R Programming LAB

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Data Management

1. Working with vectors and Matrices

Aim:

To write R program working with vectors and Matrices

Program:

```
# creating a vector and a matrix
vector_data <- c(101,102,103,104,105)
matrix_data <- matrix(1:9, nrow = 3)</pre>
```

accessing elements in a vector and matrix vector_element <- vector_data[3] matrix_element <- matrix_data[2, 2]</pre>

Performing operations on vectors and matrices
vector_sum <- sum(vector_data)
matrix_transpose <- t(matrix_data)</pre>

Output:

		_					
Data							
matrix_data	int [1:3, 1:3] 1 2 3 4 5 6 7 8 9						
matrix_transpose	int [1:3, 1:3] 1 4 7 2 5 8 3 6 9						
Values							
matrix_element	5L						
vector_data	num [1:5] 101 102 103 104 105						
vector_element	103						
vector_sum	515						

Result:

2. Sorting, Merging and Aggregating Data sets

Aim:

To write R program Sorting, Merging and Aggregating Data sets

Program:

```
vector_data <- c(106,109,133,89104,23105)
sorted_vector <- sort(vector_data)

# merging two vectors
vector1 <- c(1, 2, 3)
vector2 <- c(4, 5, 6)
merged_vector <- c(vector1, vector2)

# aggregating data with mean and sum
data <- c(12, 15, 18, 24, 9)
mean_data <- mean(data)
sum_data <- sum(data)</pre>
```

Output:

values	
data	num [1:5] 12 15 18 24 9
mean_data	15.6
merged_vector	num [1:6] 1 2 3 4 5 6
sorted_vector	num [1:5] 106 109 133 23105 89104
sum_data	78
vector_data	num [1:5] 106 109 133 89104 23105
vector1	num [1:3] 1 2 3
vector2	num [1:3] 4 5 6

Result:

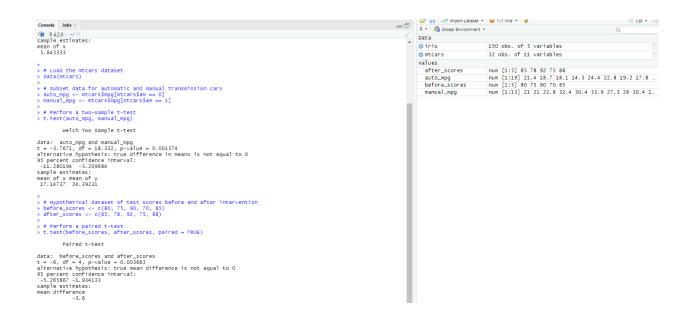
Testing Statistical Hypothesis using R

3. Test for Single, difference of mean and paired mean

Aim:

To write R program Test for Single, difference of mean and paired mean

```
# Load the iris dataset
data(iris)
# Perform a one-sample t-test
t.test(iris$Sepal.Length, mu = 5.0)
# Load the mtcars dataset
data(mtcars)
# Subset data for automatic and manual transmission cars
auto_mpg <- mtcars$mpg[mtcars$am == 0]</pre>
manual_mpg <- mtcars$mpg[mtcars$am == 1]</pre>
# Perform a two-sample t-test
t.test(auto_mpg, manual_mpg)
# Hypothetical dataset of test scores before and after intervention
before_scores <- c(80, 75, 90, 70, 85)
after_scores <- c(85, 78, 92, 75, 88)
# Perform a paired t-test
t.test(before_scores, after_scores, paired = TRUE)
```



Result:

4. Test for equality of variance

Aim:

To write R program Test for equality of variance.

Program:

Generate two sample datasets data1 <- c(23, 26, 29, 32, 35) data2 <- c(20, 25, 30, 35, 40)

Test for equality of variances
result <- var.test(data1, data2)
result</pre>

Output:

Result:

5. Applications: Chi-Square test for Goodness of fit and independence of Attributes

Aim:

To write R program Chi-Square test for Goodness of fit and independence of Attributes.

```
# Creating observed and expected frequency tables for Goodness of Fit observed <- c(35, 45, 60) expected <- c(0.3, 0.4, 0.3) # Expected frequencies should sum to 1 # Chi-Square test for goodness of fit chi_square_goodness_of_fit <- chisq.test(observed, p = expected) # Print the results for Goodness of Fit cat("Chi-Square Test for Goodness of Fit:\n") print(chi_square_goodness_of_fit) # Creating a contingency table for Independence of Attributes table_data <- matrix(c(10, 20, 15, 25, 30, 40, 35, 45), ncol = 2) # Chi-Square test for independence of attributes chi_square_independence <- chisq.test(table_data) # Print the results for Independence of Attributes cat("Chi-Square Test for Independence of Attributes:\n") chi_square_independence
```

Result:

6. Applications: One way ANOVA and two way ANOVA

Aim:

To write R program one way ANOVA and two way ANOVA.

