

# FINITE ELEMENT ANALYSIS

PROF: P.M. MOHITE

PROJECT 1: 1D BAR

PROJECT 2: 2D BEAM

BY,

J. SIRISHA (ALONE)

231010031

# PROJECT-1

①

## Introduction

Objective: To find the solution of governing equation for the bar where it is constrained at both the ends A, B and traction load  $T(x)$  applied on the bar (1D)

To develop a generic Finite element code for the governing equation of bar without using inbuilt functions to yield displacements.

## Methodology. [Question 1-a]

### Mathematical Formulation

- The governing equation of the 1-D elastic bar considered is

Case 1:  $AE \frac{d^2 u}{dx^2} - 2u = T(x)$  where  $T(x) = 4x^2 - 2x - 4$  (Quadratic)  
↓  
Traction force or source term.

Stiffness variation  $AE(x)$  ;

Case 2:  $T(x) = \text{constant}$ .

### Boundary Conditions

$$\left. \begin{array}{l} u=0, \text{ at } x=0 \\ u \frac{du}{dx} = -1; \text{ at } x=1 \end{array} \right\} \text{Dirichlet B.C.}$$

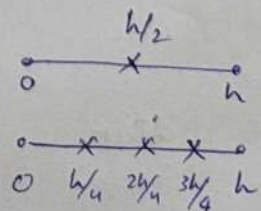
$$\left. \begin{array}{l} u=0 \text{ at } x=0 \\ \frac{du}{dx} = -3 \text{ at } x=1 \end{array} \right\} \text{Neumann B.C.}$$

## Finite element modeling

- Discretization of the domain into Elements.

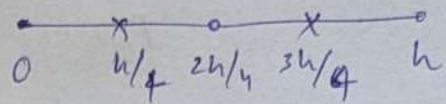
For element type (Linear) -  $e=2$   
OIF = 1

$e=4$





Element type - Quadratic  $e=2$   
 OIF=2



Trick

Pattern found

~~err~~

$$S = (\text{OIF} \times e) + 1$$

$\downarrow$  gives no. of nodes       $\uparrow$  no. of elements.  
 $\uparrow$  Orders of polynomial

for  $i=1, 2, 3$

location  $(i) = (i-1)h / s-1$  } gives the location of node.  
 end

## Shape functions

### Lagrange Interpolation Implementation

Consider

Element type - Quadratic [3 nodes]

$x_1, x_2, x_3$ .

Formula of Lagrange  $\Rightarrow$  for shape function at node 1  $\Rightarrow \frac{(x-x_2)(x-x_3)}{(x_1-x_2)(x_1-x_3)}$   
 $N_1 \Rightarrow$

Shape function at node 2  $N_2 = \frac{(x-x_1)(x-x_3)}{(x_2-x_1)(x_2-x_3)}$

$$N_3 = \frac{(x-x_1)(x-x_2)}{(x_3-x_1)(x_3-x_2)}$$

By substituting the values of  $x_1, x_2, x_3 \rightarrow -1, 0, 1$

we get,  $N_1 = \phi_1(x) = \frac{x(x-1)}{2}$

$$N_2 = \phi_2(x) = -(x-1)(x+1)$$

$$N_3 = \phi_3(x) = \frac{x(x+1)}{2}$$

For finding  $\left(\frac{dN}{dx}\right)$ , Trick Pattern found  $\Rightarrow$  Numerator is separately found and denominator as

$$\boxed{df = \frac{li}{den}} \quad (\text{refer code})$$

then  $\frac{dN_1}{dx} = d\phi_1(x) = \frac{x-1}{2}$ ;  $\frac{dN_2}{dx} = d\phi_2(x) = -2x$ ;  $\frac{dN_3}{dx} = d\phi_3(x) = \frac{x+1}{2}$



## Hierarchic shape functions

Element type - Linear  $\Rightarrow N_1 \Rightarrow \phi_1(\xi) = 0.5(1-\xi)$

$$N_2 \Rightarrow \phi_2(\xi) = 0.5(1+\xi)$$

Quadratic  $\Rightarrow N_1 \Rightarrow \phi_1(\xi) = \frac{\xi(\xi-1)}{2}$

$$N_2 \Rightarrow \phi_2(\xi) = \frac{-(\xi-1)}{\xi+1}$$

$$N_3 \Rightarrow \phi_3(\xi) = \frac{\xi(\xi+1)}{2}$$

For finding differentiation, directly manually entering.

Linear  $\Rightarrow \frac{dN_1}{d\xi} \Rightarrow d\phi_1(\xi) = -0.5$

Quadratic  $\Rightarrow \frac{dN_1}{d\xi} \Rightarrow d\phi_1(\xi) = \xi - 0.5$

$$\frac{dN_2}{d\xi} \Rightarrow d\phi_2(\xi) = 0.5$$

$$\frac{dN_2}{d\xi} \Rightarrow d\phi_2(\xi) = -2\xi$$

$$\frac{dN_3}{d\xi} \Rightarrow d\phi_3(\xi) = \xi + 0.5$$

## Implementation in MATLAB

→ Using Lagrange interpolation & hierarchic shape functions,

$$I = \int_{-1}^1 f(\xi) d\xi \approx \sum_{i=1}^N w_i f(\xi_i)$$

So, Using Gauss Quadrature, we get the values of  $\xi_i$  &  $w_i$

∴ we manually input the values such as,

$$n=1 \text{ (Linear)} ; n=2 \text{ (Quadratic)}$$

$$\phi_1(1) = 0 ;$$

$$\phi_1(1) = \sqrt{3}/3 ;$$

$$w(1) = 2 ;$$

$$\phi_2(2) = -\sqrt{3}/3 ;$$

$$w(1) = 1 ;$$

$$w(2) = 1 ;$$

So, on... till  $n=5$



These  $\phi_{1e}$  &  $\phi_{2e}$  are substituted in stiffness matrix and source term which also involves  $w(i)$  &  $p(i)$ . We find Global stiffness matrix, source term.

$$\text{Stiffness term} \Rightarrow \int_0^h \left( a \frac{d\phi_{1e}(1)}{dx} \frac{d\phi_{1e}(2)}{dx} - c \phi_{1e}(1) \phi_{1e}(2) \right) dx$$

$$\approx \sum_{e=1}^n \left[ a * (\text{diff-}\phi_{1e} \text{ values}(1)) * (\text{diff-}\phi_{1e} \text{ values}(2)) (2/h) - c * \phi_{1e} \text{ values}(1) * \phi_{1e} \text{ values}(2) (h/2) \right]$$

Similarly for force term.

These steps are same for Lagrange Interpolation and Hierarchic shape method.

## Results and Discussion

### Patch Test Results

- ① FEM solution and Exact solution of Dirichlet Boundary Conditions and Neumann Boundary Conditions have been plotted

[X, displacements]  
length of bar at each point

- ② Error in solution of FEM has been plotted (both Lagrange & Hierarchic)  
Error for FEM solution is calculated as,

$$\text{Error} = \text{abs}(\text{FEM-solution} - \text{displacement vector})$$



## Strain Energy Analysis

### ① Lagrange Interpolation

We know

$$U = \int \frac{1}{2} E \epsilon^2 dx$$

Young's modulus.

$$\epsilon = \frac{du}{dx} \rightarrow \text{displacement}$$

Strain

For our problem,  $u(x) = -2x^2 + x$ .

$$\frac{du}{dx} = -4x + 1$$

$$\epsilon^2 \Rightarrow (-4x + 1)^2$$

$$\Rightarrow U_{\text{exact}} = \int \frac{1}{2} E (-4x + 1)^2 dx$$

Using FEM

$$U(x) \approx \sum_{i=1}^n N_i(x) \cdot d_i$$

shape function      displacement

$$\epsilon \approx \sum_{i=1}^n \frac{dN_i}{dx} \cdot d_i$$

$$\epsilon^2 \Rightarrow \left( \sum_{i=1}^n \frac{dN_i}{dx} d_i \right)^2$$

$$U_{\text{FEM}} = \sum_{\substack{\text{elements} \\ i=1 \\ k=1}}^{N_{\text{elems}}} \left[ \frac{1}{2} E \left( \sum_{i=1}^n \frac{dN_i}{dx} d_i \right)^2 + c(d_i)^2 \right] dx$$

Using Gauss Quadrature, local strains = Sum [ (d shape + displacements) + c · d\_i ]

$$U_{\text{FEM}} = \frac{1}{2} E (\text{Local strains})^2 \cdot w(i) \cdot h$$

- Strain Energy is calculated over the range of elements (1 to 10) and observed the convergence.

- In Lagrange & Hierarchic shape functions, Only shape functions calculations are different.

[Quadratic]

Lagrange  $\Rightarrow$   $N_1(x) = \frac{(x-x_2)(x-x_3)}{(x_1-x_2)(x_1-x_3)}$   $N_3(x) = \frac{(x-x_1)(x-x_2)}{(x_3-x_1)(x_3-x_2)}$

$N_2(x) = \frac{(x-x_1)(x-x_3)}{(x_2-x_1)(x_2-x_3)}$

Hierarchic  $\Rightarrow$   $N_1(\xi) = 0.5 \xi (\xi - 1)$

$N_2(\xi) = 1 - \xi^2$

$N_3(\xi) = \frac{1}{2} \xi (\xi + 1)$

### Error Analysis of Convergence Rate

$$\text{Error} = \text{abs}(\text{Strain-Energy-FEM} - \text{Strain-Energy-Exact})$$

Strain Energy Exact = 1.6608

Strain Energy FEM = 1.2708

} for our problem.

Error of strain Energy at each element <sup>from</sup>  $[1:10]$  has been plotted.

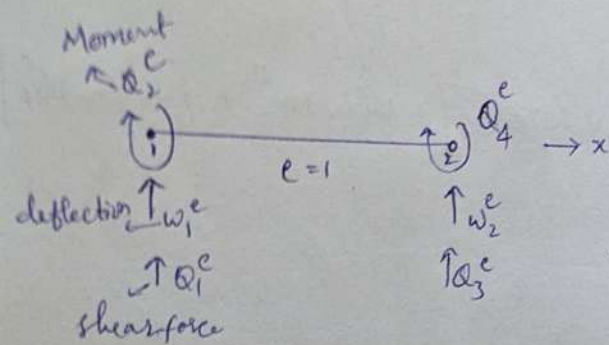
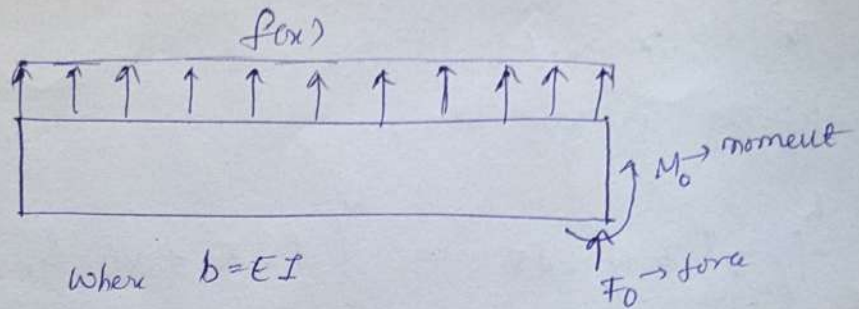
Error between FEM & Exact strain energy.



# BEAM BENDING PROBLEM

## ① Governing Equation

$$\frac{d^2}{dx^2} \left( b \frac{d^2 w}{dx^2} \right) + f(x)$$



Sign convention

$w \uparrow +ve$   
 shear force  $Q_1 \uparrow +ve$   
 Moment  $Q_2 \curvearrowright +ve$

## ② Development of a weak form

$$\int_0^l v \left[ \frac{d^2}{dx^2} \left( b \frac{d^2 w}{dx^2} \right) - f \right] dx = 0$$

(Performing Integration by Parts) (2 times)

$$\int_0^l \left[ - \frac{dv}{dx} \left[ b \frac{d^2 w}{dx^2} \right]' - v f \right] dx + \left[ v + \frac{d}{dx} \left( b \frac{d^2 w}{dx^2} \right) \right]_0^l = 0$$

$$\int_0^l \left[ \frac{d^2 v}{dx^2} b \frac{d^2 w}{dx^2} - v f \right] dx + \underbrace{\left[ v (b w'')' - b w'' v' \right]_0^l}_{\text{boundary terms}} = 0$$

Weak form of the governing Equation

Where  $v$  determines  $\rightarrow w$   
 $-v'$  determines  $\rightarrow -w' \rightarrow -\frac{dw}{dx}$  } Primary variables  
 replacing with  $v$  &  $v'$   
 Slope.

- Coefficient of primary variables helps determining boundary variables.

$(b w'')' \rightarrow$  coefficient of  $v$  (represents shear-force)

$b w'' \rightarrow$  coefficient of  $v'$  (represents moment).

So, we have four boundary conditions

$v, -v',$  shear force and moments.



$w, -dw/dx$  } called generalized displacements

$(bw'')', (bw'')$  } called generalised forces.

$$\bar{x} = x - x_{\text{node}}$$

By sub in weak form,

$$\int_0^{h_e} \underbrace{v'' \cdot b \cdot w''}_{\text{Bilinear term } B(v, w)} d\bar{x} = \int_0^{h_e} f v d\bar{x} + v(0) Q_1^e + [-v'(0) Q_2^e] + v(h_e) Q_3^e + [-v'(h_e) Q_4^e]$$

$$Q_1^e = (bw'')'|_0 \quad Q_3^e = -(bw'')'|_{h_e}$$

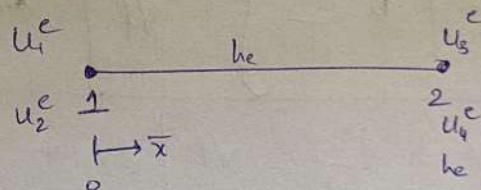
$$Q_2^e = (bw'')|_0 \quad Q_4^e = -(bw'')|_{h_e}$$

$$I(w) = \int_0^{h_e} \frac{1}{2} b(w'')^2 d\bar{x} - \int_0^{h_e} w f d\bar{x} - w(0) Q_1^e - [-w'(0) Q_2^e] - w(h_e) Q_3^e - [-w'(h_e) Q_4^e]$$

### ③ Interpolating Functions.

$u_1, u_3$  - Translational.

$u_2, u_4 \rightarrow$  Rotational.



Four D.O.F for 1 element

$$\begin{aligned} w(0) = w_1 = u_1^e &\rightarrow \text{generalized rotation for DOF} \\ -w'(0) = \theta_1 = u_2^e &\quad \left. \begin{array}{l} \text{slope} \end{array} \right\} \begin{array}{l} 2 \text{ DOF} \\ \text{at node 1} \end{array} \end{aligned}$$

$$\begin{aligned} w(h_e) = w_2 = u_3^e \\ w'(h_e) = \theta_2 = u_4^e \end{aligned} \quad \left. \begin{array}{l} \end{array} \right\} \begin{array}{l} 2 \text{ DOF} \\ \text{at node 2} \end{array}$$

$$w^e(x) = \sum_{j=1}^4 u_j^e \phi_j^e(x) \Rightarrow u_1^e \phi_1^e(\bar{x}) + u_2^e \phi_2^e(\bar{x}) + u_3^e \phi_3^e(\bar{x}) + u_4^e \phi_4^e(\bar{x})$$

Approximate function to be found.

$$w^e(0) = u_1, \quad -w^e'(0) = u_2; \quad w^e(h_e) = u_3; \quad w^e'(h_e) = u_4$$



$$\begin{bmatrix} k_{11}^e & k_{12}^e & k_{13}^e & k_{14}^e \\ k_{21}^e & k_{22}^e & k_{23}^e & k_{24}^e \\ k_{31}^e & k_{32}^e & k_{33}^e & k_{34}^e \\ k_{41}^e & k_{42}^e & k_{43}^e & k_{44}^e \end{bmatrix} \begin{Bmatrix} u_1 \\ u_2 \\ u_3 \\ u_4 \end{Bmatrix} = \begin{Bmatrix} f_1^e + Q_1^e \\ f_2^e + Q_2^e \\ f_3^e + Q_3^e \\ f_4^e + Q_4^e \end{Bmatrix}$$

due to point load & moments

due to distributed load

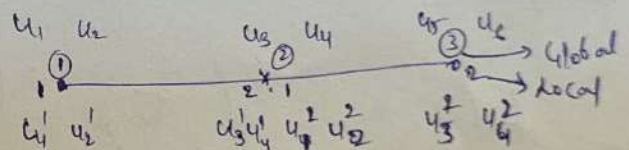
After substitution,

$$[K] = \frac{2b}{h^3} \begin{bmatrix} 6 & -3h & -6 & -3h \\ -3h & 2h^2 & 3h & h^2 \\ -6 & 3h & 6 & 3h \\ -3h & h^2 & 3h & 2h^2 \end{bmatrix} ; \{f\} = \frac{fh}{12} \begin{Bmatrix} 6 \\ -h \\ 6 \\ -h \end{Bmatrix}$$

due to symmetry, 1, 3 are same, so as 2 & 4 now.

#### ④ Assembly

$$\begin{aligned} u_3^1 &= u_1^2 = u_3 \\ u_4^1 &= u_2^2 = u_4 \end{aligned} \quad \text{Symmetry.}$$



Replacing,  $u_1^1 \Rightarrow u_1 ; u_2^1 \Rightarrow u_2 ; u_3^1 \Rightarrow u_5 ; u_4^1 = u_6$

Force balance

$$f_3^1 + f_1^2 = f_3 ; f_4^1 + f_2^2 = f_4$$

$$Q_3^1 + Q_1^2 = Q_3 ; Q_4^1 + Q_2^2 = Q_4$$

For element 1,

$$\begin{bmatrix} k_{11}^1 & k_{12}^1 & k_{13}^1 & k_{14}^1 \\ k_{21}^1 & k_{22}^1 & k_{23}^1 & k_{24}^1 \\ k_{31}^1 & k_{32}^1 & k_{33}^1 & k_{34}^1 \\ k_{41}^1 & k_{42}^1 & k_{43}^1 & k_{44}^1 \end{bmatrix} \begin{Bmatrix} u_1^1 \\ u_2^1 \\ u_3^1 \\ u_4^1 \end{Bmatrix} = \begin{Bmatrix} f_1^1 \\ f_2^1 \\ f_3^1 \\ f_4^1 \end{Bmatrix} + \begin{Bmatrix} Q_1^1 \\ Q_2^1 \\ Q_3^1 \\ Q_4^1 \end{Bmatrix}$$

Add eq 1 & 5

Add eq 2 & 6

For element 2

$$\begin{bmatrix} k_{11}^2 & k_{12}^2 & k_{13}^2 & k_{14}^2 \\ k_{21}^2 & k_{22}^2 & k_{23}^2 & k_{24}^2 \\ k_{31}^2 & k_{32}^2 & k_{33}^2 & k_{34}^2 \\ k_{41}^2 & k_{42}^2 & k_{43}^2 & k_{44}^2 \end{bmatrix} \begin{Bmatrix} u_1^2 \\ u_2^2 \\ u_3^2 \\ u_4^2 \end{Bmatrix} = \begin{Bmatrix} f_1^2 \\ f_2^2 \\ f_3^2 \\ f_4^2 \end{Bmatrix} + \begin{Bmatrix} Q_1^2 \\ Q_2^2 \\ Q_3^2 \\ Q_4^2 \end{Bmatrix}$$



$$\begin{bmatrix}
 k_{11} & k_{12} & k_{13} & k_{14} & 0 & 0 \\
 k_{21} & k_{22} & k_{23} & k_{24} & 0 & 0 \\
 k_{31} & k_{32} & k_{33} + k_{11}^2 & k_{34} + k_{12}^2 & k_{13}^2 & k_{14}^2 \\
 k_{41} & k_{42} & k_{43} + k_{21}^2 & k_{44} + k_{22}^2 & k_{23}^2 & k_{24}^2 \\
 0 & 0 & k_{31}^2 & k_{32}^2 & k_{33}^2 & k_{34}^2 \\
 0 & 0 & k_{41}^2 & k_{42}^2 & k_{43}^2 & k_{44}^2
 \end{bmatrix}
 \begin{Bmatrix}
 u_1 \\
 u_2 \\
 u_3 \\
 u_4 \\
 u_5 \\
 u_6
 \end{Bmatrix}
 =
 \begin{Bmatrix}
 f_1 \\
 f_2 \\
 f_3 \rightarrow f_3^1 + f_1^2 \\
 f_4 \rightarrow f_4^1 + f_2^2 \\
 f_5 \\
 f_6
 \end{Bmatrix}
 \begin{Bmatrix}
 Q_1 \\
 Q_2 \\
 Q_3 \rightarrow Q_3^1 + Q_1^2 \\
 Q_4 \rightarrow Q_4^1 + Q_2^2 \\
 Q_5 \\
 Q_6
 \end{Bmatrix}
 \quad (3)$$

Assembled matrix is also symmetric.

### ⑤ Applying Boundary Conditions

In our case,

At  $x=0$

$$w = w' = 0$$

At  $x=L$

$$Q_5 = F_0 ; \quad Q_6 = -M_0$$

$$Q_3^1 + Q_1^2 = 0$$

$$Q_4^1 + Q_2^2 = 0$$

$$u_1 = u_2 = 0$$

$\therefore$  we get

$$\begin{Bmatrix}
 u_3 \\
 u_4 \\
 u_5 \\
 u_6
 \end{Bmatrix}
 = \frac{h}{6EI}
 \begin{Bmatrix}
 5h^2 F_0 + 3h M_0 + 17/4 f_0 h^3 \\
 -9h F_0 - 6M_0 - 7f_0 h^2 \\
 16h^2 F_0 + 12h M_0 + 12f_0 h^3 \\
 -12h F_0 - 12M_0 - 8f_0 h^2
 \end{Bmatrix}$$

$$\therefore w(x) = \begin{cases} u_3 \phi_1' + u_4 \phi_1'' & , \quad 0 \leq x \leq h \\ u_3 \phi_1^2 + u_4 \phi_2^2 + u_5 \phi_3^2 + u_6 \phi_4^2 & , \quad h \leq x \leq 2h \end{cases}$$

$$M = EI \frac{d^2 w}{dx^2} = EI \sum_{j=1}^4 u_j^T \frac{d^2 \phi_j^T}{dx^2}$$

Bending stress  $\rightarrow$

$$\sigma_x = \mp \frac{M y}{2I} = \mp \frac{EI}{2} \sum_{j=1}^4 u_j^T \frac{d^2 \phi_j^T}{dx^2}$$

Height of the beam



## Bending Moment and shear force calculation

$$B.M \Rightarrow \boxed{M = -EI \frac{d^2 w}{dx^2}}$$

$\frac{d^2 w}{dx^2}$  is solved using central difference.

$$\text{i.e., } \frac{d^2 w}{dx^2} \Rightarrow \frac{\text{displacements } (2(n+1)) - 2 \times \text{displacements } (2n) + \text{displacements } (2(n-1)))}{2^2 \cdot el}$$

Similarly for shear force,

$$\frac{dw}{dx} = \frac{\text{displacements } (2(i+1)) - \text{displacements } (2(i-1)))}{2 \cdot el}$$

$$V = -EI \frac{dw}{dx} / el \Rightarrow \boxed{V = \frac{dM}{dx}}$$



```

clc;
clear;
syms x;

% Input parameters
OIF = 2; % order of polynomial
e = 2; % number of elements
n = OIF + 1; % number of nodes
s = (OIF * e) + 1; % Global matrix size pattern

% governing_eqn = du^2/dx^2-2*u = T(x)
T = 4*x^2 - 2*x - 4; % Source term
T_x = sym(zeros(s, 1));

exact_sol_T = -2*x^2 + x;

% Assign numerical values
a_num = 1;
c_num = -2;
x0 = 0;
xn = 1;
h = (xn - x0) / e;

GP = n; % Number of Gauss points

[w, psi] = Guass(GP);
phi_e_values = zeros(GP, n);
diff_phi_e_values = zeros(GP, n);

for i = 1:GP
    psi_values = psi(i);
    [n, phi_e, dshap] = ShapeFunctions(OIF, psi_values);

    phi_e_values(i, :) = phi_e';
    diff_phi_e_values(i, :) = dshap';
end

% w
% psi
%
phi_e_values;
diff_phi_e_values;

% storing KNM and CNM values
K_Local = zeros(n, n);
C_Local = zeros(n, n);

% Assemble local matrices using Gauss quadrature
for i = 1:n
    for j = 1:n
        K_Local(i, j) = 0;
    end
end

```



```

C_Local(i, j) = 0;
for k = 1:GP
    % Define symbolic expressions for F_xi and CNM_integrand
    F_xi = a_num * diff_phi_e_values(k, i) * diff_phi_e_values(k, j)*(2/h) - c_num * phi_e_values(k, i) * phi_e_values(k, j)*(h/2);
    CNM_integrand = -phi_e_values(k, i) * phi_e_values(k, j)*(h/2);

    % Evaluate the expressions with numerical values
    F_xi_eval = double(F_xi);
    CNM_integrand_eval = double(CNM_integrand);

    % Compute the contribution using Gauss quadrature
    F_xi_contribution = F_xi_eval * w(k);
    CNM_contribution = CNM_integrand_eval * w(k);

    % Update the local stiffness matrix and CNM matrix
    K_Local(i, j) = K_Local(i, j) + F_xi_contribution;
    C_Local(i, j) = C_Local(i, j) + CNM_contribution;
end
end

% K_Local
% C_Local

% Calculation of global stiffness matrix
global_stiffness_size = s;
s = (OIF * e) + 1;

K_global = zeros(global_stiffness_size);
C_global = zeros(global_stiffness_size);

% Loop through each element
for element = 1:n-1:s-1
    % Assemble local stiffness matrix into global stiffness matrix
    start_index = element;
    end_index = element+(n-1);
    K_global(start_index:end_index, start_index:end_index) = K_global(start_index:end_index, start_index:end_index) + K_Local(element, element+(n-1));
    C_global(start_index:end_index, start_index:end_index) = C_global(start_index:end_index, start_index:end_index) + C_Local(element, element+(n-1));
end

% K_global
% C_global

X = (x0:h/OIF:xn)'; % Defining the nodal points along the domain

for i = 1:s
    T_x(i) = subs(T, x, X(i));
end

% Dirichlet Boundary Conditions
% Boundary condition 1
C = C_global * T_x;
du_dx = -3;

```



```

K_global_Dirichlet = K_global;
g_x_Dirichlet = sym(zeros(s, 1));
GNM_Dirichlet = a_num * g_x_Dirichlet * du_dx;
RHS_Dirichlet = GNM_Dirichlet + C;

u0 = x0;
un = -xn;
K_global_Dirichlet(1, :) = [1, zeros(1, s-1)];
K_global_Dirichlet(end, :) = [zeros(1, s-1), 1];
RHS_Dirichlet(1) = u0; % RHS of equation
RHS_Dirichlet(end) = un; % RHS of equation

```

```

% finding displacements using dirichet boundar conditions

```

```

u_D_bc = inv(K_global_Dirichlet)* RHS_Dirichlet;
displacement_vector = double(u_D_bc)

```

```

displacement_vector = 5×1
    0
    0.1250
    0
   -0.3750
   -1.0000

```

```

exact_values = double(subs(exact_sol_T, x, X)); % Evaluate exact solution at node positions
errors = abs(exact_values - displacement_vector); % Calculate absolute error at each node

```

```

% Plot the FEM Solution and the actual solution

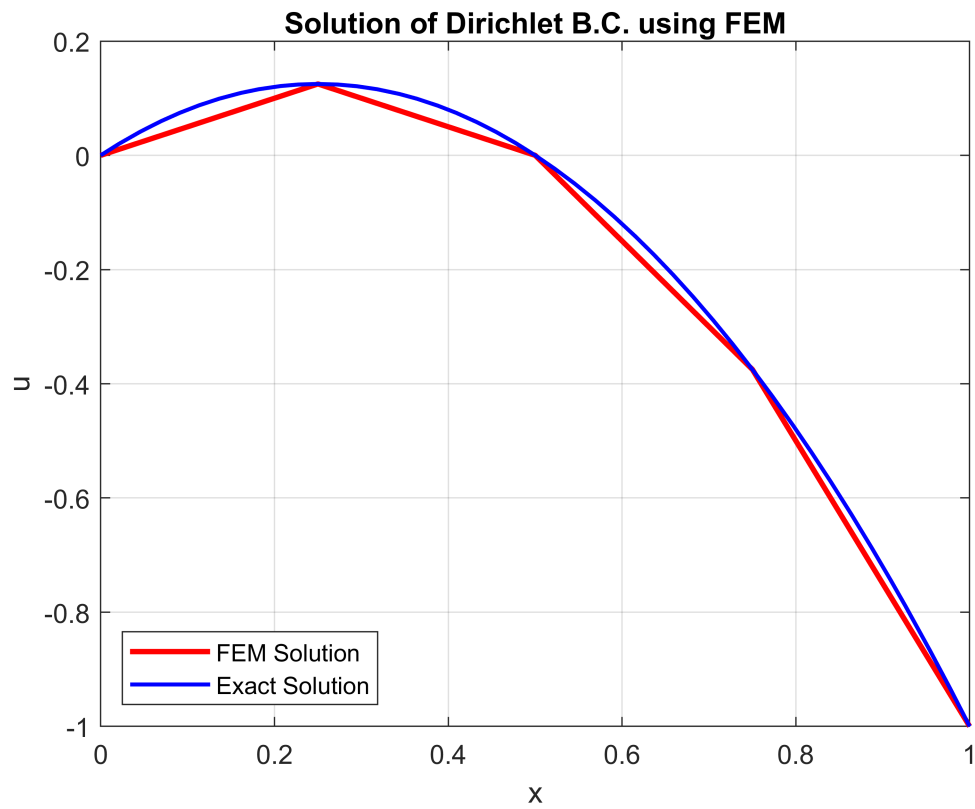
```

```

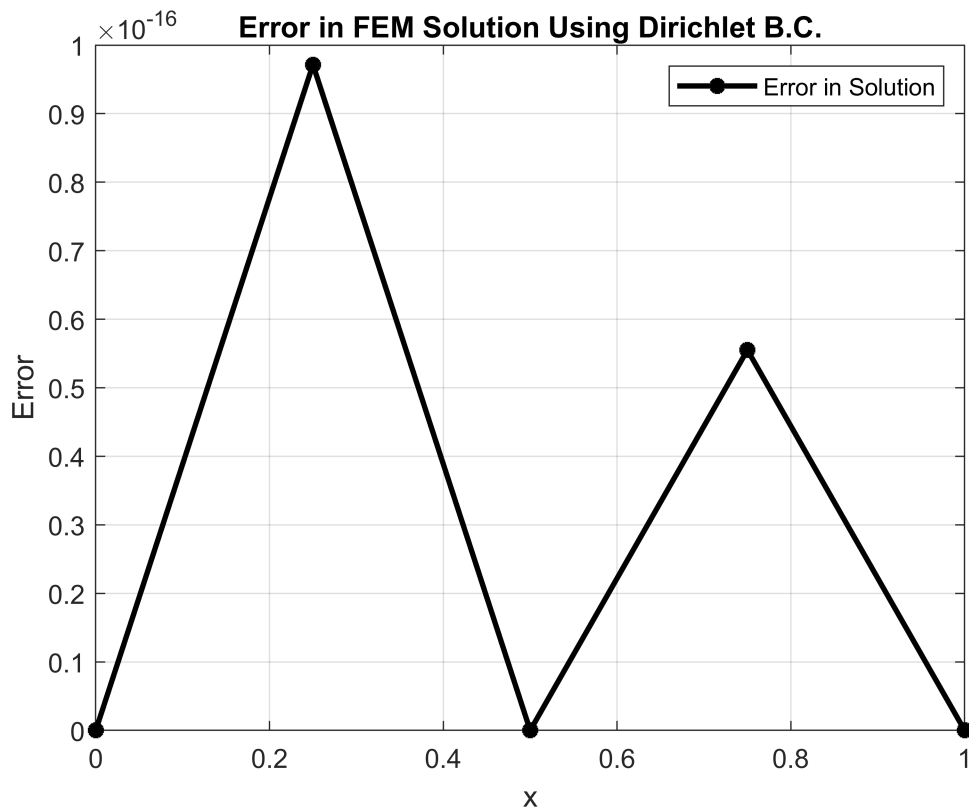
figure;
plot(X, displacement_vector, "r", "LineWidth", 2, "DisplayName", "FEM Solution");
hold on;
fplot(exact_sol_T, [0 1], "b", "LineWidth", 1.5, "DisplayName", "Exact Solution");
xlabel("x");
ylabel("u");
title("Solution of Dirichlet B.C. using FEM");
legend('Location', 'southwest');
grid on;

```





```
figure;  
plot(X, errors, 'k-*', 'LineWidth', 2, 'DisplayName', 'Error in Solution');  
xlabel('x');  
ylabel('Error');  
title('Error in FEM Solution Using Dirichlet B.C.');
```



```
% Neumann boundary Conditions
K_global_neumann = K_global;
g_x_neumann = sym(zeros(s, 1));
g_x_neumann(end)=1;

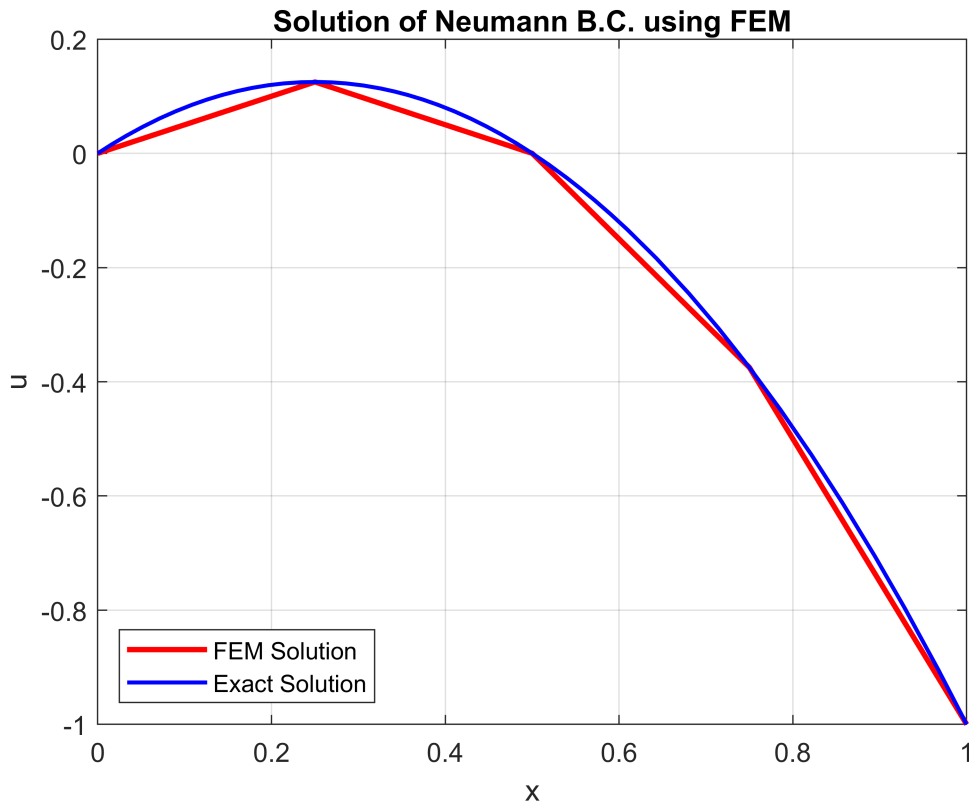
GNM_neumann= a_num * g_x_neumann * du_dx;
RHS_neumann = GNM_neumann+C;

% Boundary Conditions 2

K_global_neumann(1,:)= [1,zeros(1, s-1)];
RHS_neumann(1) = u0;
u_neumann = inv(K_global_neumann)*RHS_neumann;
displacement_vector_n = double(u_neumann);

%Plot the FEM Solution and the actual solution
figure;
plot(X, displacement_vector_n, "r", "LineWidth", 2, "DisplayName", "FEM Solution");
hold on;
fplot(x,exact_sol_T, [0 1], "b", "LineWidth", 1.5, "DisplayName", "Exact Solution");
xlabel("x");
ylabel("u");
title("Solution of Neumann B.C. using FEM");
legend('Location', 'southwest');
grid on;
```





```
function [w, psi] = Guass(p_bar)
    psi = zeros(p_bar, 1);
    w = zeros(p_bar, 1);

    if p_bar == 1
        psi(1) = 0; % for p_bar = 1
        w(1) = 2;
    elseif p_bar == 2
        psi(1) = sqrt(3)/3;
        psi(2) = -sqrt(3)/3;
        w(1) = 1;
        w(2) = 1;
    elseif p_bar == 3
        psi(1) = sqrt(15)/5;
        psi(2) = -sqrt(15)/5;
        psi(3) = 0;
        w(1) = 5/9;
        w(2) = 5/9;
        w(3) = 8/9;
    elseif p_bar == 4
        psi(1) = sqrt((15 - 2*sqrt(30))/35);
        psi(2) = -sqrt((15 - 2*sqrt(30))/35);
        psi(3) = (sqrt((14*sqrt(30))/5 + 21))/7;
        psi(4) = -(sqrt((14*sqrt(30))/5 + 21))/7; % for p_bar = 7
```

```

        w(1) = 1/2 + sqrt(30)/36;
        w(2) = 1/2 + sqrt(30)/36;
        w(3) = 1/2 - sqrt(30)/36;
        w(4) = 1/2 - sqrt(30)/36;
    elseif p_bar == 5
        psi(1) = 0.906179845938664;
        psi(2) = -0.906179845938664;
        psi(3) = 0.538469310105683;
        psi(4) = -0.538469310105683;
        psi(5) = 0;
        w(1) = 0.236926885056189;
        w(2) = 0.236926885056189;
        w(3) = 0.478628670499366;
        w(4) = 0.478628670499366;
        w(5) = 0.568888888888889;
    end
end

function [n, phi_e, dshap] = ShapeFunctions(OIF, x_loc)
    syms x;
    n = OIF + 1;
    x_local = sym(zeros(1, n));

    for i = 1:n
        x_local(i) = -1 + (i-1)*(2/OIF); % Calculate the elements of the vector
    end

    x_s = sym('x', [1, n]);
    N = sym(zeros(size(x_s)));
    df = sym(zeros(size(x_s))); % Initialize df

    for j = 1:n
        Li = 0;
        den = 1;
        product_term = 1;
        for i = 1:n
            if j ~= i
                term = 1; % Initialize product term for index j
                for k = 1:n
                    if k ~= i && k ~= j
                        term = term * (x - x_s(k));
                    end
                end
                % Add the product term for index j to Lagrange polynomial for index i
                Li = Li + term;
            end
        end
        %% creating pattern for differentiation
        if i ~= j
            product_term = product_term * (x - x_s(i)) / (x_s(j) - x_s(i));
            den = den * (x_s(j) - x_s(i));
        end
        N(j) = product_term;
    end

```



```

        df(j) = Li / den;
    end

    phi_e = subs(N, x_s, x_local);
    dshap = subs(df, x_s, x_local);

    % Evaluate the shape functions and their derivatives at the Gauss points
    xi_loc = x_loc;
    phi_e = double(subs(phi_e, x, xi_loc));
    dshap = double(subs(dshap, x, xi_loc));
end

```

```

clc;
clear;
syms x;

% Input parameters
OIF = 2; % order of polynomial
e = 2; % number of elements
n = OIF + 1; % number of nodes
s = (OIF * e) + 1; % Global matrix size

T = 4*x^2 - 2*x - 4; % Source term
T_x = sym(zeros(s, 1));

exact_sol_T = -2*x^2 + x;

% Domain definitions
a_num = 1;
c_num = -2;
x0 = 0;
xn = 1;
h = (xn - x0) / e;

% Number of Gauss points
GP = n;

% Initialize matrices for values and derivatives
phi_e_values = zeros(GP, n);
diff_phi_e_values = zeros(GP, n);

% Calculate Gauss points and weights manually
[w, psi] = Gauss(GP);

% Define hierarchical shape functions and their derivatives
for i = 1:GP
    psi_values = psi(i);
    [phi_e, dphi_e] = HierarchicalShapeFunctions(n, psi_values);
    phi_e_values(i, :) = phi_e;
    diff_phi_e_values(i, :) = dphi_e;
end

% Initialize local matrices
K_Local = zeros(n, n);
C_Local = zeros(n, n);

% Compute local matrices using numerical integration
for i = 1:n
    for j = 1:n
        for k = 1:GP
            xi = psi(k);

```



```

        F_xi = a_num * diff_phi_e_values(k, i) * diff_phi_e_values(k, j) * (2/h) - c_num *
        CNM_integrand = -phi_e_values(k, i) * phi_e_values(k, j) * (h/2);

        % Integrate using Gaussian quadrature weights
        K_Local(i, j) = K_Local(i, j) + F_xi * w(k);
        C_Local(i, j) = C_Local(i, j) + CNM_integrand * w(k);
    end
end

% Calculation of global stiffness matrix
global_stiffness_size = s;
K_global = zeros(global_stiffness_size);
C_global = zeros(global_stiffness_size);

% Loop through each element
for element = 1:e
    start_index = 1 + (element-1)*OIF;
    end_index = start_index + n - 1;
    K_global(start_index:end_index, start_index:end_index) = K_global(start_index:end_index, st
    C_global(start_index:end_index, start_index:end_index) = C_global(start_index:end_index, st
end

% Define the nodal points along the domain
X = (x0:h/OIF:xn)';

% Define the source term and calculate its values at nodes
T_x = subs(T, x, X);

exact_values = double(subs(exact_sol_T, x, X)); % Evaluate exact solution at node positions

% Dirichlet Boundary Conditions
C = C_global * T_x;
du_dx = -3;

K_global_Dirichlet = K_global;
g_x_Dirichlet = sym(zeros(s, 1));
GNM_Dirichlet = a_num * g_x_Dirichlet * du_dx;
RHS_Dirichlet = GNM_Dirichlet + C;

u0 = x0;
un = -xn;
K_global_Dirichlet(1, :) = [1, zeros(1, s-1)];
K_global_Dirichlet(end, :) = [zeros(1, s-1), 1];
RHS_Dirichlet(1) = u0;
RHS_Dirichlet(end) = un;

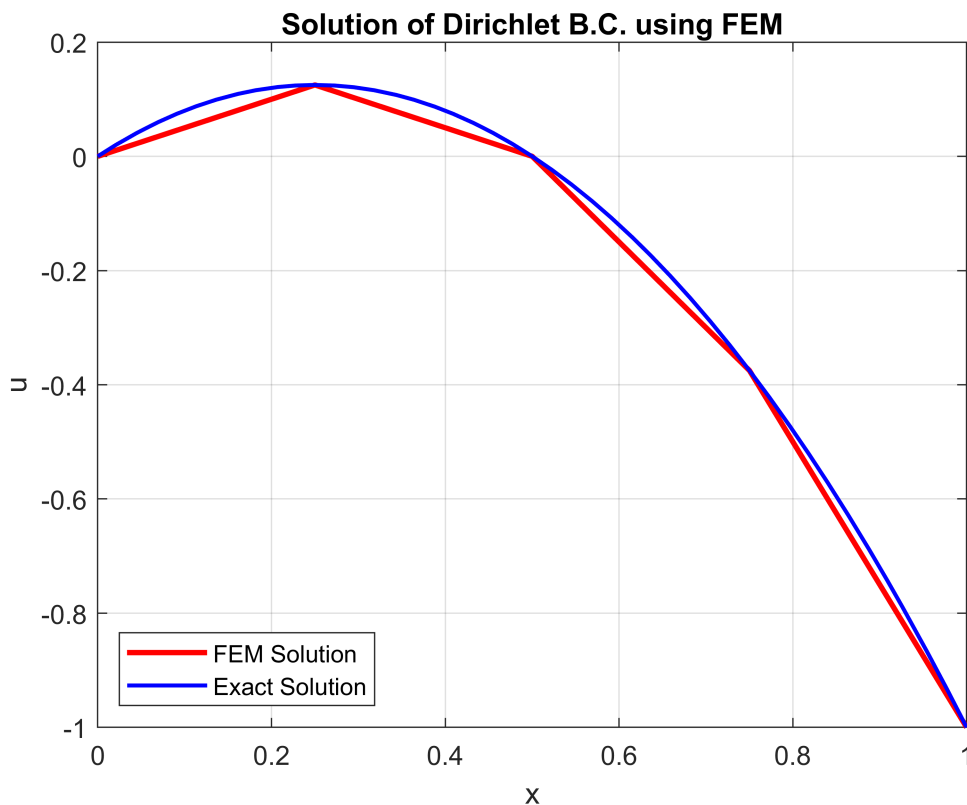
% Solving for displacements using Dirichlet boundary conditions
u_D_bc = inv(K_global_Dirichlet) * RHS_Dirichlet;
displacement_vector = double(u_D_bc)

displacement_vector = 5x1
0

```

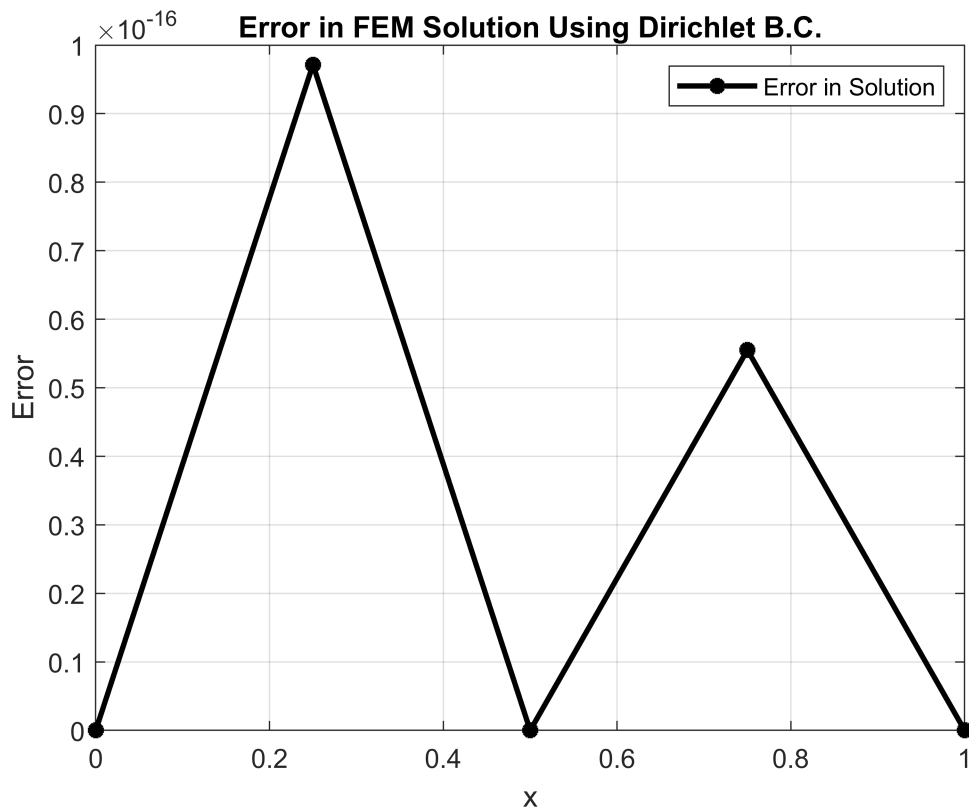
```
0.1250
0
-0.3750
-1.0000
```

```
errors = abs(exact_values - displacement_vector); % Calculate absolute error at each node
% Plot the FEM Solution and the actual solution
figure;
plot(X, displacement_vector, "r", "LineWidth", 2, "DisplayName", "FEM Solution");
hold on;
fplot(exact_sol_T, [0 1], "b", "LineWidth", 1.5, "DisplayName", "Exact Solution");
xlabel("x");
ylabel("u");
title("Solution of Dirichlet B.C. using FEM");
legend('Location', 'southwest');
grid on;
```



```
figure;
plot(X, errors, 'k-*', 'LineWidth', 2, 'DisplayName', 'Error in Solution');
xlabel('x');
ylabel('Error');
title('Error in FEM Solution Using Dirichlet B.C.');
```





```
% Neumann boundary Conditions
```

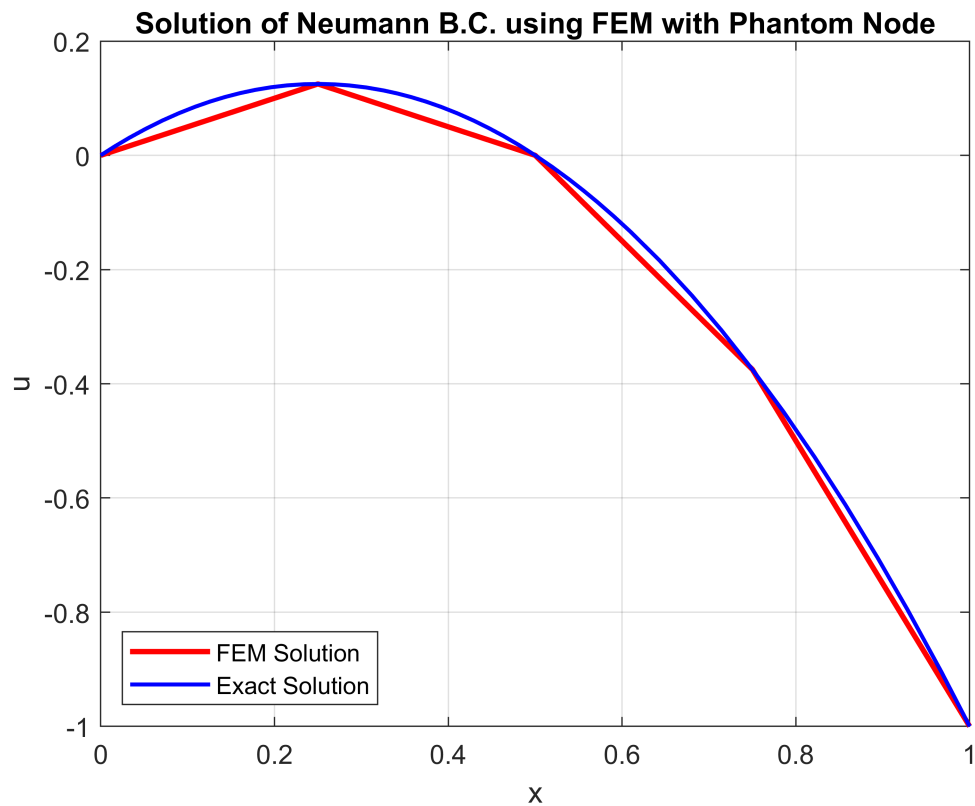
```
K_global_neumann = K_global;  
g_x_neumann = sym(zeros(s, 1));  
g_x_neumann(end) = 1;
```

```
GNM_neumann = a_num * g_x_neumann * du_dx;  
RHS_neumann = GNM_neumann + C;
```

```
K_global_neumann(1, :) = [1, zeros(1, s-1)];  
RHS_neumann(1) = u0;  
u_neumann = inv(K_global_neumann) * RHS_neumann;  
displacement_vector_n = double(u_neumann);
```

```
% Plot the FEM Solution for Neumann boundary conditions
```

```
figure;  
plot(X, displacement_vector_n, "r", "LineWidth", 2, "DisplayName", "FEM Solution");  
hold on;  
fplot(x, exact_sol_T, [0 1], "b", "LineWidth", 1.5, "DisplayName", "Exact Solution");  
xlabel("x");  
ylabel("u");  
title("Solution of Neumann B.C. using FEM with Phantom Node");  
legend('Location', 'southwest');  
grid on;
```



```
function [phi, dphi] = HierarchicalShapeFunctions(n, xi)
    phi = zeros(1, n);
    dphi = zeros(1, n);
    if n == 2 % Linear case
        phi(1) = 0.5 * (1 - xi);
        phi(2) = 0.5 * (1 + xi);
        dphi(1) = -0.5;
        dphi(2) = 0.5;
    elseif n == 3 % Quadratic case
        phi(1) = xi * (xi - 1) / 2;
        phi(2) = -(xi - 1) * (xi + 1);
        phi(3) = xi * (xi + 1) / 2;
        dphi(1) = xi - 0.5;
        dphi(2) = -2 * xi;
        dphi(3) = xi + 0.5;
    end
end

function [w, psi] = Gauss(p_bar)
    psi = zeros(p_bar, 1);
    w = zeros(p_bar, 1);

    if p_bar == 1
        psi(1) = 0; % for p_bar = 1
        w(1) = 2;
```

```

elseif p_bar == 2
    psi(1) = sqrt(3)/3;
    psi(2) = -sqrt(3)/3;
    w(1) = 1;
    w(2) = 1;
elseif p_bar == 3
    psi(1) = sqrt(15)/5;
    psi(2) = -sqrt(15)/5;
    psi(3) = 0;
    w(1) = 5/9;
    w(2) = 5/9;
    w(3) = 8/9;
elseif p_bar == 4
    psi(1) = sqrt((15 - 2*sqrt(30))/35);
    psi(2) = -sqrt((15 - 2*sqrt(30))/35);
    psi(3) = (sqrt((14*sqrt(30))/5 + 21))/7;
    psi(4) = -(sqrt((14*sqrt(30))/5 + 21))/7; % for p_bar = 7
    w(1) = 1/2 + sqrt(30)/36;
    w(2) = 1/2 + sqrt(30)/36;
    w(3) = 1/2 - sqrt(30)/36;
    w(4) = 1/2 - sqrt(30)/36;
elseif p_bar == 5
    psi(1) = 0.906179845938664;
    psi(2) = -0.906179845938664;
    psi(3) = 0.538469310105683;
    psi(4) = -0.538469310105683;
    psi(5) = 0;
    w(1) = 0.236926885056189;
    w(2) = 0.236926885056189;
    w(3) = 0.478628670499366;
    w(4) = 0.478628670499366;
    w(5) = 0.568888888888889;
end
end

```



```

clc;
clear;
syms x;

% Constants and parameters
OIF = 2; % Order of polynomial
E = 1; % Young's modulus
x0 = 0; % Start of the domain
xn = 1; % End of the domain
c_num = -2;

% Source term and exact solution
T = 4*x^2 - 2*x - 4;
exact_sol_T = -2*x^2 + x;
du_dx_exact = diff(exact_sol_T, x); % Derivative of exact solution
strain_exact = du_dx_exact^2 + c_num * exact_sol_T^2; % Strain for exact solution

% Strain energy for the exact solution
strain_energy_exact = double(int(1/2 * E * strain_exact, x, x0, xn));

% Range of elements
element_range = 1:10;
strain_energies_fem = zeros(size(element_range));
strain_energies_exact = repmat(strain_energy_exact, size(element_range));
errors = zeros(size(element_range)); % Store error values

% Initialize the arrays outside the loop
[w, psi] = gaussQuad(OIF + 1); % Assuming maximum order is related to OIF

for e = element_range
    n = OIF + 1;
    h = (xn - x0) / e; % Length of each element
    strain_energy_fem = 0;

    for element = 1:e
        local_X = linspace(x0 + (element-1)*h, x0 + element*h, n);
        local_T_x = double(subs(T, x, local_X));
        local_displacements = double(subs(exact_sol_T, x, local_X));

        for k = 1:length(w)
            [phi_e, dshap] = ShapeFunctions(OIF, psi(k));
            local_strain = dshap .* local_displacements' + c_num .* local_displacements'; % Computed strain
            strain_energy_fem = strain_energy_fem + 1/2 * E * sum(local_strain.^2) * w(k) * h;
        end
    end

    % Store the computed FEM strain energy for this element count
    strain_energies_fem(e) = strain_energy_fem;
    errors(e) = abs((strain_energy_fem - strain_energy_exact) / strain_energy_exact); % Normalized error
end

% Display and plot results
disp('Exact Strain Energy:');

```

Exact Strain Energy:

```
disp(strain_energies_exact);
```

1.0333	1.0333	1.0333	1.0333	1.0333	1.0333	1.0333	1.0333	1.0333	1.0333
--------	--------	--------	--------	--------	--------	--------	--------	--------	--------

```
disp('FEM Strain Energies:');
```

FEM Strain Energies:

```
disp(strain_energies_fem);
```

3.4667	1.8792	1.4979	1.3581	1.2922	1.2561	1.2343	1.2200	1.2103	1.2033
--------	--------	--------	--------	--------	--------	--------	--------	--------	--------

```
disp('Normalized Errors:');
```

Normalized Errors:

```
disp(errors);
```

2.3548	0.8185	0.4496	0.3143	0.2505	0.2156	0.1944	0.1807	0.1712	0.1644
--------	--------	--------	--------	--------	--------	--------	--------	--------	--------

```
% Plotting the results
```

```
figure;
```

```
subplot(2, 1, 1);
```

```
plot(element_range, strain_energies_exact, 'b-o', 'LineWidth', 2, 'DisplayName', 'Exact Strain Energy');
```

```
hold on;
```

```
plot(element_range, strain_energies_fem, 'r-o', 'LineWidth', 2, 'DisplayName', 'FEM Strain Energy');
```

```
xlabel('Number of Elements');
```

```
ylabel('Strain Energy');
```

```
title('Strain Energy Comparison using Lagrange Interpolation');
```

```
legend('Location', 'best');
```

```
grid on;
```

```
subplot(2, 1, 2);
```

```
plot(element_range, errors, 'k-*', 'LineWidth', 2, 'DisplayName', 'Normalized Error');
```

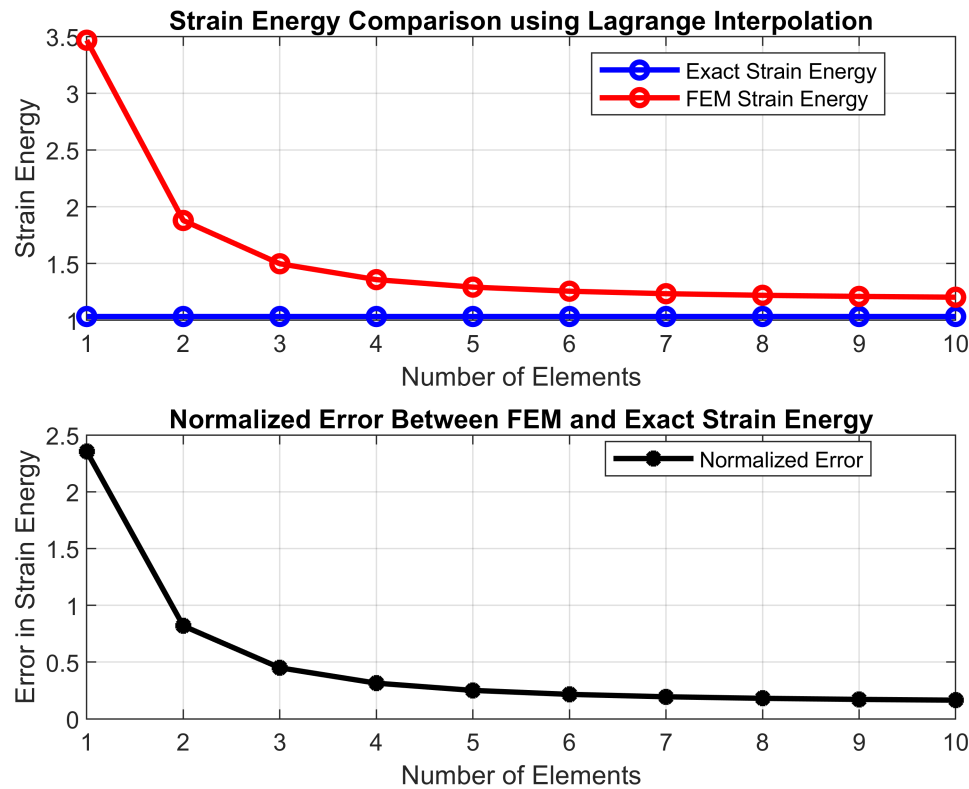
```
xlabel('Number of Elements');
```

```
ylabel('Error in Strain Energy');
```

```
title('Normalized Error Between FEM and Exact Strain Energy');
```

```
legend('Location', 'best');
```

```
grid on;
```



```
function [w, psi] = gaussQuad(p_bar)

    psi = zeros(p_bar, 1);
    w = zeros(p_bar, 1);
    if p_bar == 1
        psi(1) = 0;
        w(1) = 2;
    elseif p_bar == 2
        psi(1) = sqrt(3)/3;
        psi(2) = -sqrt(3)/3;
        w(1) = 1;
        w(2) = 1;
    elseif p_bar == 3
        psi(1) = sqrt(15)/5;
        psi(2) = -sqrt(15)/5;
        psi(3) = 0;
        w(1) = 5/9;
        w(2) = 5/9;
        w(3) = 8/9;
    end
end

function [n, phi_e, dshap] = ShapeFunctions(OIF, x_loc)
```



```

syms x;
n = OIF + 1;
x_local = sym(zeros(1, n));

for i = 1:n
    x_local(i) = -1 + (i-1)*(2/OIF); % Calculate the elements of the vector
end

x_s = sym('x', [1, n]);
N = sym(zeros(size(x_s)));
df = sym(zeros(size(x_s))); % Initialize df

for j = 1:n
    Li = 0;
    den = 1;
    product_term = 1;
    for i = 1:n
        if j ~= i
            term = 1; % Initialize product term for index j
            for k = 1:n
                if k ~= i && k ~= j
                    term = term * (x - x_s(k));
                end
            end
            % Add the product term for index j to Lagrange polynomial for index i
            Li = Li + term;
        end
    end
    %% creating pattern for differentiation
    if i ~= j
        product_term = product_term * (x - x_s(i)) / (x_s(j) - x_s(i));
        den = den * (x_s(j) - x_s(i));
    end
end

N(j) = product_term;
df(j) = Li / den;
end

phi_e = subs(N, x_s, x_local);
dshap = subs(df, x_s, x_local);

% Evaluate the shape functions and their derivatives at the Gauss points
xi_loc = x_loc;
phi_e = double(subs(phi_e, x, xi_loc))';
dshap = double(subs(dshap, x, xi_loc))';
end

```

```

clc;
clear;
syms x;

% Constants and parameters
OIF = 2; % Order of polynomial
E = 1; % Young's modulus
x0 = 0; % Start of the domain
xn = 1; % End of the domain
c_num = -2;

% Source term and exact solution
T = 4*x^2 - 2*x - 4;
exact_sol_T = -2*x^2 + x; % Exact solution for displacement
du_dx_exact = diff(exact_sol_T, x); % Derivative of exact solution
strain_exact = du_dx_exact^2 + c_num * exact_sol_T^2; % Strain for exact solution

% Strain energy for the exact solution
strain_energy_exact = double(int(1/2 * E * strain_exact, x, x0, xn))

```

```

strain_energy_exact = 1.0333

```

```

% Element range specified
element_ranges = [1, 4, 10, 50, 100];
strain_energies_fem = zeros(size(element_ranges));
errors = zeros(size(element_ranges));

for idx = 1:length(element_ranges)
    e = element_ranges(idx);
    n = OIF + 1;
    h = (xn - x0) / e; % Length of each element

    % Gauss quadrature process
    [w, psi] = gaussQuad(n); % Using number of nodes to determine Gauss points

    % Initialize variables for integration
    strain_energy_fem = 0;

    for element = 1:e
        local_X = linspace(x0 + (element-1)*h, x0 + element*h, n);
        local_T_x = double(subs(T, x, local_X));
        local_displacements = double(subs(exact_sol_T, x, local_X));

        % Local strain calculations
        for k = 1:length(w)
            [phi_e, dphi_e] = shapeFunctions(OIF, psi(k));
            local_strain = dphi_e * local_displacements' + c_num * local_displacements'; % Cor
            strain_energy_fem = strain_energy_fem + 1/2 * E * sum(local_strain.^2) * w(k) * h;
        end
    end

    % Store results
    strain_energies_fem(idx) = strain_energy_fem;

```

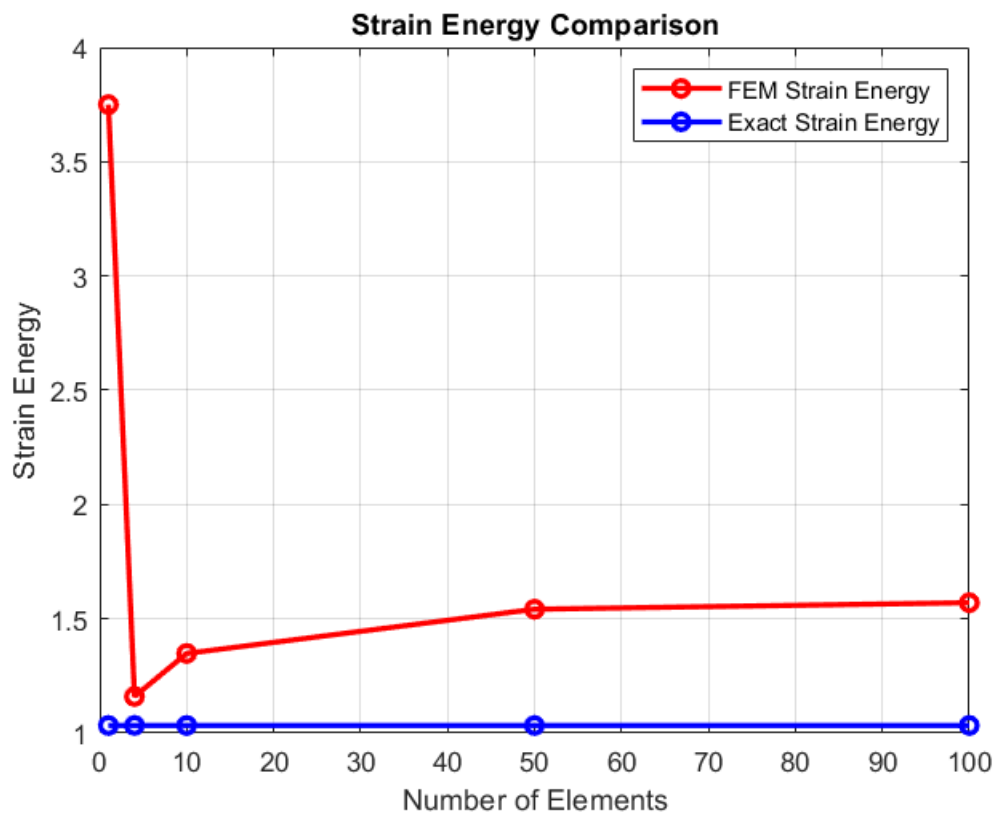
```

errors(idx) = abs(strain_energy_fem - strain_energy_exact);
end

% Plotting results
figure;
plot(element_ranges, strain_energies_fem, 'r-o', 'LineWidth', 2, 'DisplayName', 'FEM Strain Energy');
hold on;
plot(element_ranges, repmat(strain_energy_exact, 1, length(element_ranges)), 'b-o', 'LineWidth', 2, 'DisplayName', 'Exact Strain Energy');

xlabel('Number of Elements');
ylabel('Strain Energy');
title('Strain Energy Comparison');
legend('show');
grid on;

```

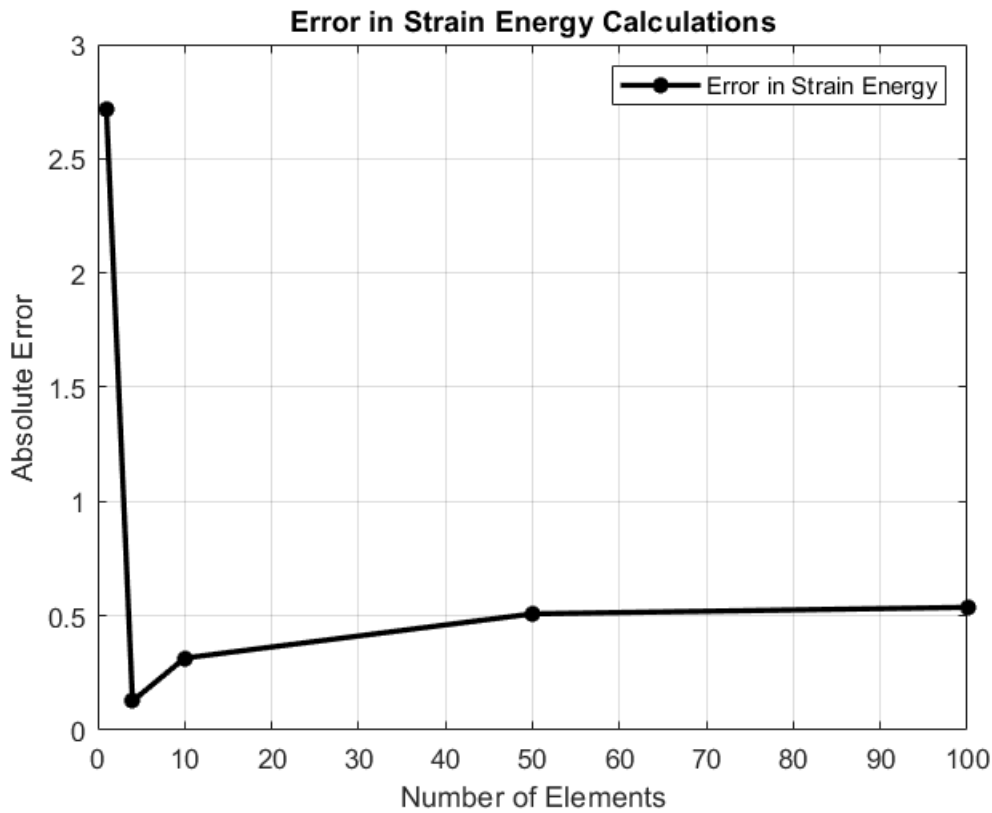


```

% Additional plotting for error comparison
figure;
plot(element_ranges, errors, 'k-*', 'LineWidth', 2, 'DisplayName', 'Error in Strain Energy');
xlabel('Number of Elements');
ylabel('Absolute Error');
title('Error in Strain Energy Calculations');
legend('show');
grid on;

```





```
function [w, psi] = gaussQuad(numPoints)
    % Returns the weights and positions for Gauss quadrature
    w = zeros(numPoints, 1);
    psi = zeros(numPoints, 1);
    if numPoints == 1
        psi(1) = 0;
        w(1) = 2;
    elseif numPoints == 2
        psi = [-sqrt(1/3), sqrt(1/3)];
        w = [1, 1];
    elseif numPoints == 3
        psi = [-sqrt(3/5), 0, sqrt(3/5)];
        w = [5/9, 8/9, 5/9];
    end
end

function [phi_e, dphi_e] = shapeFunctions(OIF, psi_val)
    % Shape functions and derivatives for Quadratic elements
    if OIF == 2
        phi_e = [0.5*psi_val*(psi_val-1), 1-psi_val^2, 0.5*psi_val*(psi_val+1)];
        dphi_e = [psi_val - 0.5, -2*psi_val, psi_val + 0.5];
    else

```

```
        error('Shape functions for given order not implemented.');
```

```
    end
```

```
end
```

```

clc;
clear;

n = 4; % Number of elements
E = 210e9; % Modulus of Elasticity in Pa
I = 2e-4; % Moment of Inertia
beam_length = 12; % Length of the beam
H = 0.5; % Height of the beam in meters, added for stress calculations

for element_count = [1, 4, 10, 50, 100] % Different element range
    n = element_count;
    el = beam_length / n; % Length of each element

    % Local stiffness matrix for one element
    ele_K = E * I / el^3 * [12    6*el  -12    6*el;
                           6*el  4*el^2 -6*el  2*el^2;
                           -12   -6*el   12   -6*el;
                           6*el  2*el^2 -6*el  4*el^2];

    % Global stiffness matrix
    K_global = zeros(2*(n+1), 2*(n+1));

    % Assemble the global stiffness matrix
    for element = 1:n
        start_index = 2*element - 1;
        end_index = start_index + 3;
        K_global(start_index:end_index, start_index:end_index) = ...
            K_global(start_index:end_index, start_index:end_index) + ele_K;
    end

    % Load vectors initialization
    F = zeros(2*(n+1), 1); % Initialize distributed load vector
    Q = zeros(2*(n+1), 1); % Initialize point load and moment vector

    % Applying distributed loads
    for element = 1:n
        start_index = 2 * element - 1;
        end_index = start_index + 3;
        F_local = 1000 * el * [1/2; el/12; 1/2; -el/12]; % 1000 N/m as an example distributed load
        F(start_index:start_index+3) = F(start_index:start_index+3) + F_local;
    end

    % Initialize parameters for loads
    %P_center = 5000; % 5000 N point load at the center
    P_end = 1500; % 1500 N point load at the end
    M_end = 2000; % 2000 Nm moment at the end

    % Apply point load at the center of the beam
    center_node_index = n / 2 + 1; % Correct index for an even number of elements
    %Q(2 * center_node_index - 1) = P_center; % Apply to the vertical displacement DOF

    % Apply point load and moment at the end of the beam

```



```

Q(end-1) = P_end; % Point load at the end
Q(end) = M_end; % Moment at the end

% Total load vector
Total_Load = F + Q;

% Apply boundary conditions
% For a fixed support at node 1 (no displacement or rotation)
K_global(1,:) = 0;
K_global(:,1) = 0;
K_global(1,1) = 1;
Total_Load(1) = 0;

K_global(2,:) = 0;
K_global(:,2) = 0;
K_global(2,2) = 1;
Total_Load(2) = 0;

% Solve for displacements
displacements = K_global \ Total_Load;

displacements

% Post-processing for shear forces and bending moments
M = zeros(n+1, 1);
V = zeros(n+1, 1);
for i = 1:n
    % Two nodes per element
    start_index = i;
    end_index = i + 1;

    % Displacement derivatives (finite differences)
    dwdx = (displacements(2*end_index) - displacements(2*start_index-1)) / 2*el;
    d2wdx2 = (displacements(2*end_index) - 2*displacements(2*start_index) + displacements(2*start_index-1)) / el^2;

    % Bending moment and shear force
    M(start_index) = -E * I * d2wdx2; % bending moment
    V(start_index) = -E * I * (dwdx / el); % shearforces
end

% Bending stress calculation
sigma = M * (H / 2) / I;

% Output results
x = linspace(0, beam_length, n+1);
figure;
subplot(3,1,1);
plot(x, displacements(1:2:end), 'b.-'); % Plotting vertical displacements
title('Transverse Displacement');

subplot(3,1,2);
plot(x, M, 'r.-');
title('Bending Moment');

```

```

subplot(3,1,3);
plot(x, V, 'g.-');
title('Shear Force');

```

```

fprintf('Analysis for %d elements completed.\n', n);

```

```

end

```

```

displacements = 4×1

```

```

0

```

```

0

```

```

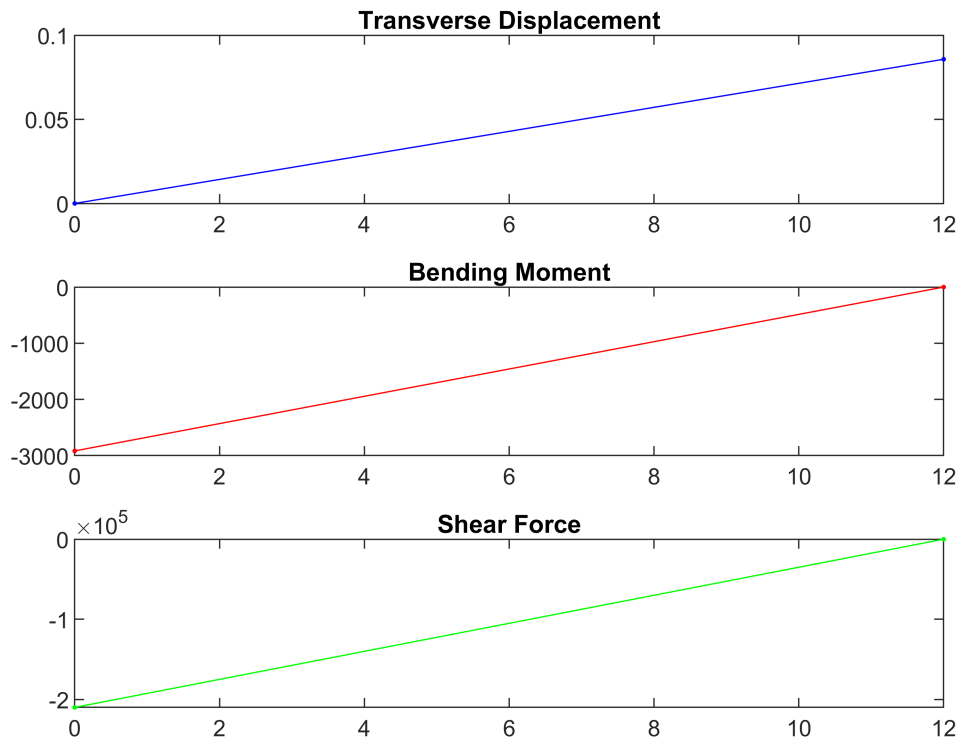
0.0857

```

```

0.0100

```



```

Analysis for 1 elements completed.

```

```

displacements = 10×1

```

```

0

```

```

0

```

```

0.0085

```

```

0.0052

```

```

0.0291

```

```

0.0082

```

```

0.0562

```

```

0.0096

```

```

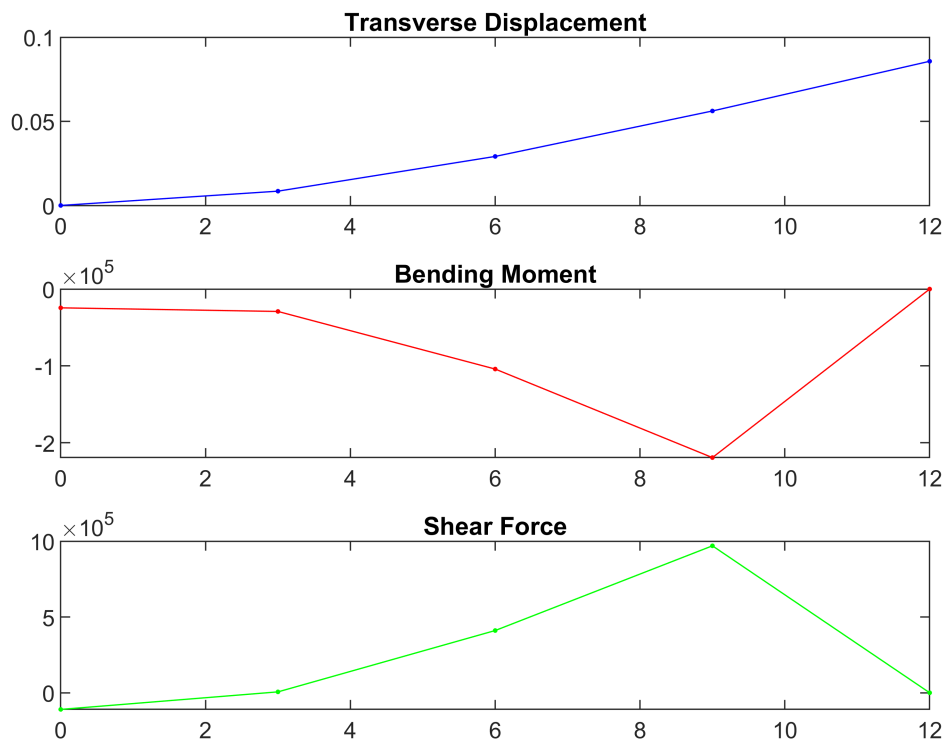
0.0857

```

```

0.0100

```



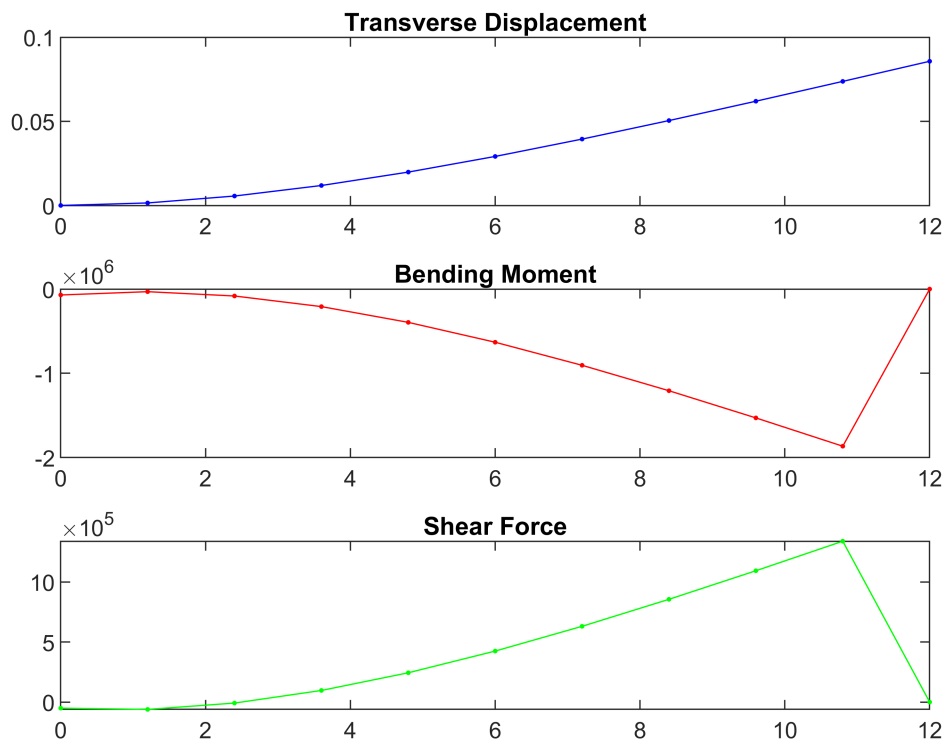
Analysis for 4 elements completed.

displacements = 22x1

```

0
0
0.0015
0.0024
0.0056
0.0044
0.0119
0.0060
0.0198
0.0073
:
:

```



Analysis for 10 elements completed.

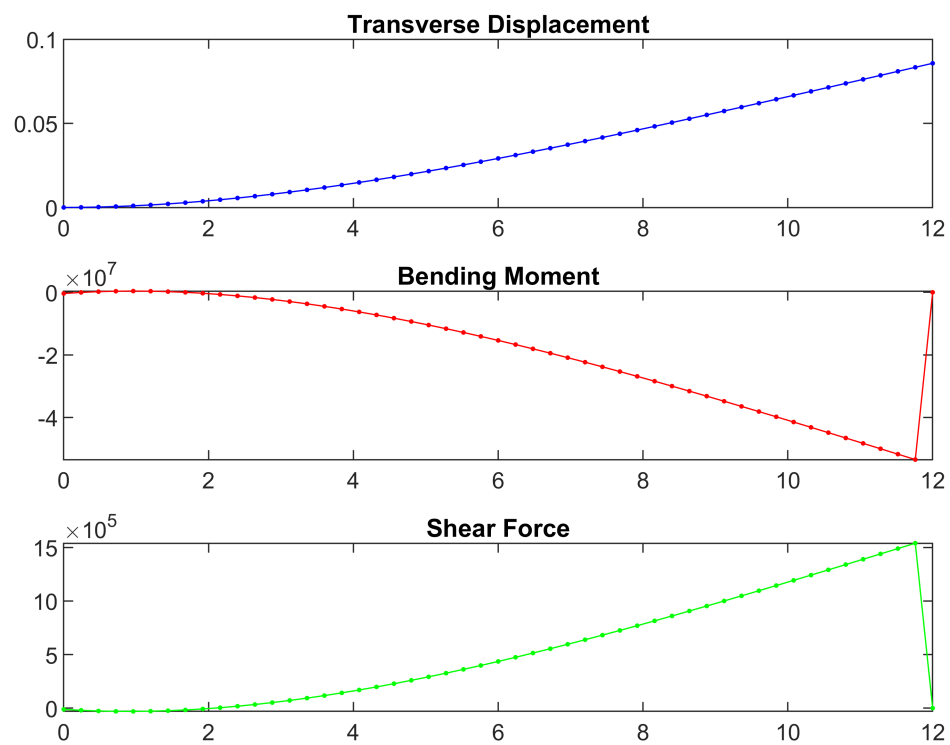
displacements = 102x1

```

0
0
0.0001
0.0005
0.0002
0.0010
0.0005
0.0015
0.0010
0.0020
:
:

```





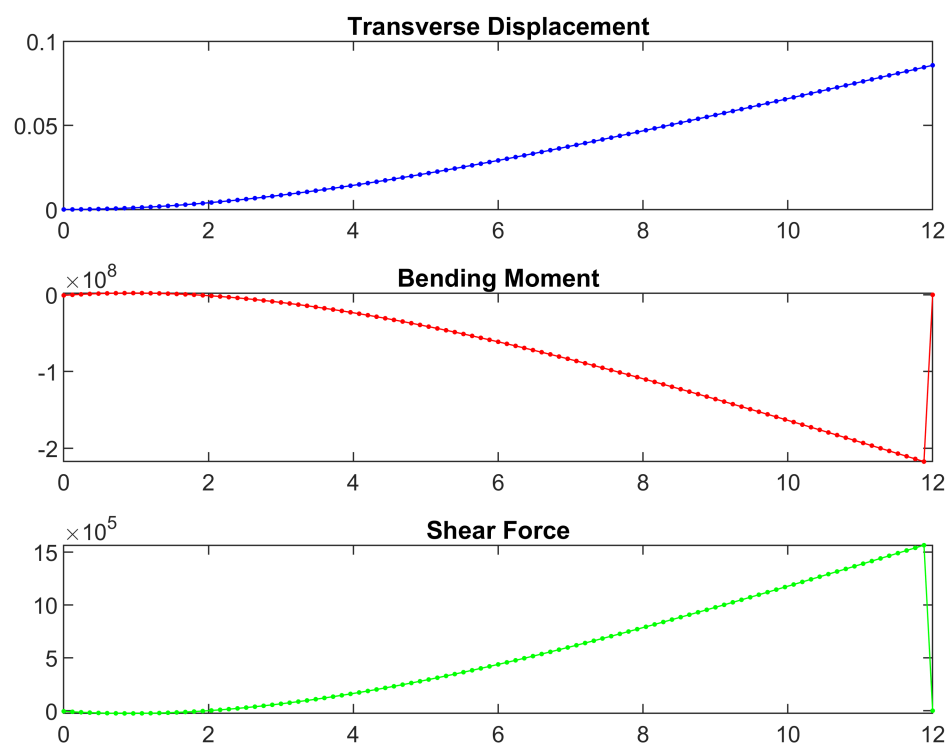
Analysis for 50 elements completed.

displacements = 202x1

```

0
0
0.0000
0.0003
0.0001
0.0005
0.0001
0.0008
0.0002
0.0010
:
:

```



Analysis for 100 elements completed.