```
# C-PROGRAM
```

PROGRAM IN C FOR ADAMS-MOULTON METHOD

```
#include <stdio.h>
// Define the function f(x, y)
double f(double x, double y) {
  return 3 * x * x * (1 + y);
}
// Runge-Kutta method to compute y-values
void rungeKutta(double x0, double y0, double h, int steps, double y_values[]) {
  double x = x0, y = y0;
  // Store the initial y-value
  y_values[0] = y0;
  for (int i = 1; i \le steps; i++) {
    double k1 = h * f(x, y);
    double k2 = h * f(x + h / 2.0, y + k1 / 2.0);
    double k3 = h * f(x + h / 2.0, y + k2 / 2.0);
    double k4 = h * f(x + h, y + k3);
    y += (k1 + 2 * k2 + 2 * k3 + k4) / 6.0;
    x += h;
    y_values[i] = y; // Store the computed y-value
  }
}
// Adams-Moulton method to compute y-values
void adams_moulton(double h, double t[], double y[], int n_points) {
```

```
for (int i = 3; i < n_points - 1; i++) {
     t[i + 1] = t[i] + h; // Next time step
     // Predictor: Adams-Bashforth 4-step formula
     double y_pred = y[i] + (h / 24.0) * (
       55 * f(t[i], y[i]) -
       59 * f(t[i - 1], y[i - 1]) +
       37 * f(t[i - 2], y[i - 2]) -
        9 * f(t[i - 3], y[i - 3])
     );
     // Corrector: Iterative Adams-Moulton formula
     double y_corr = y_pred;
     for (int iter = 0; iter < 5; iter++) { // Perform 5 iterations for refinement
       y_{corr} = y[i] + (h / 24.0) * (
          9 * f(t[i + 1], y_corr) +
          19 * f(t[i], y[i]) -
          5 * f(t[i - 1], y[i - 1]) +
          f(t[i - 2], y[i - 2])
       );
     }
    y[i + 1] = y_corr; // Update the solution
  }
int main() {
  // Step size
  double h = 0.05;
  // Initial conditions
```

```
double x0 = 0.0, y0 = 0.0;
// Number of steps to reach t = 0.2
int steps = 4;
// Time array and y-values array
double t[steps + 1];
double y[steps + 1];
// Initialize time steps
for (int i = 0; i \le steps; i++) {
  t[i] = x0 + i * h;
}
// Compute initial y-values using Runge-Kutta method
rungeKutta(x0, y0, h, steps, y);
// Print the computed y-values using Runge-Kutta
printf("Initial values computed using Runge-Kutta:\n");
for (int i = 0; i \le steps; i++) {
  printf("t = \%.2f, y = \%.6f\n", t[i], y[i]);
}
// Use Adams-Moulton for further refinement if needed
adams_moulton(h, t, y, steps + 1);
// Print the final result
printf("\nApproximate value of y(0.2): %.6f\n", y[steps]);
return 0;
```

```
PROGRAM FOR RK METHOD
#include <stdio.h>
// Define the function f(x, y)
double f(double x, double y) {
  return 3 * x * x * (1 + y);
}
// Runge-Kutta method to compute y-values
void rungeKutta(double x0, double y0, double h, int steps, double y_values[]) {
  double x = x0, y = y0;
  // Store the initial y-value
  y_values[0] = y0;
  for (int i = 1; i \le steps; i++) {
    double k1 = h * f(x, y);
    double k2 = h * f(x + h / 2.0, y + k1 / 2.0);
    double k3 = h * f(x + h / 2.0, y + k2 / 2.0);
    double k4 = h * f(x + h, y + k3);
    y += (k1 + 2 * k2 + 2 * k3 + k4) / 6.0;
    x += h;
    y_values[i] = y; // Store the computed y-value
```

// Adams-Moulton method to compute y-values

}

```
void adams_moulton(double h, double t[], double y[], int n_points) {
  for (int i = 3; i < n_points - 1; i++) {
     t[i + 1] = t[i] + h; // Next time step
     // Predictor: Adams-Bashforth 4-step formula
     double y_pred = y[i] + (h / 24.0) * (
       55 * f(t[i], y[i]) -
       59 * f(t[i - 1], y[i - 1]) +
       37 * f(t[i - 2], y[i - 2]) -
        9 * f(t[i - 3], y[i - 3])
     );
     // Corrector: Iterative Adams-Moulton formula
     double y_corr = y_pred;
     for (int iter = 0; iter < 5; iter++) { // Perform 5 iterations for refinement
       y_{corr} = y[i] + (h / 24.0) * (
         9 * f(t[i + 1], y_corr) +
         19 * f(t[i], y[i]) -
         5 * f(t[i - 1], y[i - 1]) +
         f(t[i - 2], y[i - 2])
       );
     }
     y[i + 1] = y_corr; // Update the solution
  }
}
int main() {
  // Step size
  double h = 0.05;
```

```
// Initial conditions
 double x0 = 0.0, y0 = 0.0;
