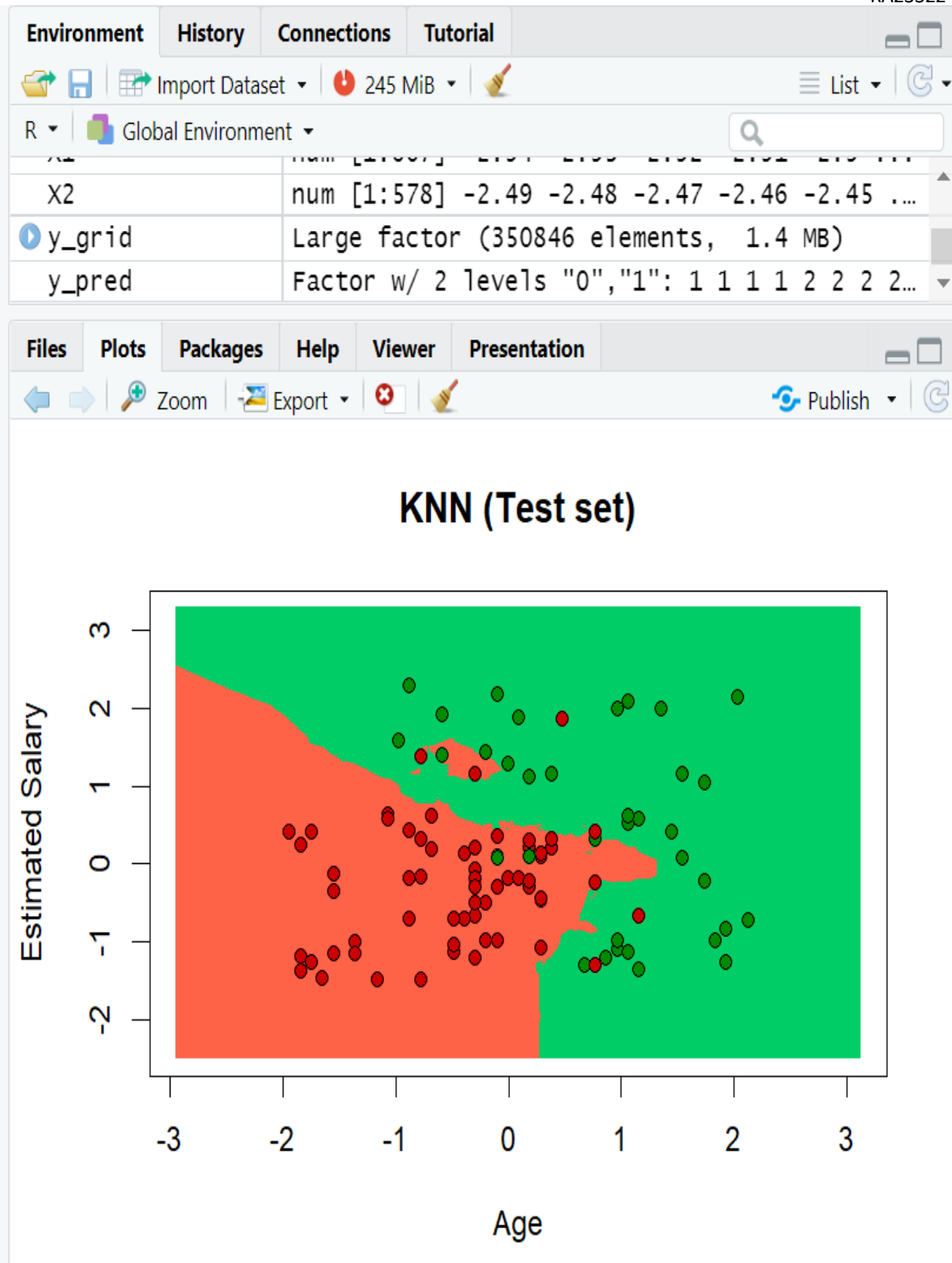
accuracy <- sum(diag(conf_matrix)) / sum(conf_matrix) print(paste("Accuracy:",
round(accuracy, 4)))
```

In this example, the Breast Cancer dataset is split into training and testing sets, and the features are scaled. The KNN model is then trained using the k n function from the class package, and the accuracy of the model is evaluated.

Remember to replace Breast Cancer \$Class with the actual column name representing the target variable in your dataset. Additionally, consider experimenting with different values of k to find the optimal number of neighbors for your specific dataset.

OUTPUT:-

```
> dataset = read.csv('social.csv')
> dataset = dataset[3:5]
>
> # Encoding the label
> dataset$Purchased = factor(dataset$Purchased, levels = c(0, 1))
>
> # Splitting the dataset
> library(caTools)
> split = sample.split(dataset$Purchased, SplitRatio = 0.75)
> training_set = subset(dataset, split == TRUE)
> test_set = subset(dataset, split == FALSE)
>
> # Feature Scaling
> training_set[, 1:2] = scale(training_set[, 1:2])
> test_set[, 1:2] = scale(test_set[, 1:2])
>
> # Fitting and predicting the classifier
> library(class)
> y_pred = knn(train = training_set[, 1:2],
+             test = test_set[, 1:2],
+             cl = training_set[, 3],
+             k = 5)
>
> # Creating the confusion matrix
> cm = table(test_set[, 3], y_pred)
>
> set = training_set
> X1 = seq(min(set[, 1]) - 1, max(set[, 1]) + 1, by = 0.01)
> X2 = seq(min(set[, 2]) - 1, max(set[, 2]) + 1, by = 0.01)
> grid_set = expand.grid(X1, X2)
> colnames(grid_set) = c('Age', 'Estimated Salary')
> y_grid = knn(train = training_set[, 1:2],
+             test = grid_set,
+             cl = training_set[, 3],
+             k = 5) # - means removing the column
> plot(set[, -3],
+      main = 'KNN (Training set)',
+      xlab = 'Age', ylab = 'Estimated Salary',
+      xlim = range(X1), ylim = range(X2))
> contour(X1, X2, matrix(as.numeric(y_grid), length(X1), length(X2)), add = TRUE)
> points(grid_set, pch = '.', col = ifelse(y_grid == 1, 'springgreen3', 'tomato'))
> points(set, pch = 21, bg = ifelse(set[, 3] == 1, 'green4', 'red3'))
>
> set = test_set
> X1 = seq(min(set[, 1]) - 1, max(set[, 1]) + 1, by = 0.01)
> X2 = seq(min(set[, 2]) - 1, max(set[, 2]) + 1, by = 0.01)
> grid_set = expand.grid(X1, X2)
> colnames(grid_set) = c('Age', 'Estimated Salary')
> y_grid = knn(train = training_set[, 1:2],
+             test = grid_set,
+             cl = training_set[, 3],
+             k = 5) # - means removing the column
> plot(set[, -3],
+      main = 'KNN (Test set)',
+      xlab = 'Age', ylab = 'Estimated Salary',
+      xlim = range(X1), ylim = range(X2))
> contour(X1, X2, matrix(as.numeric(y_grid), length(X1), length(X2)), add = TRUE)
> points(grid_set, pch = '.', col = ifelse(y_grid == 1, 'springgreen3', 'tomato'))
> points(set, pch = 21, bg = ifelse(set[, 3] == 1, 'green4', 'red3'))
```

**RESULT:-**

The code has been executed successfully.

PRACTICAL :7

AIM: Implement Naive Bayes algorithm.

PROCEDURE:

1. Convert the given dataset into frequency tables.
2. Generate Likelihood table by finding the probabilities of given features.
3. Now, use Bayes theorem to calculate the posterior probability.

CODE:

```
library(datasets)

data(iris)

install.packages("caret")

library(caret)

set.seed(42)

trainindex <- createDataPartition(iris$Species, p = 0.8, list = FALSE)

train <- iris[trainindex,]

test <- iris[-trainindex,]

library(e1071)

model <- naiveBayes(Species~ ., data = train)

predictions <- predict(model, newdata = test)

library(caret)
```

OUTPUT:

```
> library(caret)
Loading required package: ggplot2
Loading required package: lattice
> set.seed(42)
> trainindex <- createDataPartition(iris$Species, p = 0.8, list = FALSE)
> train <- iris[trainindex,]
> test <- iris[-trainindex,]
> library(e1071)
> model <- naiveBayes(Species~ ., data = train)
> predictions <- predict(model, newdata = test)
> library(caret)
> confusionMatrix(predictions, test$Species)
Confusion Matrix and Statistics
```

Confusion Matrix and Statistics

	Reference		
Prediction	setosa	versicolor	virginica
setosa	10	0	0
versicolor	0	10	2
virginica	0	0	8

Overall Statistics

Accuracy : 0.9333
95% CI : (0.7793, 0.9918)
No Information Rate : 0.3333
P-value [Acc > NIR] : 8.747e-12

Kappa : 0.9

McNemar's Test P-Value : NA

Statistics by Class:

	class: setosa	class: versicolor	class: virginica
Sensitivity	1.0000	1.0000	0.8000
Specificity	1.0000	0.9000	1.0000
Pos Pred Value	1.0000	0.8333	1.0000
Neg Pred value	1.0000	1.0000	0.9091
Prevalence	0.3333	0.3333	0.3333
Detection Rate	0.3333	0.3333	0.2667
Detection Prevalence	0.3333	0.4000	0.2667
Balanced Accuracy	1.0000	0.9500	0.9000

➤

Environment

History

Connections

Tutorial

Import Dataset

367 MiB

List

R

Global Environment

Q

Data

iris	150 obs. of 5 variables
model	List of 5
test	30 obs. of 5 variables
train	120 obs. of 5 variables
trainindex	int [1:120, 1] 1 2 3 4 5 7 8 10 11 12 ...

values

predictions	Factor w/ 3 levels "setosa","versicolor",...: 1 1 1 1 1 1 1 1 1 1 ...
-------------	---

RESULT:

The Naïve Bayes Algorithm has been implemented successfully.

PRACTICAL 8

AIM- Implement Random Forrest Algorithm.

PROCEDURE:

Step-1: Select random K data points from the training set.

Step-2: Build the decision trees associated with the selected data points (Subsets).

Step-3: Choose the number N for decision trees that you want to build.

Step-4: Repeat Step 1 & 2.

Step-5: For new data points, find the predictions of each decision tree, and assign the new data points to the category that wins the majority votes.

CODE:

```
install.packages("randomForest")

library(randomForest)

iris_1=iris[sample(150),]

View(iris_1)

train=iris_1[1:100,]

test=iris_1[101:150,]

model=randomForest(Species~.,data=train)

plot(model)

p=predict(model,test)

p

table(p,test$Species)

accuracy= (17+16+16)/50*100

accuracy
```

OUTPUT:

```

> source("~/active-rstudio-document", echo=TRUE)

> library(randomForest)

> iris_1=iris[sample(150),]

> view(iris_1)

> train=iris_1[1:100,]

> test=iris_1[101:150,]

> model=randomForest(Species~.,data=train)

> plot(model)

> p=predict(model,test)

> p
      135      96      92      145      71      49      5
50      24      82      16
versicolor versicolor versicolor virginica virginica setosa setosa
setosa      setosa versicolor setosa
      73      116      150      67      133      78      76
149      77      83      86
versicolor virginica virginica versicolor virginica virginica versicolor
virginica versicolor versicolor versicolor
      58      132      4      101      9      57      47
43      51      41      127
versicolor virginica setosa virginica setosa versicolor setosa
setosa versicolor setosa virginica
      143      28      25      17      91      21      140
108      106      130      69
virginica setosa setosa setosa versicolor setosa virginica
virginica virginica virginica versicolor
      11      95      114      22      15      81
setosa versicolor virginica setosa setosa versicolor
Levels: setosa versicolor virginica

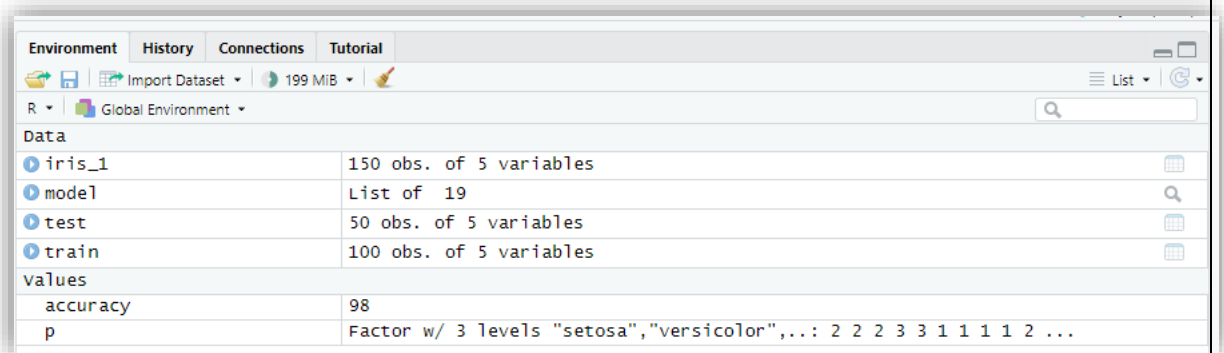
> table(p,test$Species)

p      setosa versicolor virginica
setosa      17          0          0
versicolor   0         16          1
virginica    0          2         14

> accuracy= (17+16+16)/50*100

> accuracy
[1] 98
> |

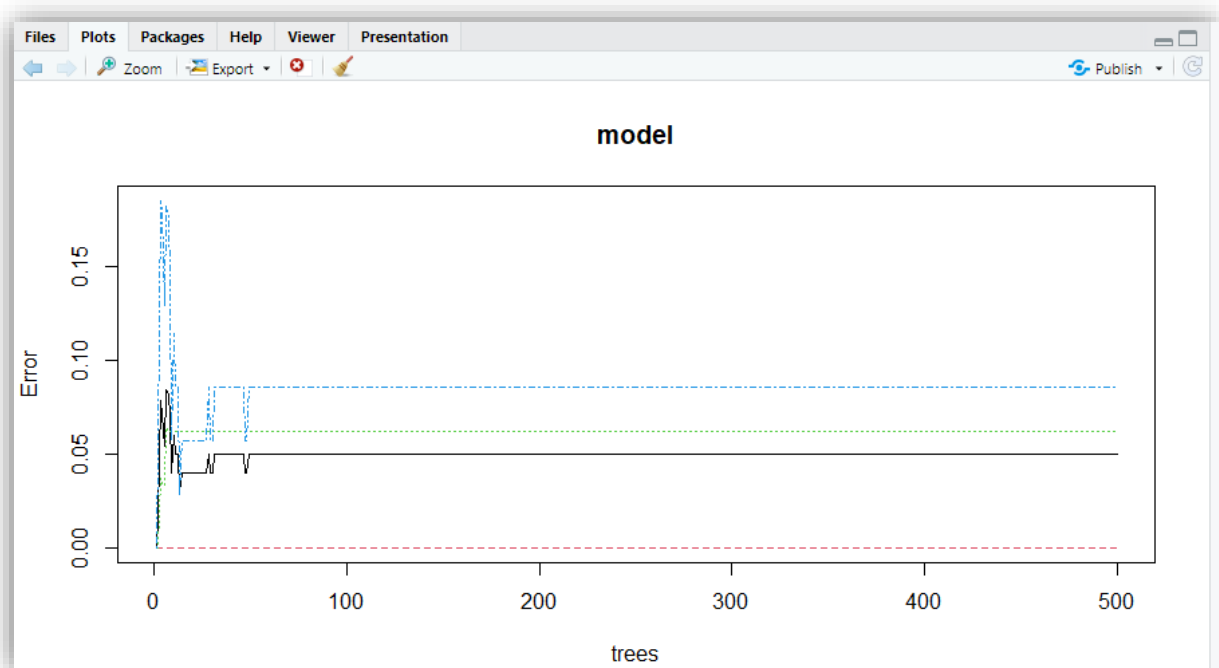
```



The RStudio Environment pane displays the following data and values:

Data	
iris_1	150 obs. of 5 variables
model	List of 19
test	50 obs. of 5 variables
train	100 obs. of 5 variables

Values	
accuracy	98
p	Factor w/ 3 levels "setosa","versicolor",...: 2 2 2 3 3 1 1 1 1 2 ...



RESULT : The Random Forest has been implemented successfully.

PRACTICAL 9

AIM : Implement K -Mean clustering in your own dataset create database using vector

PROCEDURE:

Step 1: Choose the number **K** clusters.

Step 2: Select at random K points, the centroids(Not necessarily from the given data).

Step 3: Assign each data point to closest centroid that forms K clusters.

Step 4: Compute and place the new centroid of each centroid.

Step 5: Reassign each data point to new cluster.

CODE :