Program:

```
# One-way ANOVA
data1 <- c(10, 15, 20, 25, 30)
data2 <- c(5, 10, 15, 20, 25)
data3 <- c(12, 17, 22, 27, 32)

result_one_way_anova <- aov(c(data1, data2, data3) ~ rep(c("A", "B", "C"), each = 5))
summary(result_one_way_anova)

# Two-way ANOVA
data <- data.frame(
    Treatment = rep(c("A", "B", "C"), each = 15),
    Gender = rep(c("Male", "Female"), each = 45),
    Value = rnorm(90)
)

result_two_way_anova <- aov(Value ~ Treatment * Gender, data = data)
summary(result_two_way_anova)
```

Output:

Result:

7. Applications: Latin Square Design

Aim:

To write R program Latin Square Design

Program:

```
# Create a Latin Square design
latin_square <- matrix(c(3, 2, 1, 2, 1, 3, 1, 3, 2), nrow = 3)
latin_square_design <- as.table(latin_square)</pre>
```

Perform Fisher's Exact Test on the Latin Square design fisher_exact_test <- fisher.test(latin_square_design)

Print the results
cat("Fisher's Exact Test for Latin Square Design:\n")
print(fisher_exact_test)

Output:

Result:

Numerical Solution of Equations using R

8. Newton-Raphson method

Aim:

To write R program Newton-Raphson method

```
# Define a function and its derivative
f < -function(x) x^3 - 2*x - 5
f_prime <- function(x) 3*x^2 - 2
# Implement the Newton-Raphson method
x0 <- 1 # Initial guess
tolerance <- 1e-6
max_iterations <- 100
x <- x0
for (i in 1:max_iterations) {
 x \leftarrow x - f(x) / f_prime(x)
 if (abs(f(x)) < tolerance) {
  break
 }
}
result_newton_raphson <- x
# Print the result
cat("Approximated Root:", result_newton_raphson, "\n")
```



Result:

9. Solving system of Linear Equations (Gauss elimination, Gauss Jacobi and Gauss-Seidel)

Aim:

To write R program solving system of Linear Equations (Gauss elimination, Gauss Jacobi and Gauss-Seidel)

```
# Define the coefficient matrix and right-hand side vector
A \leftarrow matrix(c(2, 1, 1, 1, 3, 2, 2, 4, 3), nrow = 3)
b <- c(7, 8, 18)
# Solve using Gauss elimination
x_gauss <- solve(A, b)</pre>
# Solve using Gauss-Jacobi method
x_jacobi <- solve(A, b, method = "Jacobi")</pre>
# Solve using Gauss-Seidel method
x_seidel <- solve(A, b, method = "Seidel")</pre>
# Print the results
cat("Solution using Gauss Elimination:\n")
print(x_gauss)
cat("Solution using Gauss-Jacobi Method:\n")
print(x_jacobi)
cat("Solution using Gauss-Seidel Method:\n")
print(x_seidel)
```

Result:

10. Power method to approximate dominant Eigen value and Eigen vector

Aim:

To write R program Power method to approximate dominant Eigen value and Eigen vector

```
# Define a matrix
A \leftarrow matrix(c(6, 2, 1, 1, 3, 1, 2, 4, 3), nrow = 3)
# Power method to approximate the dominant eigenvalue and eigenvector
power_method <- function(A, iter = 1000) {</pre>
 n <- nrow(A)
 x < -rep(1, n)
 for (i in 1:iter) {
  y <- A %*% x
  x \leftarrow y / max(y)
 lambda_max <- max(y)</pre>
 return(list(lambda_max = lambda_max, eigenvector = x))
}
result_power_method <- power_method(A)</pre>
# Print the results
cat("Approximated Dominant Eigenvalue:", result_power_method$lambda_max, "\n")
cat("Approximated Dominant Eigenvector:\n")
print(result_power_method$eigenvector)
```

```
Console Jobs x

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**Power method to approximate the dominant eigenvalue and eigenvector power_method x- function(A, iter = 1000) {

**n < - morn(A)

**Torigonal transport | mistory Connections | mistory Connections | mistory Connections | mistory Connections |

**Power method to approximate the dominant eigenvalue and eigenvector |

**power method to approximate the dominant eigenvalue and eigenvector |

**power method x- function(A, iter = 1000) {

**n < - morn(A)

**n < - morn(A)

**Torigonal transport |

**power_method (its of 2 |

**punctions |

**power_method |

**
```