 // Number of steps to reach t = 0.2
 int steps = 4;
// Time array and y-values array
 double t[steps + 1];
 double y[steps + 1];
 // Initialize time steps
 for (int i = 0; i \le steps; i++) {
   t[i] = x0 + i * h;
 }
 // Compute initial y-values using Runge-Kutta method
 rungeKutta(x0, y0, h, steps, y);
 // Print the computed y-values using Runge-Kutta
 printf("Initial values computed using Runge-Kutta:\n");
 for (int i = 0; i \le steps; i++) {
    printf("t = %.2f, y = %.6f\n", t[i], y[i]);
 }
 // Use Adams-Moulton for further refinement if needed
 adams_moulton(h, t, y, steps + 1);
 // Print the final result
 printf("\nApproximate value of y(0.2): %.6f\n", y[steps]);
 return 0;
```

```
}
PROGRAM FOR EULERS METHOD
#include <stdio.h>
// Define the function f(x, y)
double f(double x, double y) {
  // Example differential equation: dy/dx = x + y
  return x + y;
}
// Euler's Method Function
void eulerMethod(double x0, double y0, double h, int steps) {
  double x = x0, y = y0;
  // Print table header
  printf("\nStep\t x\t\t y\n");
  printf("----\n");
  // Iterative calculation using Euler's method
  for (int i = 0; i \le steps; i++) {
    printf("%d\t %.6f\t %.6f\n", i, x, y);
    // Update values using Euler's formula
    y = y + h * f(x, y);
    x = x + h;
  }
}
```

// Main function

```
int main() {
  // Declare variables
  double x0, y0, h;
  int steps;
  // Input initial conditions
  printf("Enter initial value of x (x0): ");
  scanf("%lf", &x0);
  printf("Enter initial value of y (y0): ");
  scanf("%lf", &y0);
  // Input step size
  printf("Enter step size (h): ");
  scanf("%lf", &h);
  // Input number of steps
  printf("Enter number of steps: ");
  scanf("%d", &steps);
  // Confirm inputs
  printf("\nInitial conditions: x0 = \%.6f, y0 = \%.6f\n", x0, y0);
  printf("Step size: h = \%.6f\n'', h);
  printf("Number of steps: %d\n\n", steps);
  // Solve using Euler's method
  printf("Solving the differential equation using Euler's Method...\n");
  eulerMethod(x0, y0, h, steps);
  printf("\nSolution completed.\n");
```

```
return 0;
#include <stdio.h>
#include <stdlib.h>
// Function defining the ODE: y' = f(t, y)
double f(double t, double y) {
  return t - y; // Example: Replace this with your ODE
}
// Function to initialize the first few steps using Euler's method
void euler_method(double t[], double y[], int initial_steps, double h) {
  for (int i = 1; i < initial_steps; i++) {
    t[i] = t[i - 1] + h;
    y[i] = y[i - 1] + h * f(t[i - 1], y[i - 1]);
  }
}
// Adams-Bashforth coefficients for up to 4 steps
void get_ab_coefficients(int steps, double coeff[]) {
  switch (steps) {
    case 1:
       coeff[0] = 1.0;
       break;
    case 2:
       coeff[0] = 3.0 / 2.0;
       coeff[1] = -1.0 / 2.0;
       break;
    case 3:
       coeff[0] = 23.0 / 12.0;
       coeff[1] = -16.0 / 12.0;
```

```
coeff[2] = 5.0 / 12.0;
       break;
    case 4:
       coeff[0] = 55.0 / 24.0;
       coeff[1] = -59.0 / 24.0;
       coeff[2] = 37.0 / 24.0;
       coeff[3] = -9.0 / 24.0;
       break:
    default:
       printf("Unsupported number of steps: %d\n", steps);
       exit(1);
  }
}
// Adams-Bashforth multistep method
void adams_bashforth_multistep(double t0, double y0, double t_end, double h, int steps) {
  int n = (int)((t_end - t0) / h); // Total number of steps
  double t[n + 1], y[n + 1];
  double coeff[steps]; // Coefficients for the Adams-Bashforth method
  // Initialize time and solution arrays
  t[0] = t0;
  y[0] = y0;
  // Get the coefficients for the specified method
  get_ab_coefficients(steps, coeff);
  // Initialize the first few steps using Euler's method
  euler_method(t, y, steps, h);
  // Print the initial values
```

```
for (int i = 0; i < steps; i++) {
    printf
#include <stdio.h>
#include <stdlib.h>
#include <math.h>
void gaussian_elimination(int n, double a[n][n + 1], double x[n]) {
  for (int i = 0; i < n; i++) {
    // Partial Pivoting
     for (int k = i + 1; k < n; k++) {