```
install.packages("ClusterR")
install.packages("cluster")

# Loading package
library(ClusterR)
library(cluster)

# Removing initial label of
# Species from original dataset
iris_1 <- iris[, -5]

# Fitting K-Means clustering Model
# to training dataset
set.seed(240) # Setting seed
kmeans.re <- kmeans(iris_1, centers = 3, nstart = 20)
kmeans.re

# Cluster identification for
# each observation
kmeans.re$cluster

# Confusion Matrix
cm <- table(iris$Species, kmeans.re$cluster)
cm

# Model Evaluation and visualization
plot(iris_1[c("Sepal.Length", "Sepal.Width")])
plot(iris_1[c("Sepal.Length", "Sepal.Width")],
     col = kmeans.re$cluster)
```



```

> # Confusion Matrix
> cm <- table(iris$Species, kmeans.re$cluster)

> cm
           1  2  3
setosa     50  0  0
versicolor 0 48  2
virginica  0 14 36

> # Model Evaluation and visualization
> plot(iris_1[c("Sepal.Length", "Sepal.Width")])

> plot(iris_1[c("Sepal.Length", "Sepal.Width")],
+       col = kmeans.re$cluster)

> plot(iris_1[c("Sepal.Length", "Sepal.Width")],
+       col = kmeans.re$cluster,
+       main = "K-means with 3 clusters")

> ## Plotting cluster centers
> kmeans.re$centers
Sepal.Length Sepal.Width Petal.Length Petal.Width
1      5.006000      3.428000      1.462000      0.246000
2      5.901613      2.748387      4.393548      1.433871
3      6.850000      3.073684      5.742105      2.071053

> kmeans.re$centers[, c("Sepal.Length", "Sepal.Width")]
Sepal.Length Sepal.Width
1      5.006000      3.428000
2      5.901613      2.748387
3      6.850000      3.073684

```


PRACTICAL 9

AIM : Implement K -Mean clustering in your own dataset create database using vector

PROCEDURE:

Step 1: Choose the number **K** clusters.

Step 2: Select at random K points, the centroids(Not necessarily from the given data).

Step 3: Assign each data point to closest centroid that forms K clusters.

Step 4: Compute and place the new centroid of each centroid.

Step 5: Reassign each data point to new cluster.

CODE :

```
install.packages("ClusterR")
install.packages("cluster")

# Loading package
library(ClusterR)
library(cluster)

# Removing initial label of
# Species from original dataset
iris_1 <- iris[, -5]

# Fitting K-Means clustering Model
# to training dataset
set.seed(240) # Setting seed
kmeans.re <- kmeans(iris_1, centers = 3, nstart = 20)
kmeans.re

# Cluster identification for
# each observation
kmeans.re$cluster

# Confusion Matrix
cm <- table(iris$Species, kmeans.re$cluster)
cm

# Model Evaluation and visualization
plot(iris_1[c("Sepal.Length", "Sepal.Width")])
plot(iris_1[c("Sepal.Length", "Sepal.Width")],
     col = kmeans.re$cluster)
plot(iris_1[c("Sepal.Length", "Sepal.Width")],
     col = kmeans.re$cluster,
```



```

> # Confusion Matrix
> cm <- table(iris$Species, kmeans.re$cluster)

> cm
      1  2  3
setosa 50  0  0
versicolor 0 48  2
virginica 0 14 36

> # Model Evaluation and visualization
> plot(iris_1[c("Sepal.Length", "Sepal.Width")])

> plot(iris_1[c("Sepal.Length", "Sepal.Width")],
+      col = kmeans.re$cluster)

> plot(iris_1[c("Sepal.Length", "Sepal.Width")],
+      col = kmeans.re$cluster,
+      main = "K-means with 3 clusters")

> ## Plotting cluster centers
> kmeans.re$centers
  Sepal.Length Sepal.Width Petal.Length Petal.Width
1    5.006000    3.428000    1.462000    0.246000
2    5.901613    2.748387    4.393548    1.433871
3    6.850000    3.073684    5.742105    2.071053

> kmeans.re$centers[, c("Sepal.Length", "Sepal.Width")]
  Sepal.Length Sepal.Width
1    5.006000    3.428000
2    5.901613    2.748387
3    6.850000    3.073684

```

```

> # cex is font size, pch is symbol
> points(kmeans.re$centers[, c("Sepal.Length", "Sepal.Width")],
+       col = 1:3, pch = 8, cex = 3)

> ## Visualizing clusters
> y_kmeans <- kmeans.re$cluster

> clusplot(iris_1[, c("Sepal.Length", "Sepal.Width")],
+         y_kmeans,
+         lines = 0,
+         shade = TRUE,
+         color = TRUE .... [TRUNCATED]

```

Environment

History

Connections

Tutorial

Import Dataset

164 MiB

R

Global Environment

List

Data

▶ data

3 obs. of 5 variables

▶ iris_1

150 obs. of 4 variables

▶ kmeans.re

List of 9

▶ stu_low_marks

2 obs. of 5 variables

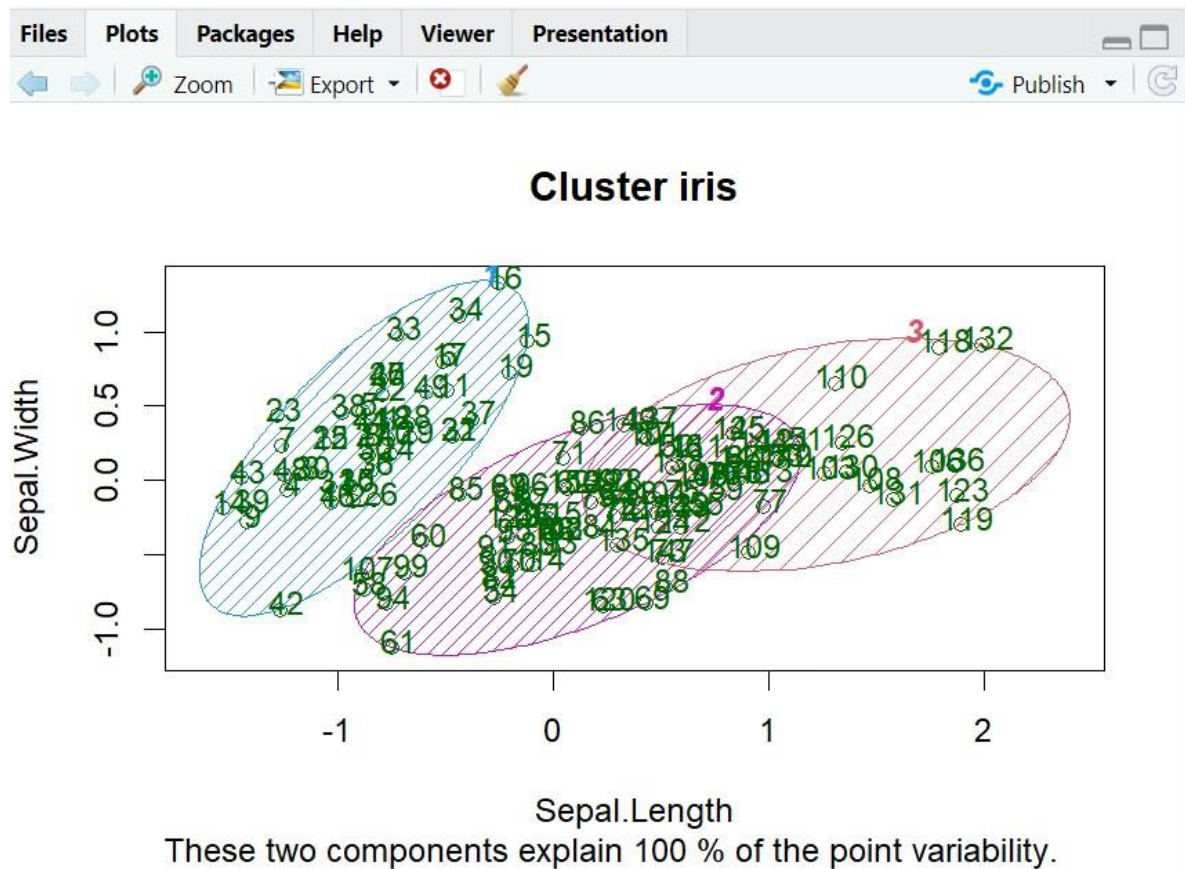
Values

cm

'table' int [1:3, 1:3] 50 0 0 0 48 14 0 2 36

y_kmeans

int [1:150] 1 1 1 1 1 1 1 1 1 1 ...



RESULT: The K-Mean Clustering Algorithm has been implemented successfully.

PRACTICAL-

10

AIM: Implement Linear Regression on iris dataset.

PROCEDURE:

- Step 1: Load the data: Load the dataset that you want to use for predicting the target variable. The dataset should have at least one dependent variable and one independent variable.
- Step 2. Split the data: Split the dataset into training and testing datasets. The training dataset is used to fit the model, and the testing dataset is used to evaluate the model's performance.
- Step 3. Prepare the data: Prepare the data by cleaning and preprocessing it. This includes removing missing values, handling categorical variables, and scaling the data if necessary.
- Step 4. Choose the type of Linear Regression: Choose the type of Linear Regression based on the problem you are trying to solve. There are two types of Linear Regression: Simple Linear Regression and Multiple Linear Regression.
- Step 5. Fit the model: Fit the Linear Regression model on the training dataset using the chosen algorithm. In the case of Simple Linear Regression, the model is a straight line that best fits the data. In the case of Multiple Linear Regression, the model is a hyperplane that best fits the data.
- Step 6. Evaluate the model: Evaluate the performance of the model on the testing dataset. This can be done using metrics such as Mean Squared Error (MSE), Root Mean Squared Error (RMSE), Mean Absolute Error (MAE), R-squared value, or Adjusted R squared value.
- Step 7. Make predictions: Use the trained Linear Regression model to make predictions on new data. This can be done using the predict() function in R.
- Step 8. Fine-tune the model: Fine-tune the Linear Regression model by tweaking the hyperparameters and optimizing the model using techniques such as regularization, feature selection, and cross-validation.
- Step 9. Deploy the model: Deploy the trained Linear Regression model to make predictions on new data in production environments

CODE:

```
require("datasets")
data("iris") str(iris)
head(iris)
Y<- iris[,"Sepal.Width"] # select Target attribute
```

```

X<- iris[,"Sepal.Length"] # select Predictor attribute head(X)
xycorr<- cor(Y,X, method="pearson") # find pearson correlation coefficient
xycorr # a value near 1 implies high correlation and that near 0 shows low correlation plot(Y~X,
col=X)
model1<- lm(Y~X)
model1 # provides regression line coefficients i.e. slope and y-intercept plot(Y~X,
col=X) # scatter plot between X and Y
abline(model1, col="blue", lwd=3)

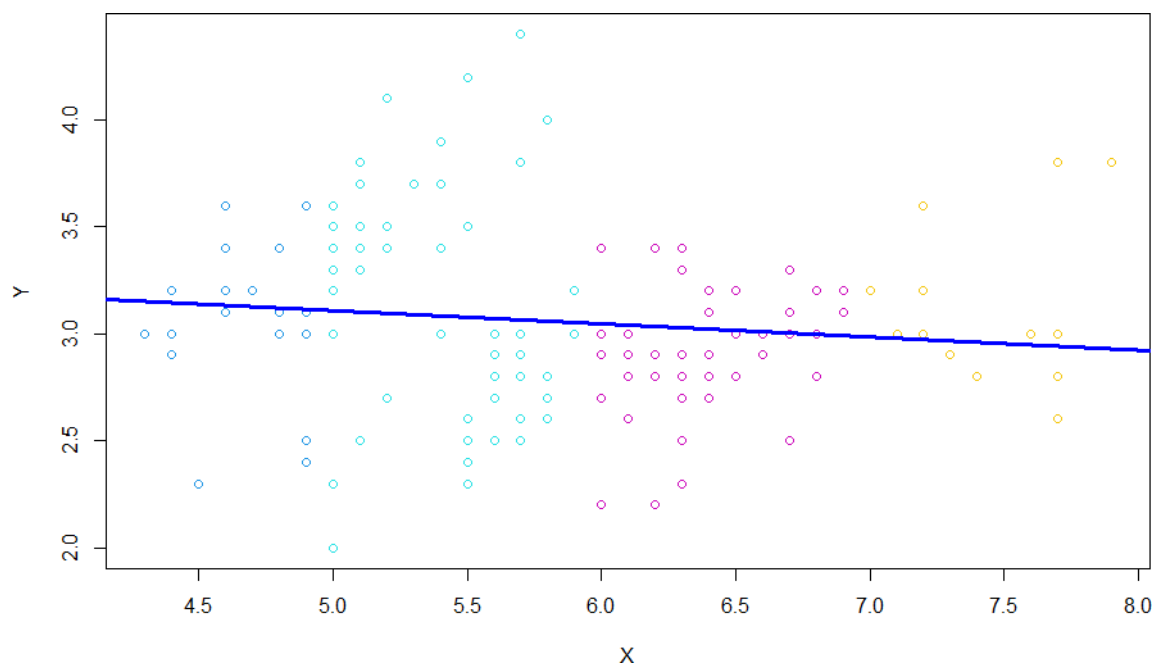
```

OUTPUT:

```

> source("~/active-rstudio-document")
'data.frame': 150 obs. of 5 variables:
 $ Sepal.Length: num 5.1 4.9 4.7 4.6 5 5.4 4.6 5 4.4 4.9 ...
 $ Sepal.Width : num 3.5 3 3.2 3.1 3.6 3.9 3.4 3.4 2.9 3.1 ...
 $ Petal.Length: num 1.4 1.4 1.3 1.5 1.4 1.7 1.4 1.5 1.4 1.5 ...
 $ Petal.Width : num 0.2 0.2 0.2 0.2 0.2 0.4 0.3 0.2 0.2 0.1 ...
 $ Species : Factor w/ 3 levels "setosa","versicolor",...: 1 1 1 1 1 1 1 1 1 1 ...
>

```



RESULT : The linear Regression has been implemented successfully.

PRACTICAL 11

AIM- Implement SVM algorithm using WDBC datasets.

PROCEDURE:

Here are the steps for implementing the SVM algorithm:

1. Load the data: Load the dataset that you want to classify or predict. The dataset should be split into a training set and a test set.
2. Normalize the data: Normalize the data by scaling the features to have zero mean and unit variance.
3. Choose the kernel function: Choose a kernel function that maps the data to a higherdimensional space. The most commonly used kernel functions are the linear kernel, the polynomial kernel, and the radial basis function (RBF) kernel.
4. Train the model: Train the SVM model on the training data using the chosen kernel function.
5. Choose the regularization parameter: Choose the regularization parameter (C) that controls the trade-off between maximizing the margin and minimizing the classification error.
6. Test the model: Test the SVM model on the test data to evaluate its performance.
7. Tune the hyperparameters: Tune the hyperparameters of the SVM model using techniques such as grid search or random search to find the best combination of kernel function, regularization parameter, and other hyperparameters.
8. Deploy the model: Deploy the trained SVM model to make predictions on new data.

CODE:

```
dataset =
read.csv('social.csv')dataset
= dataset[3:5]

dataset$Purchased = factor(dataset$Purchased, levels =
c(0, 1))install.packages('caTools')

library(caTools) set.seed(123)

split = sample.split(dataset$Purchased, SplitRatio = 0.75)
```



```

training_set = subset(dataset, split == TRUE)
test_set = subset(dataset, split == FALSE training_set[-3]

= scale(training_set[-3]) test_set[-3] =

scale(test_set[-3])install.packages('e1071')

library(e1071)
classifier = svm(formula =
Purchased ~ .,data = training_set,
type = 'C- classification' kernel =
'linear')

y_pred = predict(classifier,
newdata =test_set[-3]) cm =
table(test_set[, 3], y_pred) set =
test_set

X1 = seq(min(set[, 1]) - 1, max(set[, 1]) + 1, by = 0.01) X2
= seq(min(set[, 2]) - 1, max(set[, 2]) + 1, by =
0.01)grid_set = expand.grid(X1, X2)

colnames(grid_set) = c('Age',
'EstimatedSalary') y_grid =
predict(classifier, newdata = grid_set)
plot(set[, -3], main = 'SVM (Test set)',
xlab =
'Age', ylab = 'Estimated Salary',xlim =
range(X1), ylim = range(X2))

contour(X1, X2, matrix(as.numeric(y_grid), length(X1), length(X2)), add =
TRUE) points(grid_set, pch = '.', col = ifelse(y_grid == 1, 'coral1',
'aquamarine')) points(set,pch = 21, bg = ifelse(set[, 3] == 1, 'green4', 'red3'))

```

OUTPUT:-

```

> training_set[-3] = scale(training_set[-3])
> test_set[-3] = scale(test_set[-3])
> install.packages('e1071')
WARNING: Rtools is required to build R packages but is not currently installed. Please download and install the appropriate version of Rtools before proceeding:

https://cran.rstudio.com/bin/windows/Rtools/
Installing package into 'C:/Users/akash/AppData/Local/R/win-library/4.2'
(as 'lib' is unspecified)
trying URL 'https://cran.rstudio.com/bin/windows/contrib/4.2/e1071_1.7-13.zip'
Content type 'application/zip' length 652584 bytes (637 KB)
downloaded 637 KB

package 'e1071' successfully unpacked and MD5 sums checked

The downloaded binary packages are in
C:\Users\akash\AppData\Local\Temp\RtmpMLO1q5\downloaded_packages
> library(e1071)
>
> classifier = svm(formula = Purchased ~ .,
+                  data = training_set,
+                  type = 'C-classification',
+                  kernel = 'linear')
> y_pred = predict(classifier, newdata = test_set[-3])
> cm = table(test_set[, 3], y_pred)
> set = test_set
> X1 = seq(min(set[, 1]) - 1, max(set[, 1]) + 1, by = 0.01)
> X2 = seq(min(set[, 2]) - 1, max(set[, 2]) + 1, by = 0.01)
>
> grid_set = expand.grid(X1, X2)

> colnames(grid_set) = c('Age', 'EstimatedSalary')
> y_grid = predict(classifier, newdata = grid_set)
>
> plot(set[, -3], main = 'SVM (Test set)',
+      xlab = 'Age', ylab = 'Estimated Salary',
+      xlim = range(X1), ylim = range(X2))
>
> contour(X1, X2, matrix(as.numeric(y_grid), length(X1), length(X2)), add = TRUE)
>
> points(grid_set, pch = '.', col = ifelse(y_grid == 1, 'coral1', 'aquamarine'))
>
> points(set, pch = 21, bg = ifelse(set[, 3] == 1, 'green4', 'red3'))
>

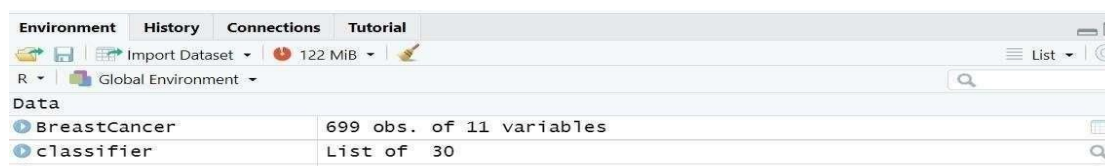
> dataset = read.csv('social.csv')
> dataset = dataset[3:5]
> dataset$Purchased = factor(dataset$Purchased, levels = c(0, 1))
> install.packages('caTools')
WARNING: Rtools is required to build R packages but is not currently installed. Please download and install the appropriate version of Rtools before proceeding:

https://cran.rstudio.com/bin/windows/Rtools/
Installing package into 'C:/Users/akash/AppData/Local/R/win-library/4.2'
(as 'lib' is unspecified)
trying URL 'https://cran.rstudio.com/bin/windows/contrib/4.2/caTools_1.18.2.zip'
Content type 'application/zip' length 246159 bytes (240 KB)
downloaded 240 KB

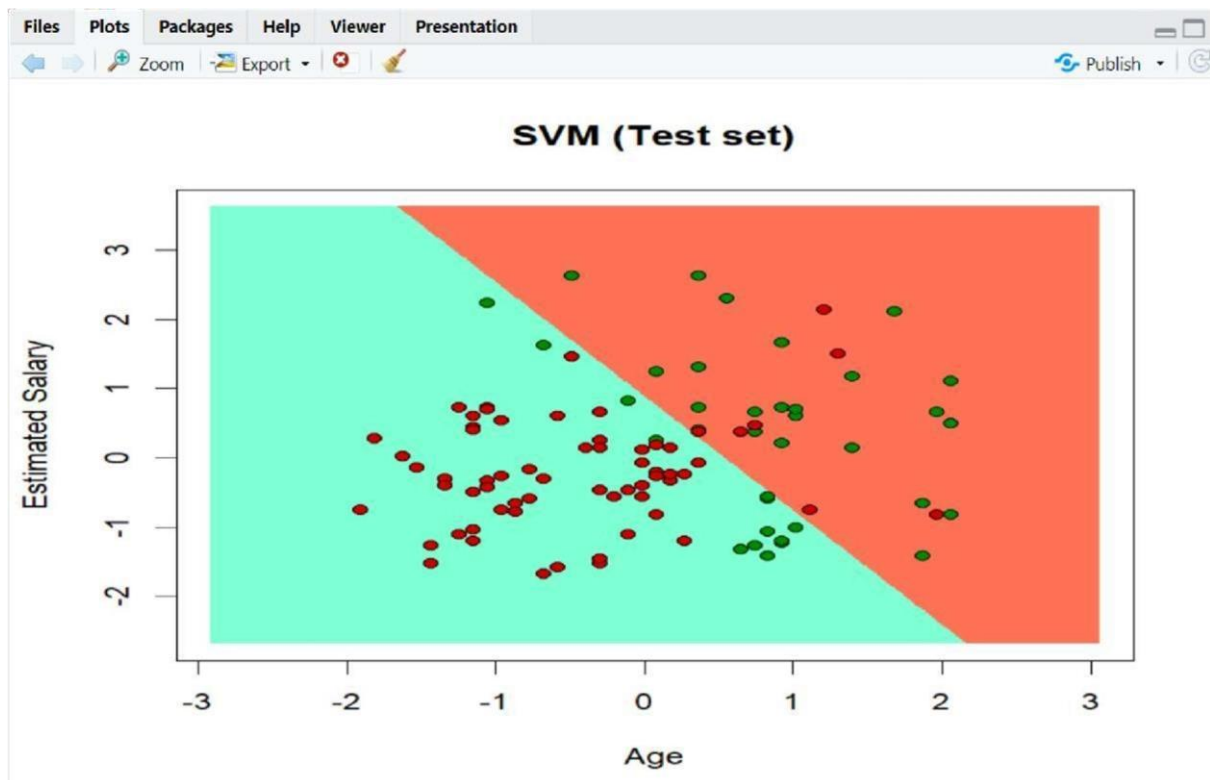
package 'caTools' successfully unpacked and MD5 sums checked

The downloaded binary packages are in
C:\Users\akash\AppData\Local\Temp\RtmpMLO1q5\downloaded_packages
> library(caTools)
>
> set.seed(123)
> split = sample.split(dataset$Purchased, splitRatio = 0.75)
>
> training_set = subset(dataset, split == TRUE)
> test_set = subset(dataset, split == FALSE)

```



Environment	History	Connections	Tutorial
<div> <div>Import Dataset</div> <div>122 MiB</div> </div> <div> <div>R</div> <div>Global Environment</div> </div>			
Data			
BreastCancer	699 obs. of 11 variables		
classifier	List of 30		



Result:- The SVM Algorithm have been implemented successfully.

PRACTICAL 12

Aim: Implement Logistic regression on IRIS datasets.

PROCEDURE:

1. Import the dataset: Import the dataset which contains the input variables (or the independent variables) and the output variable (or the dependent variable).
2. Data Pre-processing: Check for missing values and handle them, handle outliers if any, and remove or transform variables that do not contribute to the model.
3. Split the dataset: Split the dataset into a training set and a testing set. The training set is used to train the model, while the testing set is used to evaluate the performance of the model.
4. Feature Scaling: Scale the input variables (or the independent variables) to have the same range to ensure that no variable dominates the others in the modeling process.
5. Model Training: Train the logistic regression model using the training set. The logistic regression model calculates the probability of the output variable taking a particular value given the input variables (or the independent variables).
6. Model Evaluation: Evaluate the performance of the model using the testing set. There are several metrics to evaluate the performance of the model, such as accuracy, precision, recall, F1 score, and area under the ROC curve.
7. Model Improvement: If the model does not perform well, try to improve the model by changing the hyperparameters, selecting the relevant features, or using other modeling techniques.
8. Model Deployment: Once the model is finalized and performs well on the testing set, deploy it to the production environment for making predictions on new data.

CODE:

```
install.packages("dplyr")
install.packages("caTools")
install.packages("ROCR")

# loading packages

library(dplyr)
library(caTools
)
```

```

library(ROCR) dataSet =
mtcars View(dataSet)

split <- sample.split(mtcars, SplitRatio = 0.8)
table(split)

train_reg <- subset(mtcars, split ==
"TRUE")test_reg <- subset(mtcars, split ==
"FALSE") # Training model

logistic_model <- glm(vs ~ wt +
disp,data = train_reg,

family = "binomial")

logistic_model
summary(logistic_model)
predict_reg
<
-
predict(logistic_model,

test_reg, type = "response") table(predict_reg)
predict_reg <- ifelse(predict_reg >0.5, 1, 0)
table(test_reg$vs, predict_reg)
missing_classerr <- mean(predict_reg != test_reg$vs)
print(paste('Accuracy =', 1 - missing_classerr))
ROCPred <- prediction(predict_reg, test_reg$vs)
ROCPer <- performance(ROCPred, measure = "tpr", x.measure = "fpr")

auc <- performance(ROCPred, measure = "auc")
auc <- auc@y.values[[1]]

auc plot(ROCPer) plot(ROCPer, colorize = TRUE,
print.cutoffs.at = seq(0.1, by = 0.1),

main = "ROC CURVE")

abline(a = 0, b = 1) auc <- round(auc, 4)

legend(.6, .4, auc, title = "AUC", cex = 1)

```

```

> library(caTools)
> library(ROCR)
> # Summary of dataset in package
> dataSet = mtcars
> View(dataSet)
> # Splitting dataset
> split <- sample.split(mtcars, SplitRatio = 0.8)
> table(split)
split
FALSE TRUE
   3    8
> train_reg <- subset(mtcars, split == "TRUE")
> test_reg <- subset(mtcars, split == "FALSE")
> # Training model
> logistic_model <- glm(vs ~ wt + disp,
+                       data = train_reg,
+                       family = "binomial")
> logistic_model

Call:  glm(formula = vs ~ wt + disp, family = "binomial", data = train_reg)

Coefficients:
(Intercept)          wt          disp
    0.84002      2.67137     -0.04907

Degrees of Freedom: 22 Total (i.e. Null);  20 Residual
Null Deviance:      31.49
Residual Deviance: 13.43      AIC: 19.43
> # Summary

> summary(logistic_model)

Call:
glm(formula = vs ~ wt + disp, family = "binomial", data = train_reg)

Deviance Residuals:
    Min       1Q   Median       3Q      Max
-1.67702  -0.08231  -0.02704   0.55160   1.60126

Coefficients:
              Estimate Std. Error z value Pr(>|z|)
(Intercept)  0.84002    2.75234   0.305  0.7602
wt           2.67137    1.98927   1.343  0.1793
disp        -0.04907    0.02637  -1.861  0.0628 .
---
Signif. codes:  0 '***' 0.001 '**' 0.01 '*' 0.05 '.' 0.1 ' ' 1

(Dispersion parameter for binomial family taken to be 1)

    Null deviance: 31.492  on 22  degrees of freedom
Residual deviance: 13.427  on 20  degrees of freedom
AIC: 19.427

Number of Fisher Scoring iterations: 7

> # Predict test data based on model

> predict_reg <- predict(logistic_model,
+                       test_reg, type = "response")
> table(predict_reg)
predict_reg
0.000249778611633657 0.00739119585512458 0.0122220979368047 0.0380270689031347
      1              1              1              1
0.139427936790023   0.497091675739643   0.894056832644504   0.904939540010952
      1              1              1              1
0.912637391844415
      1
> # Changing probabilities
> predict_reg <- ifelse(predict_reg > 0.5, 1, 0)
> table(test_reg$vs, predict_reg)
      predict_reg
      0 1
0 5 0
1 1 3

```



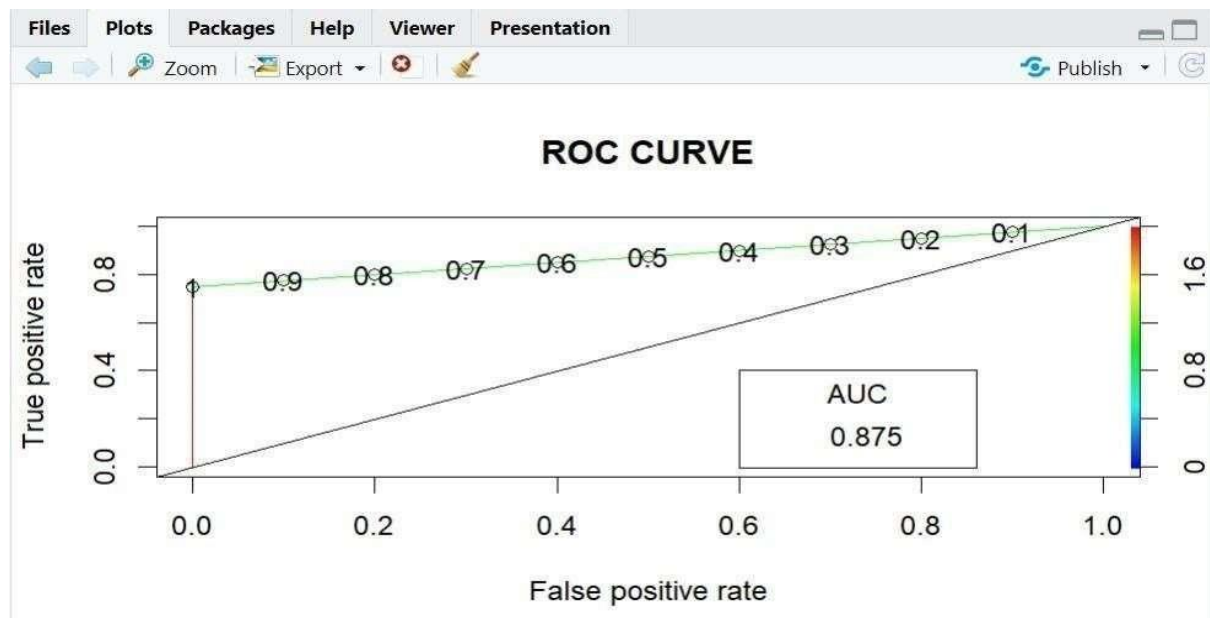
```

> missing_classerr <- mean(predict_reg != test_reg$vs)
> print(paste('Accuracy =', 1 - missing_classerr))
[1] "Accuracy = 0.888888888888889"
> # ROC-AUC Curve
> ROCPred <- prediction(predict_reg, test_reg$vs)
> ROCPer <- performance(ROCPred, measure = "tpr", x.measure
+                       = "fpr")
> auc <- performance(ROCPer, measure = "auc")
> auc <- auc@y.values[[1]]
> auc
[1] 0.875
> # Plotting curve
> plot(ROCPer)

> plot(ROCPer, colorize = TRUE,
+      print.cutoffs.at = seq(0.1, by = 0.1),
+      main = "ROC CURVE")
> abline(a = 0, b = 1)
> auc <- round(auc, 4)
> legend(.6, .4, auc, title = "AUC", cex = 1)

```

Environment	History	Connections	Tutorial
<div> <div>Import Dataset</div> <div>330 MiB</div> <div>List</div> </div>			
R Global Environment			
Data			
dataSet	32 obs. of 11 variables		
logistic_mod...	List of 30		
ROCPer	Formal class performance		
ROCPred	Formal class prediction		
test_reg	9 obs. of 11 variables		



Result: The logistic regression have been implemented successfully

PRACTICAL 13

AIM: Implement Apriori algorithm.

PROCEDURE:

1. Read each item in the transaction.
2. Calculate the support of every item.
3. If support is less than minimum support, discard the item. Else, insert it into frequent itemset.
4. Calculate confidence for each non- empty subset.
5. If confidence is less than minimum confidence, discard the subset. Else, it into strong rules.

CODE:

```
install.packages("arules")
install.packages("arulesViz")
install.packages("RColorBrewer")
library(arules) library(arulesViz)
library(RColorBrewer)
data("Groceries") rules <-
apriori(Groceries, parameter =
list(supp = 0.01, conf = 0.2))
inspect(rules[1:10])
arules::itemFrequencyPlot

(Groceries, topN = 20,

col = brewer.pal(8, 'Pastel2'),

main = 'Relative Item Frequency Plot',
type = "relative" ylab = "Item Frequency (Relative)")
```


Output:

```

> library(arules)
> library(arulesViz)
> library(RColorBrewer)
>
> # import dataset
> data("Groceries")
>
> # using apriori() function

> rules <- apriori(Groceries,
+                 parameter = list(supp = 0.01, conf = 0.2))
Apriori

Parameter specification:
 confidence minval  smax arem  aval originalSupport maxtime support minlen
      0.2      0.1   1 none FALSE               TRUE     5   0.01      1
maxlen target  ext
     10  rules TRUE

Algorithmic control:
 filter tree heap memopt load sort verbose
  0.1 TRUE TRUE  FALSE TRUE    2    TRUE

Absolute minimum support count: 98

set item appearances ...[0 item(s)] done [0.00s].
set transactions ...[169 item(s), 9835 transaction(s)] done [0.00s].
sorting and recoding items ... [88 item(s)] done [0.00s].
creating transaction tree ... done [0.00s].
checking subsets of size 1 2 3 4 done [0.00s].
writing ... [232 rule(s)] done [0.00s].
creating S4 object ... done [0.00s].

> # using inspect() function
> inspect(rules[1:10])

```

	lhs	rhs	support	confidence	coverage
[1]	{}	=> {whole milk}	0.25551601	0.2555160	1.00000000
[2]	{hard cheese}	=> {whole milk}	0.01006609	0.4107884	0.02450432
[3]	{butter milk}	=> {other vegetables}	0.01037112	0.3709091	0.02796136
[4]	{butter milk}	=> {whole milk}	0.01159126	0.4145455	0.02796136
[5]	{ham}	=> {whole milk}	0.01148958	0.4414062	0.02602949
[6]	{sliced cheese}	=> {whole milk}	0.01077783	0.4398340	0.02450432
[7]	{oil}	=> {whole milk}	0.01128622	0.4021739	0.02806304
[8]	{onions}	=> {other vegetables}	0.01423488	0.4590164	0.03101169
[9]	{onions}	=> {whole milk}	0.01209964	0.3901639	0.03101169
[10]	{berries}	=> {yogurt}	0.01057448	0.3180428	0.03324860

```


```

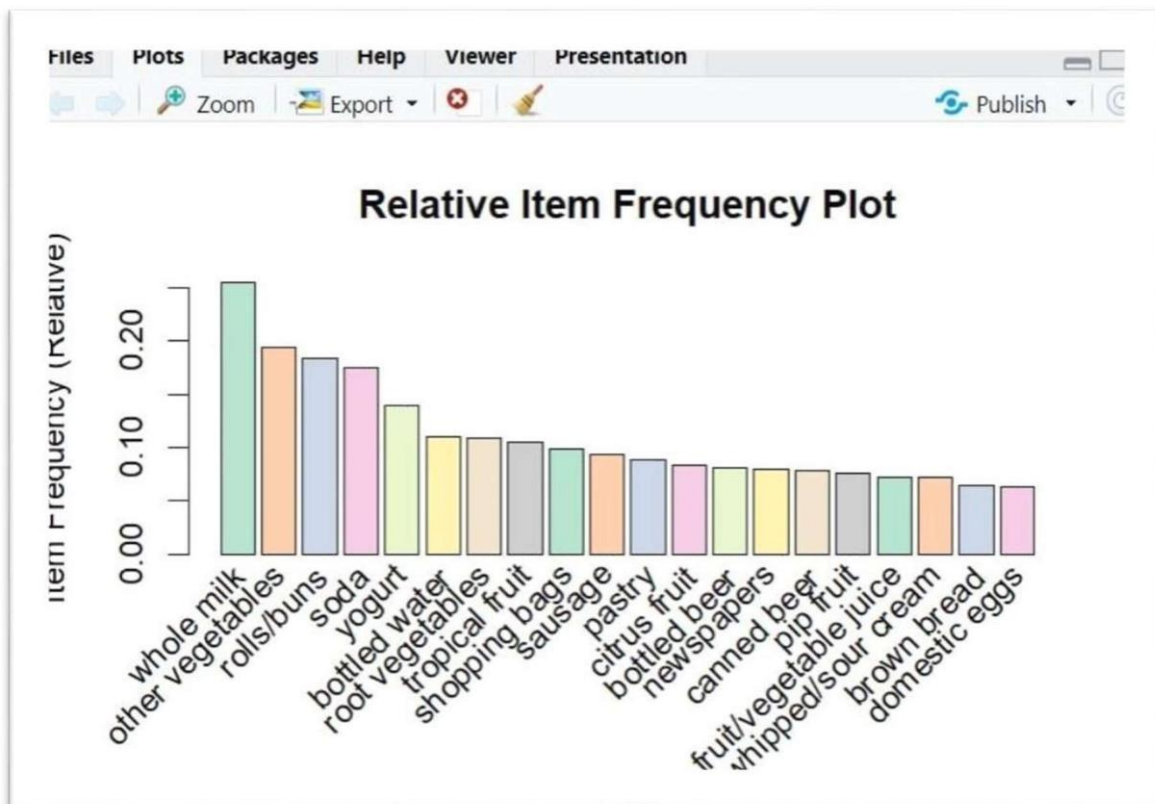
	lift	count
[1]	1.000000	2513
[2]	1.607682	99
[3]	1.916916	102
[4]	1.622385	114
[5]	1.727509	113
[6]	1.721356	106
[7]	1.573968	111
[8]	2.372268	140
[9]	1.526965	119
[10]	2.279848	104

```

>
> # using itemFrequencyPlot() function
> arules::itemFrequencyPlot(Groceries, topN = 20,
+                           col = brewer.pal(8, 'Pastel2'),
+                           main = 'Relative Item Frequency Plot',
+                           type = "relative",
+                           ylab = "Item Frequency (Relative)")

```

Environment	History	Connections	Tutorial
<div> <div> <div>Import Dataset</div> <div>19 MiB</div> </div> <div>List</div> </div>			
<div> <div>R</div> <div>Global Environment</div> </div>			
Data			
Groceries	Formal class	transactions	
rules	Formal class	rules	



Result: The Apriori algorithm have been implemented successfully

PRACTICAL-14

AIM: Input two 3x3 Matrices and display both and their addition matrix.

PROCEDURE:

```
# Input first 3x3 matrix
mat1 <- matrix(numeric(), nrow = 3, ncol = 3)
cat("Enter elements of the first 3x3 matrix:\n")
for (i in 1:3) {
  for (j in 1:3) {
    mat1[i, j] <- as.numeric(readline(prompt = paste("Enter element [", i, ", ", j, "]: ")))
  }
}

# Input second 3x3 matrix
mat2 <- matrix(numeric(), nrow = 3, ncol = 3)
cat("Enter elements of the second 3x3 matrix:\n")
for (i in 1:3) {
  for (j in 1:3) {
    mat2[i, j] <- as.numeric(readline(prompt = paste("Enter element [", i, ", ", j, "]: ")))
  }
}

# Display both matrices
cat("\nFirst Matrix:\n")
print(mat1)
cat("\nSecond Matrix:\n")
print(mat2)

# Addition of matrices
add_mat <- mat1 + mat2
# Display addition matrix
cat("\nAddition Matrix:\n")
print(add_mat)
```

INPUT/OUTPUT:

```

Console Terminal × Background Jobs ×
R 4.3.3 · ~/
Enter element [ 1 , 1 ]: 9
Enter element [ 2 , 3 ]: 4
Enter element [ 3 , 1 ]: 3
Enter element [ 3 , 2 ]: 2
Enter element [ 3 , 3 ]: 1

> # Display both matrices
> cat("\nFirst Matrix:\n")

First Matrix:

> print(mat1)
      [,1] [,2] [,3]
[1,]    1    2    3
[2,]    4    5    6
[3,]    7    8    9

> cat("\nSecond Matrix:\n")

Second Matrix:

> print(mat2)
      [,1] [,2] [,3]
[1,]    9    8    7
[2,]    6    5    4
[3,]    3    2    1

> # Addition of matrices
> add_mat <- mat1 + mat2

> # Display addition matrix
> cat("\nAddition Matrix:\n")

Addition Matrix:

> print(add_mat)
      [,1] [,2] [,3]
[1,]   10   10   10
[2,]   10   10   10
[3,]   10   10   10
> |

```

RESULT: The Code Has Been Executed Successfully.

PRACTICAL -15**AIM:** Import IRIS dataset and display First, 3rd and last columns only.**PROCEDURE:**

Import IRIS dataset

data(iris)

Display all data

print(iris)

Display first, third and last columns

iris_subset <- iris[, c(1, 3, ncol(iris))]

print(iris_subset)

INPUT/OUTPUT:

Console Terminal x Background Jobs x

R 4.3.3 · ~/

```
> print(iris)
```

	Sepal.Length	Sepal.Width	Petal.Length	Petal.Width	Species
1	5.1	3.5	1.4	0.2	setosa
2	4.9	3.0	1.4	0.2	setosa
3	4.7	3.2	1.3	0.2	setosa
4	4.6	3.1	1.5	0.2	setosa
5	5.0	3.6	1.4	0.2	setosa
6	5.4	3.9	1.7	0.4	setosa
7	4.6	3.4	1.4	0.3	setosa
8	5.0	3.4	1.5	0.2	setosa
9	4.4	2.9	1.4	0.2	setosa
10	4.9	3.1	1.5	0.1	setosa
11	5.4	3.7	1.5	0.2	setosa
12	4.8	3.4	1.6	0.2	setosa
13	4.8	3.0	1.4	0.1	setosa
14	4.3	3.0	1.1	0.1	setosa
15	5.8	4.0	1.2	0.2	setosa
16	5.7	4.4	1.5	0.4	setosa
17	5.4	3.9	1.3	0.4	setosa
18	5.1	3.5	1.4	0.3	setosa
19	5.7	3.8	1.7	0.3	setosa
20	5.1	3.8	1.5	0.3	setosa
21	5.4	3.4	1.7	0.2	setosa
22	5.1	3.7	1.5	0.4	setosa
23	4.6	3.6	1.0	0.2	setosa
24	5.1	3.3	1.7	0.5	setosa
25	4.8	3.4	1.9	0.2	setosa
26	5.0	3.0	1.6	0.2	setosa
27	5.0	3.4	1.6	0.4	setosa
28	5.2	3.5	1.5	0.2	setosa
29	5.2	3.4	1.4	0.2	setosa
30	4.7	3.2	1.6	0.2	setosa
31	4.8	3.1	1.6	0.2	setosa
32	5.4	3.4	1.5	0.4	setosa
33	5.2	4.1	1.5	0.1	setosa
34	5.5	4.2	1.4	0.2	setosa

```

Console Terminal × Background Jobs ×
R 4.3.3 · ~/
> data(iris)
> iris_subset <- iris[, c(1, 3, ncol(iris))]
> print(iris_subset)
  Sepal.Length Petal.Length Species
1          5.1          1.4   setosa
2          4.9          1.4   setosa
3          4.7          1.3   setosa
4          4.6          1.5   setosa
5          5.0          1.4   setosa
6          5.4          1.7   setosa
7          4.6          1.4   setosa
8          5.0          1.5   setosa
9          4.4          1.4   setosa
10         4.9          1.5   setosa
11         5.4          1.5   setosa
12         4.8          1.6   setosa
13         4.8          1.4   setosa
14         4.3          1.1   setosa
15         5.8          1.2   setosa
16         5.7          1.5   setosa
17         5.4          1.3   setosa
18         5.1          1.4   setosa
19         5.7          1.7   setosa
20         5.1          1.5   setosa
21         5.4          1.7   setosa
22         5.1          1.5   setosa
23         4.6          1.0   setosa
24         5.1          1.7   setosa
25         4.8          1.9   setosa
26         5.0          1.6   setosa
27         5.0          1.6   setosa
28         5.2          1.5   setosa
29         5.2          1.4   setosa
30         4.7          1.6   setosa
31         4.8          1.6   setosa
32         5.4          1.5   setosa
33         5.2          1.5   setosa
34         5.5          1.4   setosa

```

RESULT: The Code Has Been Executed Successfully.