Result:

Numerical Interpolations Using R

11. Lagrange Interpolation

Aim:

To write R program Lagrange Interpolation

Program:

```
# Define known data points
x_values <- c(1, 2, 4, 5)
y_values <- c(3, 5, 9, 11)

# Define the point at which to interpolate
x_interpolate <- 3

# Perform Lagrange interpolation
lagrange_interpolation <- approxfun(x_values, y_values)
y_interpolated <- lagrange_interpolation(x_interpolate)

# Print the result
cat("Interpolated Value at x = ", x_interpolate, ":", y_interpolated, "\n")</pre>
```

Output:

Result:

12. Newton's forward and Backward Interpolation

Aim:

To write R program Newton's forward and Backward Interpolation

```
# Define known data points
x \text{ values} <- c(1, 2, 4, 5)
y_values <- c(3, 5, 9, 11)
# Define the point at which to interpolate
x_interpolate <- 3
# Perform Newton's forward interpolation
newton_forward_interpolation <- approxfun(x_values, y_values, method = "linear")</pre>
y_interpolated_forward <- newton_forward_interpolation(x_interpolate)</pre>
# Perform Newton's backward interpolation
newton_backward_interpolation <- approxfun(x_values, y_values, method = "linear", f = 1)
y interpolated backward <- newton backward interpolation(x interpolate)
# Print the results
cat("Interpolated Value using Newton's Forward Interpolation at x =", x_interpolate, ":",
y_interpolated_forward, "\n")
cat("Interpolated Value using Newton's Backward Interpolation at x =", x_interpolate, ":",
y_interpolated_backward, "\n")
```

```
Source
                                                                                                                                                                                                                      Environment History Connections Tutorial
                                                                                                                                                                                                                                     R • Global Environment • 172 MiB • 4
 R R4.20 · ~/ ∞
> # Define known data points
> x_values <- c(1, 2, 4, 5)
> y_values <- c(3, 5, 9, 11)</pre>
                                                                                                                                                                                                                                    x_interpolate 3
x_values 7
y_interpolated 7
y_interpolated_bac_7
y_interpolated_for_7
y_values nu
Functions
                                                                                                                                                                                                                                                                             num [1:4] 1 2 4 5
   # Define the point at which to interpolate x_interpolate <- 3 \,
  >
> # Perform Lagrange interpolation
> lagrange_interpolation <- approxfun(x_values, y_values)
> y_interpolated <- lagrange_interpolation(x_interpolate)
                                                                                                                                                                                                                                                                              num [1:4] 3 5 9 11
                                                                                                                                                                                                                                        lagrange_interpola_ function (v)
newton_backward_in_ function (v)
newton_forward_int... function (v)
>" Print the result 
> cat("Interpolated Value at x =", x_interpolate, ":", y_interpolated, "\n")
Interpolated value at x = 3 : 7
> # Define known data points
> x_values <- c(1, 2, 4, 5)
> y_values <- c(3, 5, 9, 11)
   # Define the point at which to interpolate x_{interpolate} < -3
   # Perform Newton's forward interpolation
newton_forward_interpolation <- approxfun(x_values, y_values, method = "linear")
y_interpolated_forward <- newton_forward_interpolation(x_interpolate)</pre>
 > 
 # Perform Newton's backward interpolation
> newton_backward_interpolation <- approxfun(x_values, y_values, method = "linear", f = 1)
> y_interpolated_backward <- newton_backward_interpolation(x_interpolate)
 > 
# Print the results
> cat("interpolated value using Newton's Forward Interpolation at x =", x_interpolate, ":", y_interpolated_forward,
"""
"\n")
Interpolated Value using Newton's Forward Interpolation at x = 3 : 7
> cat("Interpolated Value using Newton's Backward Interpolation at x =", x_interpolate, ":", y_interpolated_backward,
 Interpolated Value using Newton's Backward Interpolation at x = 3 : 7
```

Result:

Numerical integration using R

13. Numerical integration using Trapezoidal and Simpson's 1/3rd and 3/8th rules

Aim:

To write R program Numerical integration using Trapezoidal and Simpson's 1/3rd and 3/8th rules

Program:

```
# Define a function to be integrated f <- function(x) x^2 + 1

# Define integration limits a <- 0
b <- 2
```

Perform numerical integration using trapezoidal rule
n_intervals <- 4
trapezoidal_integral <- integrate(f, lower = a, upper = b, subdivisions = n_intervals)