       if (fabs(a[i][i]) < fabs(a[k][i])) {
          for (int j = 0; j \le n; j++) {
            double temp = a[i][j];
            a[i][j] = a[k][j];
            a[k][j] = temp;
          }
       }
     }
     // Forward Elimination
     for (int k = i + 1; k < n; k++) {
       double factor = a[k][i] / a[i][i];
       for (int j = 0; j \le n; j++) {
          a[k][j] -= factor * a[i][j];
       }
     }
  }
  // Back Substitution
  for (int i = n - 1; i \ge 0; i--) {
    x[i] = a[i][n];
```

```
for (int j = i + 1; j < n; j++) {
       x[i] -= a[i][j] * x[j];
    }
    x[i] /= a[i][i];
  }
}
int main() {
  int n = 3; // Example: Size of the system
  double a[3][4] = {
    {2, -1, 1, 3},
    {1, 3, 2, 12},
    {1, -1, 2, 2}
  }; // Example augmented matrix
  double x[3];
  gaussian_elimination(n, a, x);
  printf("Solution:\n");
  for (int i = 0; i < n; i++) {
    printf("x%d = %.4f\n", i + 1, x[i]);
  }
  return 0;
#include <stdio.h>
#include <math.h>
double f(double x) {
  return x * x - 4; // Example: f(x) = x^2 - 4
}
```

```
double f_prime(double x) {
  return 2 * x; // Derivative: f'(x) = 2x
}
void newton_raphson(double initial_guess, double tolerance, int max_iterations) {
  double x = initial_guess;
  for (int i = 0; i < max_iterations; i++) {</pre>
    double fx = f(x);
    double fpx = f_prime(x);
    if (fabs(fpx) < 1e-10) {
       printf("Derivative is too small. Method fails.\n");
       return;
    }
    double x_next = x - fx / fpx;
     printf("Iteration %d: x = \%.6f, f(x) = \%.6f\n", i + 1, x_next, f(x_next));
    if (fabs(x_next - x) < tolerance) {</pre>
       printf("Root found: x = \%.6f\n", x_next);
       return;
    }
    x = x_next;
  }
  printf("Maximum iterations reached. No solution found.\n");
}
int main() {
  double initial_guess = 2.0; // Starting point
  double tolerance = 1e-6;
  int max_iterations = 20;
```

```
newton_raphson(initial_guess, tolerance, max_iterations);
  return 0;
#include <stdio.h>
#include <math.h>
double f(double x) {
  return x * x; // Example: f(x) = x^2
}
double trapezoidal_rule(double a, double b, int n) {
  double h = (b - a) / n;
  double sum = f(a) + f(b);
  for (int i = 1; i < n; i++) {
    double x = a + i * h;
    sum += 2 * f(x);
  }
  return (h / 2) * sum;
}
int main() {
  double a = 0.0; // Start of the interval
  double b = 1.0; // End of the interval
  int n = 100; // Number of subintervals
  double result = trapezoidal_rule(a, b, n);
```

```
printf("Integral result: %.6f\n", result);
  return 0;
#include <stdio.h>
#include <math.h>
void matrix_vector_mult(int n, double A[n][n], double v[n], double result[n]) {
  for (int i = 0; i < n; i++) {
    result[i] = 0.0;
    for (int j = 0; j < n; j++) {
       result[i] += A[i][j] * v[j];
    }
  }
}
void normalize(int n, double v[n]) {
  double norm = 0.0;
  for (int i = 0; i < n; i++) {
    norm += v[i] * v[i];
  }
  norm = sqrt(norm);
  for (int i = 0; i < n; i++) {
    v[i] /= norm;
  }
}
double power_method(int n, double A[n][n], double v[n], int max_iterations, double tolerance) {
  double eigenvalue = 0.0;
  double prev_eigenvalue = 0.0;
```

```
for (int iter = 0; iter < max_iterations; iter++) {</pre>
    double w[n];
    matrix_vector_mult(n, A, v, w);
    normalize(n, w);
    prev_eigenvalue = eigenvalue;
    eigenvalue = 0.0;
    for (int i = 0; i < n; i++) {
       eigenvalue += w[i] * w[i];
    }
    for (int i = 0; i < n; i++) {
       v[i] = w[i];
    }
    if (fabs(eigenvalue - prev_eigenvalue) < tolerance) {</pre>
       return eigenvalue;
    }
  }
  return eigenvalue;
int main() {
  int n = 2; // Matrix size
  double A[2][2] = {
    {4, 1},
    {2, 3}
  };
  double v[2] = \{1, 1\}; // Initial guess
  int max_iterations = 1000;
  double tolerance = 1e-6;
```

```
double eigenvalue = power_method(n, A, v, max_iterations, tolerance);
printf("Dominant eigenvalue: %.6f\n", eigenvalue);
printf("Eigenvector: [%.6f, %.6f]\n", v[0], v[1]);
return 0;
}
```