Perform numerical integration using Simpson's 1/3 rule (automatic method selection) simpson_13_integral <- integrate(f, lower = a, upper = b, subdivisions = n_intervals)

Print the results cat("Trapezoidal Rule Integral:", trapezoidal_integral\$value, "\n") cat("Simpson's 1/3 Rule Integral:", simpson_13_integral\$value, "\n")

Output:

```
Console Jobs x

R R420 · J Perine a function to be integrated

f <- function(X) x^2 + 1

# Define integration limits

a <- 0

b Perform numerical integration using trapezoidal rule

n_intervals <- 4

trapezoidal_integral <- integrate(f, lower = a, upper = b, subdivisions = n_intervals)

# Perform numerical integration using simpson's 1/3 rule (automatic method selection)

s impson_13_integral <- integrate(f, lower = a, upper = b, subdivisions = n_intervals)

# Perform numerical integration using simpson's 1/3 rule (automatic method selection)

s impson_13_integral <- integrate(f, lower = a, upper = b, subdivisions = n_intervals)

# Perform numerical integration using simpson's 1/3 rule (automatic method selection)

s impson_13_integral <- integrate(f, lower = a, upper = b, subdivisions = n_intervals)

# Perint the results

cat("Trapezoidal Rule Integral:", trapezoidal_integralSvalue, "\n")

Trapezoidal Rule Integral: 4.666667

| Integral | Int
```

Result:

Solution of Ordinary differential equations using R

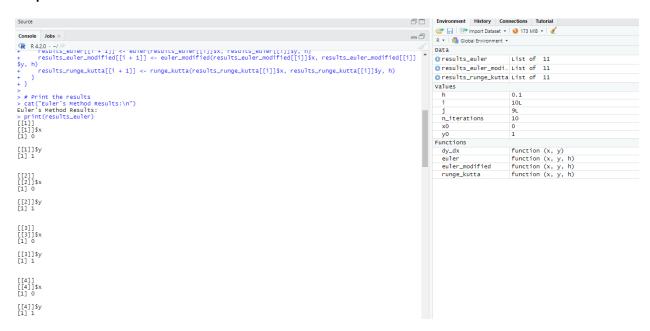
14. Euler's method, Euler's modified method, Runge-Kutta methods

Aim:

To write R program Euler's method, Euler's modified method, Runge-Kutta methods

```
# Define a differential equation
dy_dx \leftarrow function(x, y) - 2 * x * y
# Define initial values
x0 <- 0
y0 < -1
h <- 0.1 # Step size
# Euler's method
euler <- function(x, y, h) {
y_new <- y + h * dy_dx(x, y)
return(list(x_new = x + h, y_new = y_new))
}
# Euler's modified method
euler_modified <- function(x, y, h) {</pre>
y_prime <- y + h * dy_dx(x, y)
y_new <- y + 0.5 * h * (dy_dx(x, y) + dy_dx(x + h, y_prime))
 return(list(x_new = x + h, y_new = y_new))
}
# Runge-Kutta method (4th order)
runge_kutta <- function(x, y, h) {</pre>
k1 <- h * dy_dx(x, y)
 k2 <- h * dy_dx(x + 0.5 * h, y + 0.5 * k1)
 k3 <- h * dy_dx(x + 0.5 * h, y + 0.5 * k2)
 k4 <- h * dy_dx(x + h, y + k3)
y_new <- y + (1/6) * (k1 + 2 * k2 + 2 * k3 + k4)
 return(list(x_new = x + h, y_new = y_new))
}
# Perform iterations using each method
n iterations <- 10
results_euler <- list()
results_euler_modified <- list()
results_runge_kutta <- list()
```

```
for (i in 1:n_iterations) {
 results_euler[[i]] <- list(x = x0, y = y0)
 results euler modified[[i]] <- list(x = x0, y = y0)
 results_runge_kutta[[i]] <- list(x = x0, y = y0)
 for (j in 1:(n_iterations - 1)) {
  results_euler[[i + 1]] <- euler(results_euler[[i]]$x, results_euler[[i]]$y, h)
  results_euler_modified[[i + 1]] <- euler_modified(results_euler_modified[[i]]$x,
results_euler_modified[[i]]$y, h)
  results_runge_kutta[[i + 1]] <- runge_kutta(results_runge_kutta[[i]]$x, results_runge_kutta[[i]]$y, h)
 }
}
# Print the results
cat("Euler's Method Results:\n")
print(results_euler)
cat("Euler's Modified Method Results:\n")
print(results_euler_modified)
cat("Runge-Kutta Method Results:\n")
print(results_runge_kutta)
```



Result: