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MATLAB FEM Code - From Elasticity to Plasticity

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<p>Abstract:</p> <p>A MATLAB Finite Element code for plane strain analysis of footings on an Elasto-plastic material using the Mohr Coulomb failure criteria has been developed. The first step is to develop codes for mesh generation and Gaussian numerical integration. Next, the force matrix, the stiffness matrix and the self weight matrix are assembled. After that functions for non linear analysis such as the plastic potential derivatives are formed. Finally plots of the mesh, displacement shadings, stress shadings and stress-strain curves are developed.</p> <p>For the purpose of verification results from the code for biaxial test are compared with the theoretical solution. Additionally comparison is made between the code and prandtl's bearing capacity solutions for a footing problem. These results show that accuracy depends on two factors: - the type of the element and the number of elements used. The three node triangular element and the four node rectangular element give less accurate results when compared to higher order element types. And for a relatively accurate result the number of elements should be too high.</p>
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Keywords:

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2. FEM
3. MATLAB
4. Elasto-Plastic

MASTER THESIS

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MATLAB FEM code – from elasticity to plasticity

BACKGROUND

Supported excavations and other comparably complex geotechnical problems were first studied with the finite element method (FEM) in the early 1970s. Since then, the method has been considerably refined and developed into a versatile design tool. The conditioning parameters of FEM analyses are well understood through many case studies presented in literature, often including a comparison of measured and calculated performance. However, with increasing versatility and development, the FEM has also become a tool that is increasingly difficult to understand in all its facets for the practicing engineer as well as for students new to the subject. A well structured, easy to use FEM code with limited features is therefore a desirable starting point in teaching the method. Such a code shall be developed within this thesis.

TASK DESCRIPTION

The aim of the thesis is to generate a structured FEM code in MATLAB that can be applied to basic geotechnical problems. The code shall be structured so that it is easy to understand and that it can be easily expanded by other students. A good documentation of the code is essential.

The FEM code generated within this project shall be limited to plain strain. An elasto-plastic material model with Mohr Coulomb failure criterion shall be implemented. Results of the development shall be validated against a commercially available FE code and/or analytical solutions. It shall be also focused on the performance of various element types and integration techniques.

The objectives of the thesis are defined as follows:

1. To summarize the theory of non-linear FEM analysis
2. To realize and document a plane strain non linear FEM code
3. To implement an elasto-plastic model with Mohr Coulomb failure criterion.
4. To validate the implementation against analytical solutions and other FEM codes
5. To discuss important aspects of the implementation, such as explicit and implicit time integration, in more depth
6. To investigate into the performance of various element types
7. To conclude on the generated results

It is acknowledged that the given task, especially the implicit version of the code is a highly advanced task and that due to unforeseen programming issues not all objectives may be fulfilled in the given timeframe. The student may base his work on a 1-point version of the FEM code Plaxis which the student got access to.

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Date: 06.06.2012



Professor in charge

Dedicated to Dr. Miftah Nesru Sherif

Preface

The purpose of this paper is to compile a FEM code in MATLAB that can be applied to basic geotechnical problems. It is structured in an easy way so it can be understandable and can be used as a platform for future work for other geotechnical problems. The FEM code generated within this paper is limited to plain strain problems with an elasto-plastic material model using the Mohr Coulomb failure criterion. The final outcome of this work is two codes that can be used for biaxial test and bearing capacity problem. In a previous project work a similar code for linear elastic materials has been done.

Acknowledgements

I would like to sincerely acknowledge Professor Thomas Benz, Department of Civil and Transport Engineering, NTNU for his full cooperation during the progress of the work and for providing a one point FORTRAN code that made this work to become a reality. I would also like to thank all my friends and family for their advises and moral supports.

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CHAPTER 1

INTRODUCTION

1.1 Background

The finite element method is nowadays the most frequently used computational method in engineering problems. In this numerical technique all complexities of a problem such as shape, boundary and loading conditions are kept the same but the results obtained are approximate. When using this method, calculations are robust due to the large number of unknowns leading to a large pile of simultaneous equations for the user to solve. Hence the use of computer programs to take care of these equations is one face of the method.

The evaluation of these simultaneous equations is done by using one of the matrix methods of solving simultaneous equations. The MATLAB programming language is useful in illustrating how to program the Finite element method due to the fact that it allows one to very quickly code Numerical methods and has a vast predefined mathematical library suitable for handling matrices.

1.2 Scope and limitation of the problem

The MATLAB Finite Element code presented here analyzes the stresses, strains and displacements and gives the bearing capacity of a uniformly loaded strip footing on an Elasto-plastic soil material in plane strain condition. Additionally a code for biaxial test is also included. Bearing capacity analysis begins with the selection of the dimensions of the region which is usually a rectangular region that is restrained vertically at the left and right sides and totally restrained at the bottom end. When dividing the region in to smaller pieces five element types are provided, three quadrilaterals (with 9,8 or 4 nodes) and two triangular (with 3 or 6 nodes).

There are two Load types to be considered; the first one is an external uniform load applied at the top of the region, which the user needs to provide its length of distribution and magnitude. The second type is the self weight of the soil which is generated by inserting the unit weight of the soil material. When analysis begins all the functions developed work in the background to provide the outputs for the user to see.

1.3 Organization of the paper

Any finite element analysis has the following main steps

- Define a set of elements connected at nodes (discretization of the domain).
- Assemble the global system, $[K]\{u\} = \{f\}$
- Modify the global system by imposing essential (displacements) boundary conditions.
- Solve the global system and obtain the global displacements, $\{u\}$
- For each element, evaluate the strains and stresses at the nodes.

Chapter 2 covers about the types of elements and how the discretization is done and which boundary conditions to be applied. In chapter 3 the basic equations of finite element analysis in plane strain elements are presented in matrix form. Chapter 4 discusses about plane strain and stress invariants. In chapter 5 the main features of an Elasto-plastic Mohr Coulomb material is presented. The explicit integration scheme in load stepping for an Elasto-plastic material is shown in chapter 6. In chapter 7 the main features of the code and how the user can implement it is presented. Finally in chapter 8 the code is tested for validity with a biaxial test and a bearing capacity problem and the results are compared with results from the PLAXIS software and with theoretically solutions. Additionally the effect of element number and element type on these results is discussed. Finally summary and future work based on this code are outlined. For quick reference, all the functions and script files in the code are provided as an appendix. Additionally a code for linear elastic materials from a previous work is attached as a reference in the appendix.

CHAPTER 2

DISCRETIZATION OF THE DOMAIN

2.1 Types of Elements

The first step in the finite element analysis involves the division of the body into smaller Pieces, known as finite elements. This is equivalent to replacing the body with an infinite number of degrees of freedom by a system having finite number of degrees of freedom.

The shapes, number, and configurations of the elements are chosen in a way that the resulting body resembles as closely as possible to that of the original body. The choice of the type of element is dictated by the geometry of the body and the number of independent coordinates necessary to describe the system. In plain strain analysis the geometry, material properties, and the field variable of the system can be described in terms of two spatial coordinates i.e. x and y . Elements are considered to be interconnected at specific points called nodes.

Nodes are the selected finite points at which basic unknowns are to be determined in the finite element analysis .There are two types of nodes, external nodes and internal nodes. External nodes are those which occur on the edges or surface of the elements and they may be common to two or more elements. These nodes may be further classified as Primary nodes and Secondary nodes. Primary nodes occur at the corners of elements. Secondary nodes occur along the side of an element but not at corners. Internal nodes are the one which occur inside an element and they are unique to each element. Based on this elements are categorized in to two groups.

- 1) Basic elements
 - Three node triangular element (T3)
 - Four node rectangular element (Q4)

- 2) Higher order elements
 - Six node triangular element (T6)
 - Eight node rectangular element (Q8)
 - Nine node rectangular element (Q9)

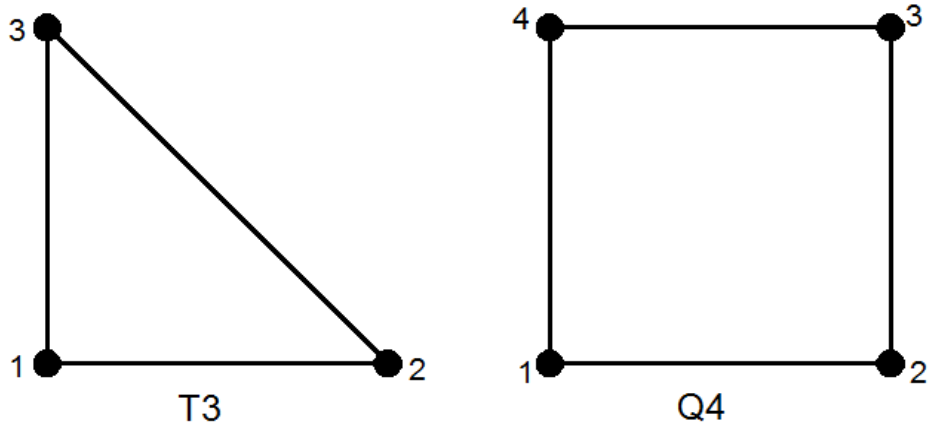


Figure 2.1:- Three node triangular element and four node rectangular element

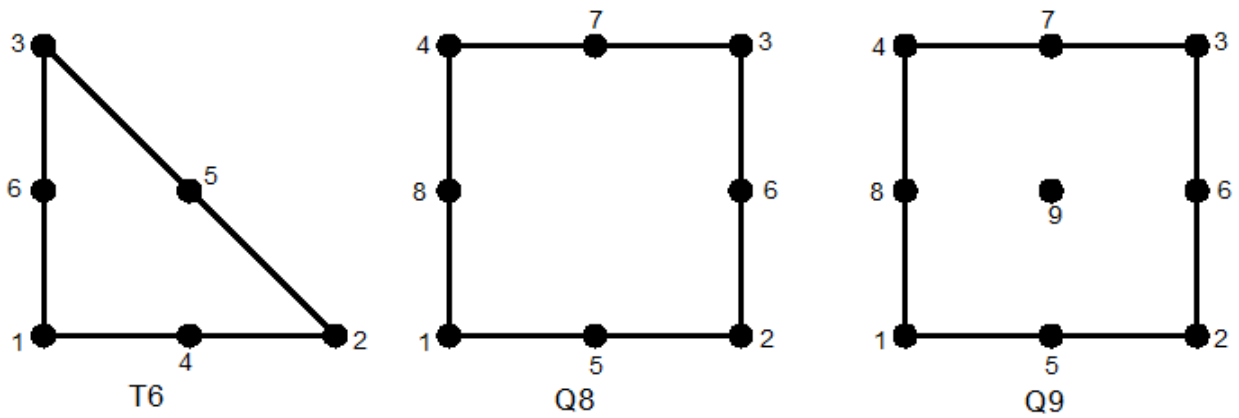


Figure 2.2:- six node triangular element, eight and nine node rectangular element

2.2 Mesh generation

Mesh generation is the process of determining the nodal and element connectivity in the domain. Nodes are numbered with an increasing order in the x-direction starting from the bottom left corner of the domain, which is a rectangle with length L and depth D whose four corner points are given by:-

$$\text{Point 1} = (0, -D/2)$$

$$\text{Point 2} = (L, -D/2)$$

$$\text{Point 3} = (L, D/2)$$

$$\text{Point 4} = (0, D/2)$$

Nodal connectivity is given by a matrix containing the coordinate of each node relative to the origin. Whereas the element connectivity defines how the elements are connected to each other. It is given by a matrix with the node identification in each element.

For mesh generation the number of elements in the region is user specified by the number of elements in x and y direction. For rectangular elements it is generated by dividing the original rectangle in to smaller equal rectangles whose size depend on the dimensions L and D of the domain. In case of triangular elements, for the same number of elements in x and y direction, these rectangles are further divided in to two triangles giving a denser mesh with number of 'elements' twice that of the rectangular ones.

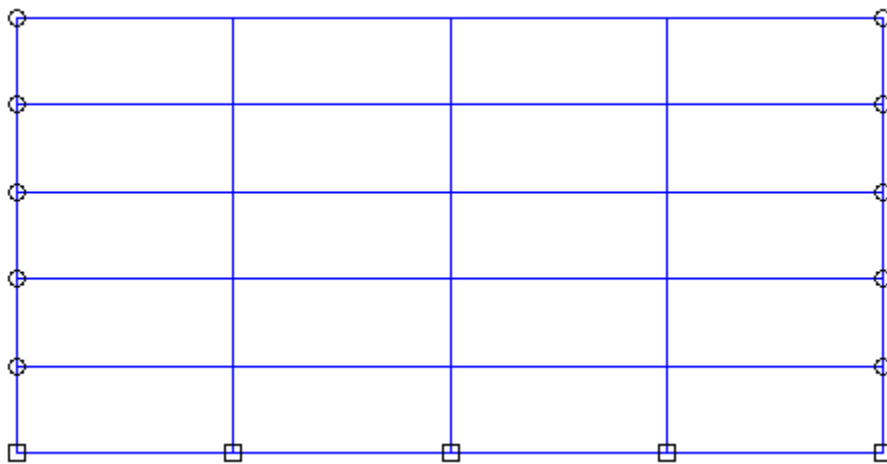


Figure 2.3:- Typical region with four node rectangular elements with support condition

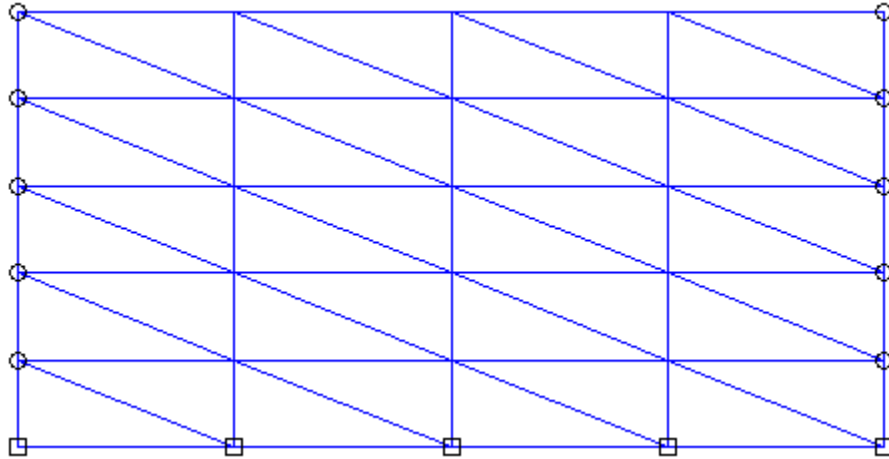


Figure 2.4:- Typical region with three node triangular elements with support condition

2.3 Support condition

The restrained degrees of freedom in the system are two types

- Translation in x and y direction for the nodes at the bottom of the domain which are represented by black squares in the mesh and
- Translation in x direction for the nodes at the left and right edge of the domain, represented by black circles in the mesh.

These prescribed degrees of freedom, being zero, usually reduce the size of the stiffness matrix and hence the calculation effort. Representation of support conditions in the finite element mesh is shown in figures 2.3 and 2.4 above.

However, for the case of biaxial test the restrained degrees of freedom are,

- Translation in x and y direction for the node at the bottom left of the domain which is represented by black squares in the mesh.
- Translation in x direction for the nodes at the left edge of the domain, represented by black circles in the mesh and
- Translation in y direction for the nodes at the bottom of the domain, represented by black circles in the mesh.

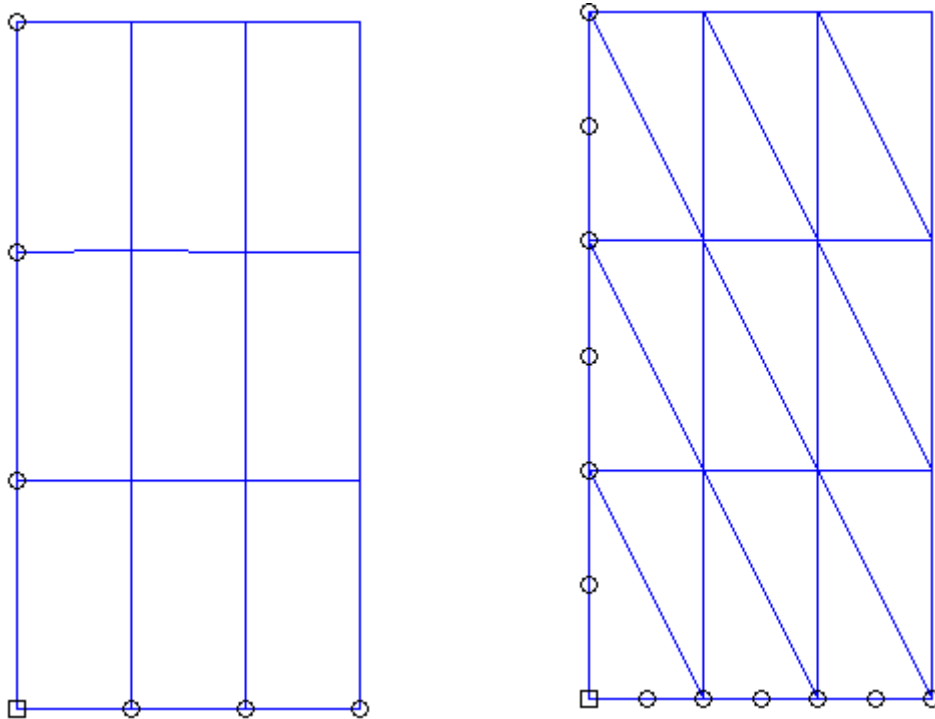


Figure 2.5:- Typical biaxial test finite mesh for Q4 and T6 element with support condition

CHAPTER 3

BASIC EQUATIONS IN FEM

3.1 Shape functions

In the finite element analysis the aim is to find the field variables at nodal points by rigorous analysis, assuming that at any point inside the element basic variable is a function of values at nodal points of the element. This function which relates the field variable at any point within the element to the field variables of nodal points is called shape function or interpolation function. Taking displacement as the field variable this relationship can be expressed as:-

$$u = \sum_{i=1}^n N_i u_i \dots\dots\dots 3.1$$

$$v = \sum_{i=1}^n N_i v_i \dots\dots\dots 3.2$$

Where

u= horizontal displacement

v= vertical displacement

u_i = horizontal displacement at node i

v_i = vertical displacement at node i

N_i = shape function expression at node i

Summation is over the number of nodes (n) of the element.

The shape functions are always expressed in terms of the natural coordinate system. A natural coordinate system is a coordinate system which permits the specification of a point within the element by a set of dimensionless numbers, whose magnitude never exceeds unity. It is obtained by assigning weights to the nodal coordinates in defining the coordinate of any point inside the element. Hence such system has the property that ith coordinate has unit value at node i of the element and zero value at all other nodes. As an illustration the use of shape functions is shown for three element categories.

For a two node line element

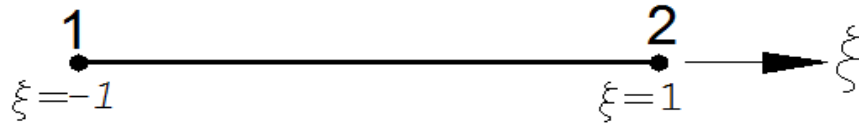


Figure 3.1:- Natural coordinates in two node line element

From equation 3.1,

$$\mathbf{u} = \sum_{i=1}^n N_i u_i = N_1 u_1 + N_2 u_2$$

Where,

$$N_1 = \frac{(1-\xi)}{2}, \quad N_2 = \frac{(1+\xi)}{2}$$

$$\mathbf{u} = [N_1 \quad N_2] \begin{Bmatrix} u_1 \\ u_2 \end{Bmatrix} \dots\dots\dots 3.3$$

For a three node triangular element

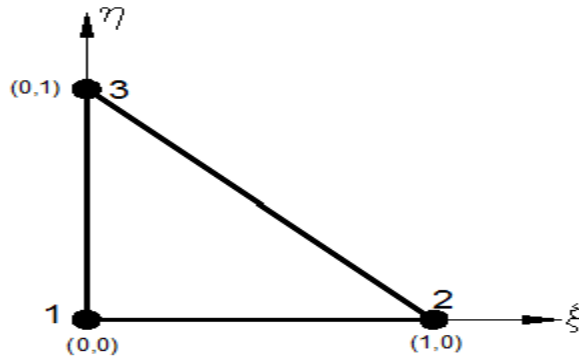


Figure 3.2:- Natural coordinates in three node triangular element

From equations 3.1 and 3.2,

$$\mathbf{u} = \sum_{i=1}^n N_i u_i = N_1 u_1 + N_2 u_2 + N_3 u_3$$

$$\mathbf{v} = \sum_{i=1}^n N_i v_i = N_1 v_1 + N_2 v_2 + N_3 v_3$$

$$N_1 = 1 - \xi - \eta \quad , \quad N_2 = \xi \quad \text{and} \quad N_3 = \eta$$

$$\begin{Bmatrix} \mathbf{u} \\ \mathbf{v} \end{Bmatrix} = \begin{pmatrix} N_1 & 0 & N_2 & 0 & N_3 & 0 \\ 0 & N_1 & 0 & N_2 & 0 & N_3 \end{pmatrix} \begin{Bmatrix} u_1 \\ v_1 \\ u_2 \\ v_2 \\ u_3 \\ v_3 \end{Bmatrix} \dots\dots\dots 3.4$$

For a four node rectangular element

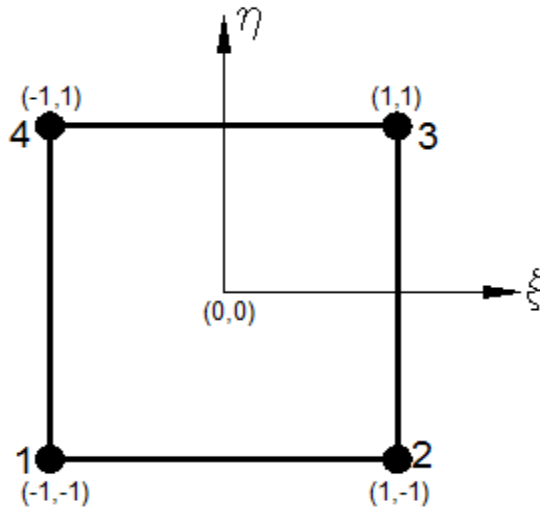


Figure 3.3:- Natural coordinates in four node rectangular element

From equations 3.1 and 3.2,

$$u = \sum_{i=1}^n N_i u_i = N_1 u_1 + N_2 u_2 + N_3 u_3 + N_4 u_4$$

$$v = \sum_{i=1}^n N_i v_i = N_1 v_1 + N_2 v_2 + N_3 v_3 + N_4 v_4$$

Where,

$$N_1 = \frac{(1-\xi)(1-\eta)}{4}, \quad N_2 = \frac{(1+\xi)(1-\eta)}{4}$$

$$N_3 = \frac{(1+\xi)(1+\eta)}{4}, \quad N_4 = \frac{(1-\xi)(1+\eta)}{4}$$

$$\begin{Bmatrix} \mathbf{u} \\ \mathbf{v} \end{Bmatrix} = \begin{pmatrix} N_1 & 0 & N_2 & 0 & N_3 & 0 & N_4 & 0 \\ 0 & N_1 & 0 & N_2 & 0 & N_3 & 0 & N_4 \end{pmatrix} \begin{Bmatrix} u_1 \\ v_1 \\ \cdot \\ \cdot \\ u_4 \\ v_4 \end{Bmatrix} \dots\dots\dots 3.5$$

Using similar procedure, the shape functions for higher order elements can be formulated as presented in the function *shape_func* in appendix C.

3.2 The Strain displacement matrix

Relationship between strains at any point in the element with nodal displacement can be formed using strain displacement matrices.

$$\{\varepsilon\} = \begin{Bmatrix} \varepsilon_{xx} \\ \varepsilon_{yy} \\ \varepsilon_{xy} \\ \varepsilon_{zz} \end{Bmatrix} = [B]\{u\} \dots\dots\dots 3.6$$

Where

- {ε} = strain at any point in the element.
- {u} = displacement vector of nodal values for the element
- [B] = strain displacement matrix

The strain displacement matrix is a function of the partial derivatives of shape functions with respect to the Cartesian coordinates x and y . Because shape functions are not directly functions of x and y but of the natural coordinates ξ and η , the chain rule can be used for differentiation. For example considering a four node quadrilateral element:-

$$B = \begin{pmatrix} \frac{\partial N_1}{\partial x} & 0 & \frac{\partial N_2}{\partial x} & 0 & \frac{\partial N_3}{\partial x} & 0 & \frac{\partial N_4}{\partial x} & 0 \\ 0 & \frac{\partial N_1}{\partial y} & 0 & \frac{\partial N_2}{\partial y} & 0 & \frac{\partial N_3}{\partial y} & 0 & \frac{\partial N_4}{\partial y} \\ \frac{\partial N_1}{\partial y} & \frac{\partial N_1}{\partial x} & \frac{\partial N_2}{\partial y} & \frac{\partial N_2}{\partial x} & \frac{\partial N_3}{\partial y} & \frac{\partial N_3}{\partial x} & \frac{\partial N_4}{\partial y} & \frac{\partial N_4}{\partial x} \end{pmatrix} \dots\dots\dots 3.7$$

But, $\frac{\partial N}{\partial x} = \frac{\partial N}{\partial \xi} \frac{\partial \xi}{\partial x} + \frac{\partial N}{\partial \eta} \frac{\partial \eta}{\partial x} \dots\dots\dots 3.8$

and,

$$\frac{\partial N}{\partial y} = \frac{\partial N}{\partial \xi} \frac{\partial \xi}{\partial y} + \frac{\partial N}{\partial \eta} \frac{\partial \eta}{\partial y} \dots\dots\dots 3.9$$

Which can be expressed at any node in matrix form as:-

$$\begin{bmatrix} \frac{\partial N_i}{\partial x} \\ \frac{\partial N_i}{\partial y} \end{bmatrix} = \begin{pmatrix} \frac{\partial \xi}{\partial x} & \frac{\partial \eta}{\partial x} \\ \frac{\partial \xi}{\partial y} & \frac{\partial \eta}{\partial y} \end{pmatrix} \begin{bmatrix} \frac{\partial N_i}{\partial \xi} \\ \frac{\partial N_i}{\partial \eta} \end{bmatrix} \dots\dots\dots 3.10$$

Introducing the Jacobian matrix which relates derivatives of the function in local natural coordinate system to derivatives in global coordinate system,

$$[J] = \begin{pmatrix} \frac{\partial x}{\partial \xi} & \frac{\partial y}{\partial \xi} \\ \frac{\partial x}{\partial \eta} & \frac{\partial y}{\partial \eta} \end{pmatrix} = \begin{pmatrix} \frac{\partial N_1}{\partial \xi} & \frac{\partial N_2}{\partial \xi} & \frac{\partial N_3}{\partial \xi} & \frac{\partial N_4}{\partial \xi} \\ \frac{\partial N_1}{\partial \eta} & \frac{\partial N_2}{\partial \eta} & \frac{\partial N_3}{\partial \eta} & \frac{\partial N_4}{\partial \eta} \end{pmatrix} \begin{pmatrix} x_1 & y_1 \\ x_2 & y_2 \\ x_3 & y_3 \\ x_4 & y_4 \end{pmatrix} \dots\dots\dots 3.11$$

Therefore

$$\begin{bmatrix} \frac{\partial N_i}{\partial x} \\ \frac{\partial N_i}{\partial y} \end{bmatrix} = \mathbf{J}^{-1} \begin{bmatrix} \frac{\partial N_i}{\partial \xi} \\ \frac{\partial N_i}{\partial \eta} \end{bmatrix} \dots\dots\dots 3.12$$

Hence the strain displacement matrix can be calculated for any type of element using these expressions.

3.3 The Stiffness Matrix

Displacements in an element are a result of externally applied loads or self weight. Relationship between these parameters can be formed using what is called stiffness. Consider a small portion of linearly elastic material subjected to externally applied nodal force, {dF}, resulting displacements {du}, strains {dε} and stresses {dσ} at the nodes. The relationship between these variables can be plotted graphically as shown below.

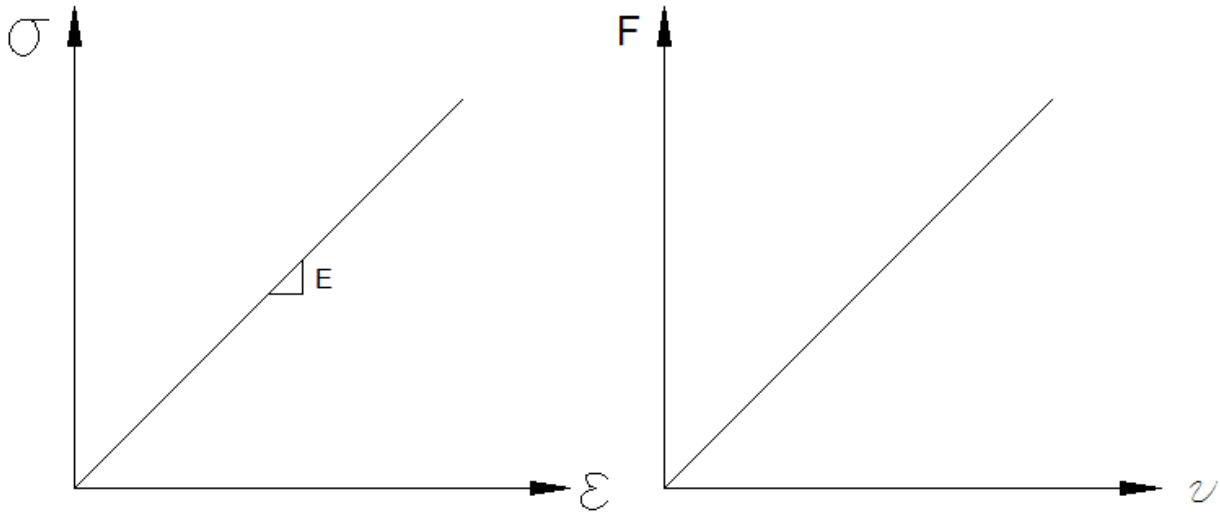


Figure 3.4:- stress-strain and load-displacement relations in linearly-elastic material

From the principle of minimum potential energy which states that “Work done by external applied loads is equal to internal strain energy” the following equations can be written.

The external work, W_e , is equal to the area under the force displacement graph,

$$W_e = \frac{1}{2} du dF \quad \text{or in matrix form,} \quad W_e = \frac{1}{2} \{du\}^T \{dF\} \dots\dots\dots 3.13$$

And the internal work, W_i , is equal to the area under the stress strain graph, integrated over the volume of the element.

$$W_i = \frac{1}{2} \int_{vol} d\epsilon d\sigma dV \quad \text{or in matrix form,} \quad W_i = \frac{1}{2} \int_{vol} \{d\epsilon\}^T \{d\sigma\} dV \dots\dots\dots 3.14$$

Substituting equations 3.1 and 3.3 in to equation 3.14,

$$W_i = \frac{1}{2} \int_{vol} \{du\}^T [B]^T [C][B] \{du\} dV \dots\dots\dots 3.15$$

Equating the external and internal work and simplifying yields,

$$\{dF\} = \int_{vol} [B]^T [C][B] dV \{du\} \dots\dots\dots 3.16$$

The element stiffness matrix, $[K_e]$, relating nodal forces $\{dF\}$ to nodal displacements $\{du\}$, is therefore:

$$[K_e] = \int_{vol} [B]^T [C][B] dV \dots\dots\dots 3.17$$

Combining equations 3.16 and 3.17 gives the generalized equation of displacement based finite element equation

$$\{F\} = [K]\{u\} \dots\dots\dots 3.18$$

From which the nodal displacements are evaluated using

$$\{u\} = [K^{-1}]\{F\} \dots\dots\dots 3.19$$

The stiffness matrix for the whole system which is called the global stiffness matrix (size= total unknowns x total unknowns) can be assembled first by making all elements zero and then by placing the stiffness matrix of each element in the “place” corresponding to the degree of freedom of each point in the global system. The integral can be evaluated using the Gauss numerical integration method.

3.4 The Force Matrix

Forces acting on an element can either be externally applied loads or due to the self weight. In either case these loads can only be applied at the nodes as a point load, hence they have to be distributed to the corresponding nodes using the shape functions using the expressions below.

$$\{F\} = \int_{vol} [N]^T \{X_b\} dV + \int [N]^T \{T\} dl \dots\dots\dots 3.20$$

While, $\{X_b\} = \begin{Bmatrix} 0 \\ -\gamma \end{Bmatrix}$

Where

γ = self weight of material $\{T\}$ = externally applied uniform load

N = shape function

3.5 Numerical integration

During the evaluation of the stiffness matrix and the force or self weight matrix, integrations over the volume or the area of the element are encountered. Numerical integration is essential for practical evaluation of these integrals over the domain of the element. The common method of integration is the Gauss integration method since it uses a minimal number of sample points to achieve a desired level of accuracy.

Before integration, the number of integration points (integration order) must be selected. From which the corresponding coordinates and weights of each point is selected and the integration is performed using:-

$$\int_{-1}^1 \int_{-1}^1 f(\xi, \eta) d\xi d\eta = \sum_{j=1}^n \sum_{i=1}^n W_j W_i f(\xi, \eta) \dots\dots\dots 3.21$$

Where

ξ = natural x coordinate of the Gauss sample point

η = natural y coordinate of the Gauss sample point

W = weight of the Gauss sample point

n = number of integration points (integration order)

$f(\xi, \eta)$ = value of the function at the sample points

The selection of the integration order depends on the accuracy desired and the type of element selected for analysis. For example using six point integration for a Q4 element is waste of memory, as it will take longer time for integration, while compared to the accuracy achieved using two or three point integration. But it will be a good choice for higher order elements like Q8 and Q9 element. The minimum number of sample points that can be taken for integration is one which is for a line element. This is applied during the distribution of the external applied load to the corresponding nodes. For quadrilateral elements the integration order is the between two and seven, for example If $n=2$ there are $2 \times 2=4$ integration points and if $n=3$ there are $3 \times 3=9$ integration points which are symmetric about the natural coordinate axis with ξ and η coordinates having the same magnitude. Whereas in case of triangular elements the definition is different as illustrated below.

n	$\xi_i = \eta_i$	W_i
1	$\xi_1 = 0$	$W_1 = 2$
2	$\xi_1 = \xi_2 = \pm 0.577350269189626$	$W_1 = W_2 = 1$
3	$\xi_1 = \xi_3 = \pm 0.774596669241483$ $\xi_2 = 0$	$W_1 = W_3 = 0.5555555555555556$ $W_2 = 0.8888888888888889$
4	$\xi_1 = \xi_4 = \pm 0.861134311594053$ $\xi_2 = \xi_3 = \pm 0.339981043584856$	$W_1 = W_4 = 0.347854845137454$ $W_2 = W_3 = 0.652145154862546$
5	$\xi_1 = \xi_5 = \pm 0.906179845938664$ $\xi_2 = \xi_4 = \pm 0.538469310105683$ $\xi_3 = 0$	$W_1 = W_5 = 0.236926885056189$ $W_2 = W_4 = 0.478628670499366$ $W_3 = 0.5688888888888889$
6	$\xi_1 = \xi_6 = \pm 0.932469514203152$ $\xi_2 = \xi_5 = \pm 0.661209386466265$ $\xi_3 = \xi_4 = \pm 0.238619186003152$	$W_1 = W_6 = 0.171324492379170$ $W_2 = W_5 = 0.360761573048139$ $W_3 = W_4 = 0.467913934572691$
7	$\xi_1 = \xi_7 = \pm 0.949107912342759$ $\xi_2 = \xi_6 = \pm 0.741531185599394$ $\xi_3 = \xi_5 = \pm 0.405845151377397$ $\xi_4 = 0$	$W_1 = W_7 = 0.129484966168870$ $W_2 = W_6 = 0.279705391489277$ $W_3 = W_5 = 0.381830050505119$ $W_4 = 0.417959183673469$

Table 3.1:- Gauss point coordinates and weights in rectangular elements

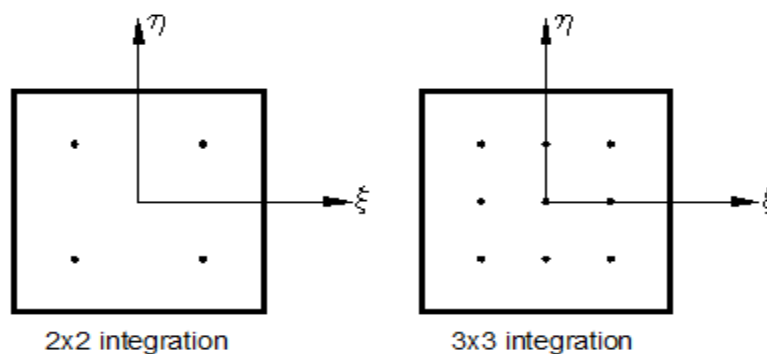


Figure 3.5:- Integration order in rectangular elements

n	ξ_i or η_i	W_i	Remark
1	$\xi_1 = \eta_1 = 0.33333333333333$	$W_1 = 1$	
3	$\xi_1 = \xi_3 = \eta_1 = \eta_2 = 0.16666666666667$ $\xi_2 = \eta_3 = 0.66666666666667$	$W_1 = W_2 =$ $W_3 = 0.33333333333333$	Used in the code if integration order is 2
7	$\xi_1 = \xi_3 = \eta_1 = \eta_2 = 0.1012865073235$ $\xi_2 = \eta_3 = 0.7974269853531$ $\xi_4 = \xi_5 = \eta_5 = \eta_6 = 0.4701420641051$ $\xi_6 = \eta_4 = 0.0597158717898$ $\xi_7 = \eta_7 = 0.33333333333333$	$W_1 = W_2 =$ $W_3 = 0.1259391805448$ $W_4 = W_5 =$ $W_6 = 0.1323941527885$ $W_7 = 0.225$	Used in the code if selected integration order is less than or equal to 5 ($2 < n \leq 5$)
13	$\xi_1 = \xi_3 = \eta_1 = \eta_2 = 0.0651301029022$ $\xi_2 = \eta_3 = 0.8697397941956$ $\xi_4 = \xi_8 = \eta_5 = \eta_9 = 0.3128654960049$ $\xi_5 = \xi_7 = \eta_6 = \eta_8 = 0.6384441885698$ $\xi_6 = \xi_9 = \eta_4 = \eta_7 = 0.0486903154253$ $\xi_{10} = \xi_{12} = \eta_{10} = \eta_{11} = 0.2603459660790$ $\xi_{13} = \eta_{13} = 0.33333333333333$	$W_1 = W_2 =$ $W_3 = 0.0533472356088$ $W_4 = W_5 = W_6 =$ $W_7 = W_8 =$ $W_9 = 0.0771137608903$ $W_{10} = W_{11} =$ $W_{12} = 0.1756152576332$ $W_{13} = -0.149570044467$	Used in the code if selected integration order is less than or equal to 7 ($5 < n \leq 7$)

Table 3.2:- Gauss point coordinates and weights in triangular elements

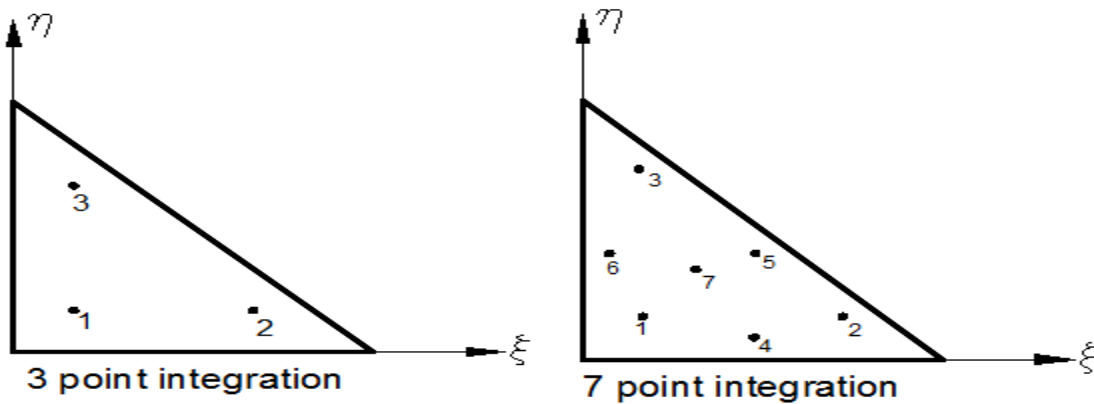


Figure 3.6:- Integration order in triangular elements

CHAPTER 4

PLANE STRAIN AND STRESS INVARIANTS

4.1 Plane Strain

In geotechnical analysis, problems that have one very large spatial dimension compared to the others are often encountered. In these situations, it is often reasonable to assume that the primary field variables in the long direction are zero, such analysis is called plane strain analysis. In the displacement based finite element method, the primary field variable is the displacement, which may vary throughout the region being modeled. Stresses and strains are secondary variables and can be evaluated from the displacements.

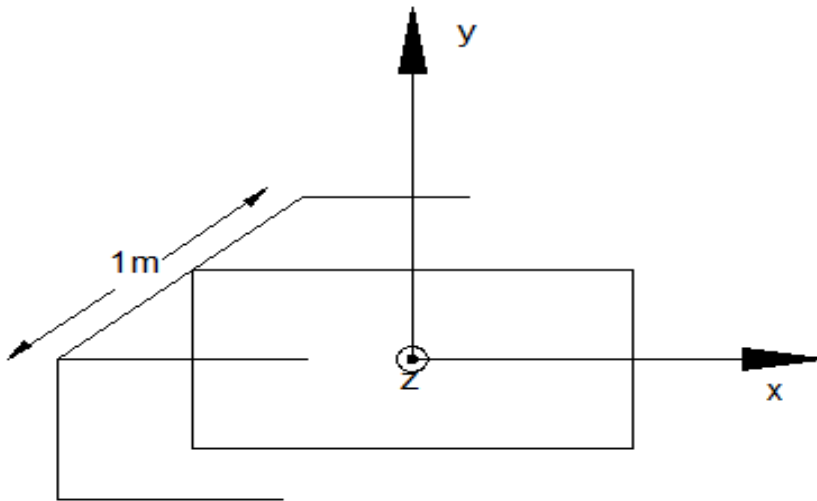


Figure 4.1:- A plane strain element in the xy plane

The stresses and strains in a plain strain element are given in matrix form as:-

$$\{\epsilon\} = \begin{Bmatrix} \epsilon_{xx} \\ \epsilon_{yy} \\ \epsilon_{xy} \\ \epsilon_{zz} \end{Bmatrix} = [B]\{u\} \dots\dots\dots 4.1$$

In plane strain case the strain in the z direction is zero, $\epsilon_{zz} = 0$.

$$\{\sigma\} = \begin{Bmatrix} \sigma_{xx} \\ \sigma_{yy} \\ \tau_{xy} \\ \sigma_{zz} \end{Bmatrix} = [C]\{\varepsilon\} \dots\dots\dots 4.2$$

While, $\sigma_{zz} = \nu(\sigma_{xx} + \sigma_{yy}) \dots\dots\dots 4.3$

$$\{u\} = \begin{Bmatrix} u \\ v \end{Bmatrix} \dots\dots\dots 4.4$$

$$C = \frac{E}{(1+\nu)(1-2\nu)} \begin{pmatrix} 1-\nu & \nu & 0 & \nu \\ \nu & 1-\nu & 0 & \nu \\ 0 & 0 & \frac{1-2\nu}{2} & 0 \\ \nu & \nu & 0 & 1-\nu \end{pmatrix} \dots\dots\dots 4.5$$

Where

ε = the strain vector

ν = displacement in y direction

B = the strain displacement matrix

u = displacement in x direction

σ = the stress vector

C = the linear elastic constitutive matrix

u= the displacement vector

ν = poisson's ratio

E = young's modulus

During the calculation of the stiffness matrix, internal nodal reactions and others, the strain displacement matrix (B) which was originally a 3X8 matrix should be modified into a 4X8 matrix so that matrix multiplication is possible because the stress, the strain and the constitutive matrix for plane strain case have a fourth row for the accommodation of stresses and strains in the z direction.

4.2 Stress Invariants

As shown above, a plane strain state of stress can be specified with four components with a fixed coordinate system (σ_{xx} , σ_{yy} , τ_{xy} , σ_{zz}). The magnitude of these stresses depends on the orientation of the chosen coordinate system. Another way of specifying stress state is using the principal stresses (σ_1 , σ_2 , σ_3), which act in the same direction and have the same magnitude for a given stress state, regardless of the chosen orientation of coordinate axes. Principal stresses are therefore frequently used to show a stress point in space. The plot of the principal stresses on a three dimensional graph, known as principal stress space, shows the stress state and stress paths. Figure 3.2 shows a principal stress space.

In the figure, a line representing a hydrostatic stress state ($\sigma_1 = \sigma_2 = \sigma_3$) is shown, this line is called the space diagonal. The plane normal to the space diagonal, represented by $\sigma_1 + \sigma_2 + \sigma_3 = \text{Constant}$ and bound by the condition that all principal stress are greater than zero, is called the deviatoric plane. The shaded region on the deviatoric plane represents the region where $\sigma_1 \geq \sigma_2 \geq \sigma_3$. Another way to describe the state of stress geometrically is to use stress invariants in the principal stress space.

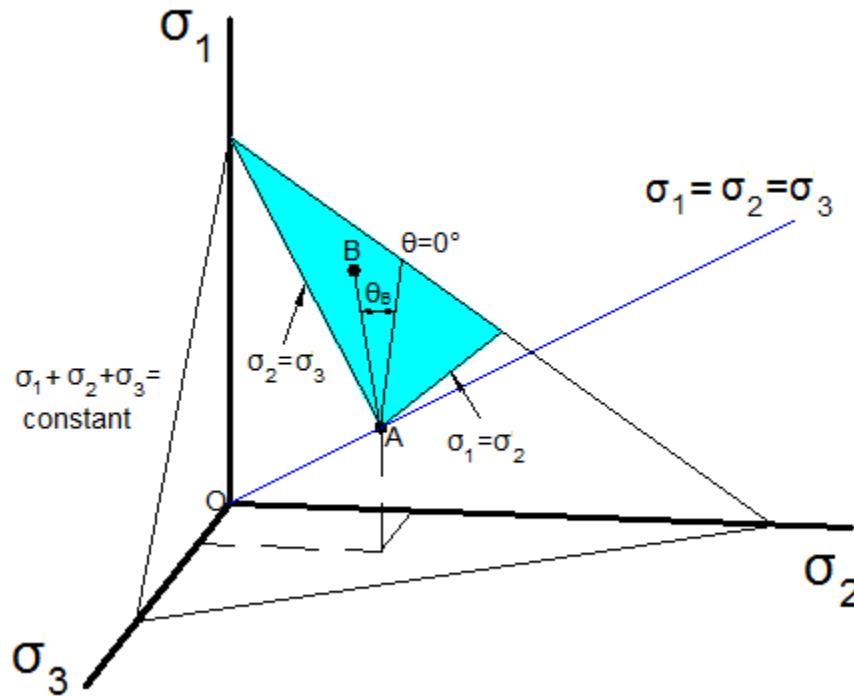


Figure 4.2:- Principal stress space

The location of a stress point (let point B, in figure 4.2) in the stress space can be described numerically using the following invariants

$$s = \frac{\sigma_1 + \sigma_2 + \sigma_3}{\sqrt{3}} \dots\dots\dots 4.6$$

$$t = \frac{1}{\sqrt{3}} \sqrt{((\sigma_1 - \sigma_2)^2 + (\sigma_2 - \sigma_3)^2 + (\sigma_3 - \sigma_1)^2)} \dots\dots\dots 4.7$$

$$\theta = \tan^{-1} \left(\frac{1}{\sqrt{3}} \left(\frac{2(\sigma_2 - \sigma_3)}{(\sigma_1 - \sigma_3)} - 1 \right) \right) \dots\dots\dots 4.8$$

The distance along the space diagonal from the origin to the deviatoric plane (OA) is represented by the stress invariant s, while t is the distance from the space diagonal to the stress point (AB) and Lode angle (Θ) is the angular position of the stress point in the deviatoric plane. The value of Lode’s angle on the plane varies between the extremities of the intermediate principal stresses as:-

Θ=+30°, triaxial compression, $|\sigma'_1| \geq |\sigma'_2| = |\sigma'_3|$ and
 Θ=-30°, triaxial extension, $|\sigma'_1| = |\sigma'_2| \geq |\sigma'_3|$

In Geotechnical Engineering, the most widely used invariants are mean and deviatoric stress which can be expressed in terms of Cartesian stresses as:-

$$\sigma_m = p = \frac{\sigma_{xx} + \sigma_{yy} + \sigma_{zz}}{3} \dots\dots\dots 4.9$$

Rewriting equation 4.7 in Cartesian form,

$$t = \frac{1}{\sqrt{3}} \sqrt{((\sigma_{xx} - \sigma_{yy})^2 + (\sigma_{yy} - \sigma_{zz})^2 + (\sigma_{zz} - \sigma_{xx})^2 + 6\tau_{xy}^2)} \dots\dots\dots 4.10$$

From which, $q = \sqrt{\frac{3}{2}} \cdot t \dots\dots\dots 4.11$

Similarly the Lode angle is given by

$$\theta = \frac{1}{3} \sin^{-1} \left(\frac{-3\sqrt{6}J_3}{t^3} \right) \dots\dots\dots 4.12$$

where

$$J_3 = s_x s_y s_z - s_z \tau_{xy}^2 \dots\dots\dots 4.13$$

$$s_x = \sigma_{xx} - p \dots\dots\dots 4.14$$

$$s_y = \sigma_{yy} - p \dots\dots\dots 4.15$$

$$s_z = \sigma_{zz} - p \dots\dots\dots 4.16$$

The principal stresses can be expressed in terms of these invariants as:-

$$\sigma_1 = p + \frac{2}{3} q \sin\left(\theta - \frac{2\pi}{3}\right) \dots\dots\dots 4.17$$

$$\sigma_2 = p + \frac{2}{3} q \sin(\theta) \dots\dots\dots 4.18$$

$$\sigma_3 = p + \frac{2}{3} q \sin\left(\theta + \frac{2\pi}{3}\right) \dots\dots\dots 4.19$$

Note that, since compression is negative, in these equations the value of p is negative and that of q is positive.

CHAPTER 5

ELASTO PLASTICITY

5.1 Introduction

Due to the complexity of real soil behavior, a single constitutive model that can describe all facts of behavior, with a reasonable number of input parameters, has not been achieved. Consequently, there are many soil models available, each of which has different advantages and disadvantages. The Mohr-Coulomb material model is one of the well known elastic perfectly plastic soil models. The main features of this model are discussed below.

5.2 Yield Function

Based on the current stress state, an Elasto-plastic material behaves either as an elastic solid or a plastic solid. The transition from elasticity to plasticity is described by the yield criterion which forms a surface in three dimensional principle stress space. Stress states lying within the yield surface are regarded as elastic, while stress states lying on the yield surface are plastic. As the material deforms plastically, the stresses must remain on the yield surface and so stress states lying outside the yield surface are inadmissible and must be redistributed through an iterative procedure. Algebraically, these surfaces are expressed in terms of yield or failure function (F). This function, which has units of stress, is dependent on material properties and invariant combinations of the stress components. The derivation of the Mohr Coulomb yield function is discussed below.

The shear strength of soil at failure (τ_f) is given by the Mohr Coulomb failure criterion, which can be expressed as:-

$$\tau_f = c + \sigma_n \tan \phi \dots\dots\dots 5.1$$

Or alternatively

$$\tau_f = (\sigma_n + a) \tan \phi \dots\dots\dots 5.2$$

Where

$$a = c \cdot \cot \phi \dots\dots\dots 5.3$$

The Mohr Coulomb failure criterion can be plotted on shear strength versus normal stress graph as shown in figure 5.1.

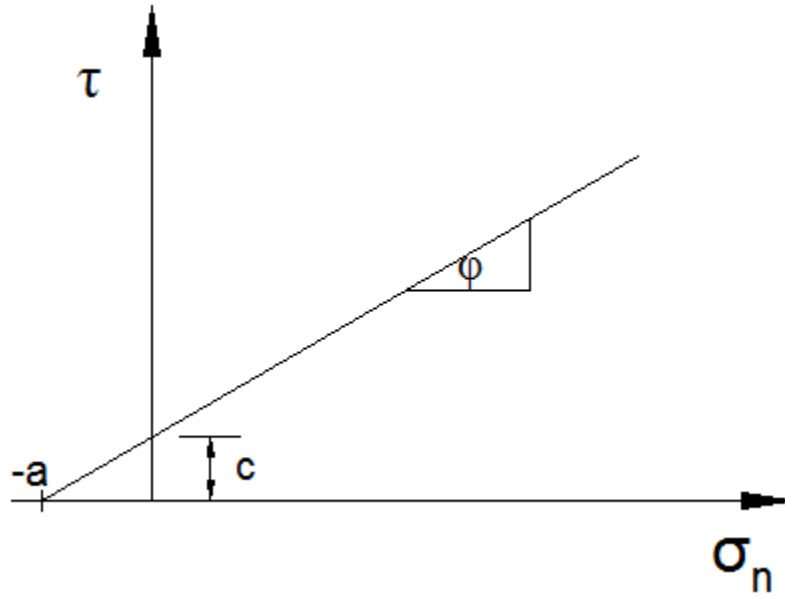


Figure 5.1:- The Mohr Coulomb failure criterion

In three dimensional principal stress space the plot of this failure criterion is an irregular hexagonal pyramid.

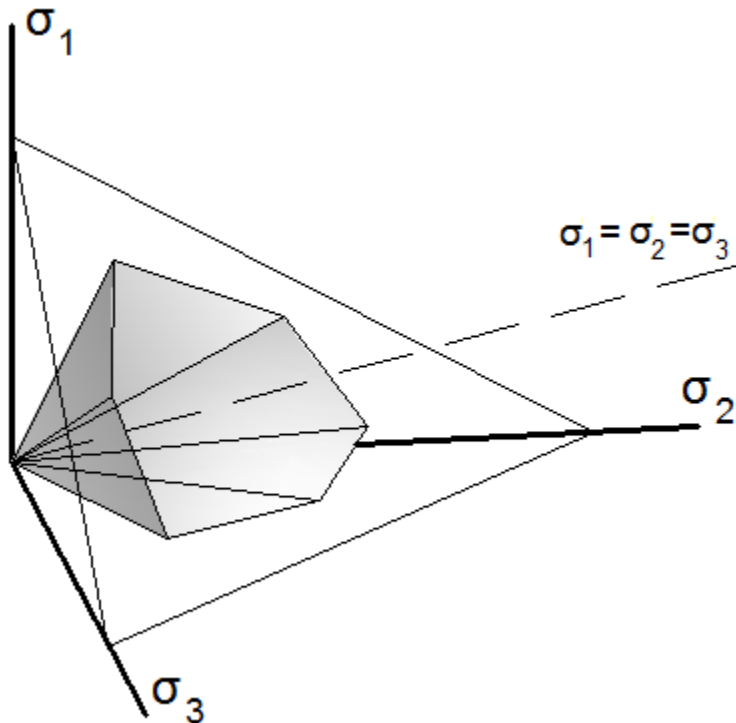


Figure 5.2:- The Mohr Coulomb yield surface in 3 dimensional principal stress space

Since the intermediate principal stress (σ_2) does not affect the Coulomb criterion, and including the effect of cohesion (attraction) it can also be plotted in two dimensional principal stress space.

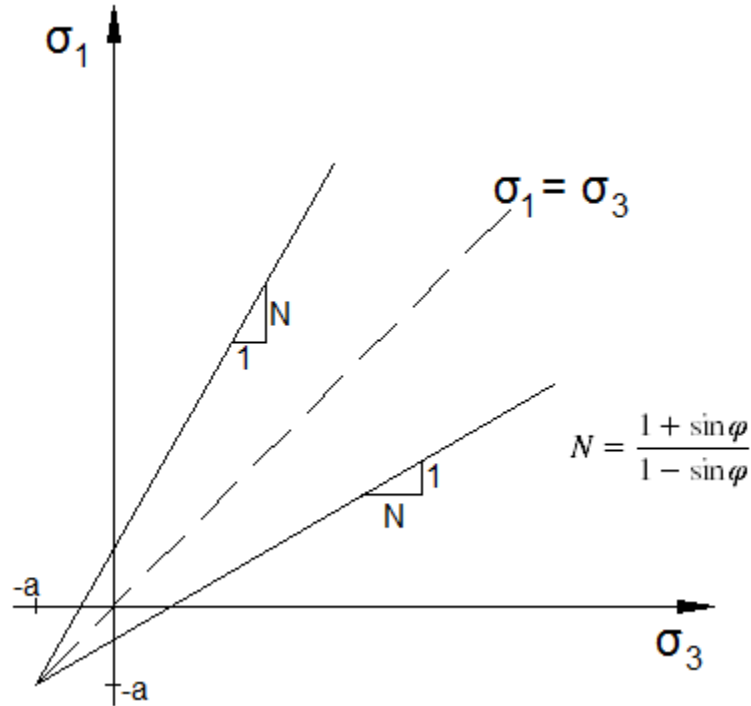


Figure 5.3:- The Mohr Coulomb yield surface in 2 dimensional principal stress space

From figure 5.3, the yield function defining the Mohr-Coulomb failure criterion in terms of the maximum and minimum principal stresses will be

$$F = \left(\frac{\sigma_1 + \sigma_3}{2} \right) \sin \phi - \left(\frac{\sigma_1 - \sigma_3}{2} \right) - c \cdot \cos \phi = 0 \dots\dots\dots 5.4$$

Substituting equation 4.17 and 4.19 in to equation 5.4, a general invariant form of the yield function will become,

$$F = p \sin \phi + q \left(\frac{\cos \theta}{\sqrt{3}} - \frac{\sin \theta \sin \phi}{3} \right) - c \cdot \cos \phi = 0 \dots\dots\dots 5.5$$

For a triaxial compression case, where $\Theta=30$ this can be written as

$$F = p \sin \phi + q \left(0.5 - \frac{\sin \phi}{6} \right) - c \cdot \cos \phi = 0 \dots\dots\dots 5.6$$

$$\text{Since, } c \cdot \cos \phi = a \cdot \sin \phi$$

With attraction (a = negative), Equation 5.6 can be rearranged as

$$q = (p - a) \left(\frac{6 \sin \phi}{3 - \sin \phi} \right) \dots\dots\dots 5.7$$

$$\text{With, } M = \frac{6 \sin \phi}{3 - \sin \phi} \dots\dots\dots 5.8$$

Equation 5.7 simplifies to

$$q = M(p - a) \dots\dots\dots 5.9$$

Equation 5.9 is the plot of the Mohr-Coulomb failure criterion in a triaxial stress space commonly known as, p-q diagram.

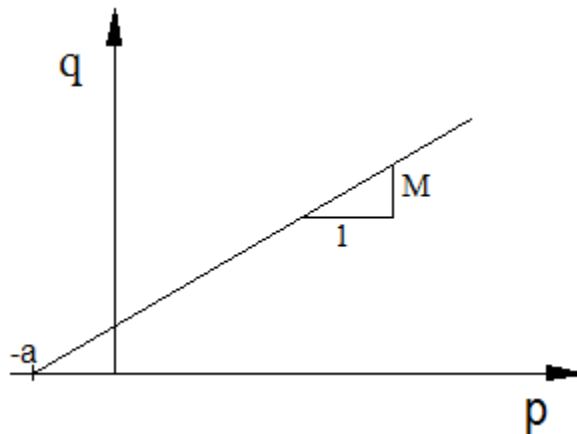


Figure 5.4:- The Mohr Coulomb yield criterion in a p-q diagram for triaxial compression case

5.3 Plastic potential Function and Flow rule

The plastic potential function, G , specifies the direction of the plastic strain increment, which is proportional to the derivative of G with respect to the corresponding stress.

$$\{\Delta\varepsilon^p\} = \Delta\lambda \left(\frac{\partial G}{\partial \sigma} \right) \dots\dots\dots 5.10$$

Where $\Delta\lambda$ is a scalar quantity called plastic multiplier. This scalar is not a property of the material but rather a property of the flow conditions. The plastic potential function is defined by

$$G = p \sin \psi + q \left(\frac{\cos \theta}{\sqrt{3}} - \frac{\sin \theta \sin \psi}{3} \right) - c \cdot \cos \psi = 0 \dots\dots\dots 5.11.$$

When the friction angle (ϕ) and dilatancy angle (ψ) are equal, the flow condition is said to be associated flow and the yield and plastic potential function are identical and equation 5.10 can be written as

$$\{\Delta\varepsilon^p\} = \Delta\lambda \left(\frac{\partial F}{\partial \sigma} \right) \dots\dots\dots 5.12$$

Otherwise the flow is said to be non-associated flow.

5.4 The Plastic potential Derivative

During the calculation of plastic strain increments, the plastic potential derivative should be evaluated according to chain rule.

$$\left(\frac{\partial G}{\partial \sigma} \right) = \left(\frac{\partial G}{\partial p} \right) \left(\frac{\partial p}{\partial \sigma} \right) + \left(\frac{\partial G}{\partial q} \right) \left(\frac{\partial q}{\partial \sigma} \right) + \left(\frac{\partial G}{\partial \theta} \right) \left(\frac{\partial \theta}{\partial \sigma} \right) \dots\dots\dots 5.13$$

Expressing the quantities q and θ in terms of J_2 and J_3 simplifies the derivative process,

$$\left(\frac{\partial G}{\partial \sigma} \right) = \left(\frac{\partial G}{\partial p} \right) \left(\frac{\partial p}{\partial \sigma} \right) + \left(\frac{\partial G}{\partial J_2} \right) \left(\frac{\partial J_2}{\partial \sigma} \right) + \left(\frac{\partial G}{\partial J_3} \right) \left(\frac{\partial J_3}{\partial \sigma} \right) \dots\dots\dots 5.14$$

Where

$$J_2 = \frac{q^2}{3} \quad \text{and} \quad J_3 = -0.385 \sin(3\theta)$$

After differentiation, the terms q and Θ are restored in the equations. Expressing equation 5.14 in separate form for simplicity,

$$\left(\frac{\partial G}{\partial \sigma}\right) = (dg_1 \cdot m_1 + dg_2 \cdot m_2 + dg_3 \cdot m_3)\{\sigma\} \dots\dots\dots 5.15$$

Where

$$\left(\frac{\partial G}{\partial p}\right) = dg_1 = \sin(\psi) \dots\dots\dots 5.16$$

$$\left(\frac{\partial G}{\partial J_2}\right) = dg_2 = \frac{\sqrt{3} \cos(\theta)}{2q} \left[1 + (\tan \theta \tan 3\theta) + \left(\frac{\sin \psi}{\sqrt{3}} (\tan 3\theta - \tan \theta)\right)\right] \dots\dots 5.17$$

$$\left(\frac{\partial G}{\partial J_3}\right) = dg_3 = 1.5 \left(\frac{\sqrt{3} \sin \theta + \sin \psi \cos \theta}{q^2 \cos 3\theta}\right) \dots\dots\dots 5.18$$

$$\left(\frac{\partial p}{\partial \sigma}\right) = m_1\{\sigma\} \dots\dots\dots 5.19$$

$$\left(\frac{\partial J_2}{\partial \sigma}\right) = m_2\{\sigma\} \dots\dots\dots 5.20$$

$$\left(\frac{\partial J_3}{\partial \sigma}\right) = m_3\{\sigma\} \dots\dots\dots 5.21$$

$$m_1 = \frac{1}{3(\sigma_{xx} + \sigma_{yy} + \sigma_{zz})} \begin{bmatrix} 1 & 1 & 0 & 1 \\ 1 & 1 & 0 & 1 \\ 0 & 0 & 0 & 0 \\ 1 & 1 & 0 & 1 \end{bmatrix} \dots\dots\dots 5.22$$

$$m_2 = \frac{1}{3} \begin{bmatrix} 2 & -1 & 0 & -1 \\ -1 & 2 & 0 & -1 \\ 0 & 0 & 6 & 0 \\ -1 & -1 & 0 & 2 \end{bmatrix} \dots\dots\dots 5.23$$

$$m_3 = \frac{1}{3} \begin{bmatrix} s_x & s_z & \tau_{xy} & s_z \\ s_z & s_y & \tau_{xy} & s_x \\ \tau_{xy} & \tau_{xy} & -3s_z & -2\tau_{xy} \\ s_z & s_x & -2\tau_{xy} & s_z \end{bmatrix} \dots\dots\dots 5.24$$

$$\{\sigma\} = \begin{Bmatrix} \sigma_{xx} \\ \sigma_{yy} \\ \tau_{xy} \\ \sigma_{zz} \end{Bmatrix} \dots\dots\dots 5.25$$

$$s_x = \sigma_{xx} - p \dots\dots\dots 5.26$$

$$s_y = \sigma_{yy} - p \dots\dots\dots 5.27$$

$$s_z = \sigma_{zz} - p \dots\dots\dots 5.28$$

The yield and plastic potential functions of a Mohr Coulomb material are a function of the Lode angle, which in turn is variable depending on the stress state, as shown in equations 5.5 and 5.11. During the calculation of the plastic strains the plastic potential derivative becomes indeterminate when the lode angle approaches the corners of the hexagonal surface ($\pm 30^\circ$).

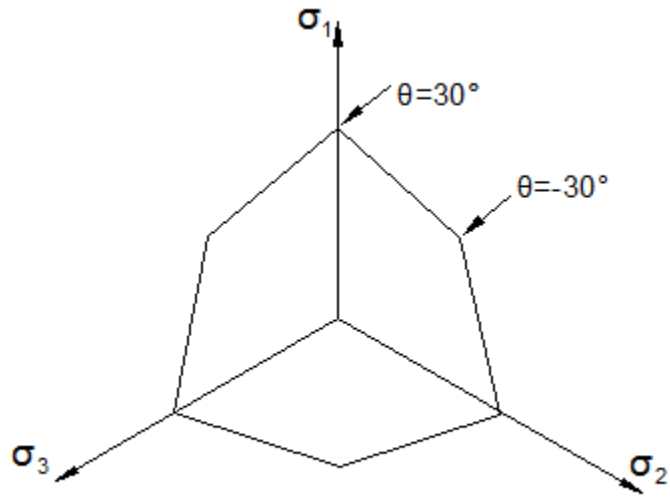


Figure 5.5:- The corners of Mohr Coulomb yield surface

This problem can be resolved by smoothening the corners of the yield surface if the lode angle gets close enough to $\pm 30^\circ$. This is achieved by substituting $\Theta = \pm 30^\circ$ in equation 5.11 depending on whether it is getting close to triaxial compression or extension. In these cases the terms dg_2 and dg_3 in equations 5.17 and 5.18 respectively should be modified as,

For $\Theta \approx 30^\circ$ triaxial compression case

$$dg_2 = \frac{0.25}{q} (3 - \sin \psi) \dots\dots\dots 5.29$$

$$dg_3 = 0 \dots\dots\dots 5.30$$

For $\Theta \approx -30^\circ$ triaxial extension case

$$dg_2 = \frac{0.25}{q} (3 + \sin \psi) \dots\dots\dots 5.31$$

$$dg_3 = 0 \dots\dots\dots 5.32$$

CHAPTER 6

ELASTO PLASTICITY IN FINITE ELEMENT ANALYSIS

6.1 Introduction

Non-linear problems in finite element analysis can be grouped into the following three categories, based on the sources of non-linearity:

1. Material Nonlinearity Problems
2. Geometric Nonlinearity Problems and
3. Both material and Geometric Nonlinearity Problems.

Here, only the case of Material Nonlinearity is discussed. Nonlinearity due to an Elasto Plastic material causes the governing finite element equations to be reduced to the following incremental form,

$$[K]\{\Delta u\}^i = \{\Delta f\}^i \dots\dots\dots 6.1$$

Where

- [K] = the global system stiffness matrix
- { Δu } = the vector of incremental nodal displacements
- { Δf } = the vector of incremental nodal forces and
- i = the increment number

To obtain a solution to a boundary value problem, the change in boundary conditions is applied in a series of increments and for each increment equation 6.1 must be solved. The final solution is obtained by summing the results of each increment. Due to the nonlinear constitutive behavior of the Mohr Coulomb material, the global stiffness matrix is dependent on the current stress and strain levels and therefore is not constant, but varies over an increment.

In practical Finite Element Analysis, two main types of solution strategies can be used to solve this equation. The first method, known as the Tangent Stiffness Method, considers the reduction of the material stiffness as failure is approached. Hence the global stiffness matrix is updated at each iteration to achieve convergence. In this method the numbers of iterations needed to achieve convergence are relatively small but the cost of computer memory, due to the formation of the global stiffness matrix at each iteration, is very large and hence is not used here.

In the second method, known as the Constant Stiffness Method, the stiffness matrix is formed only once and is kept constant throughout which makes the memory cost small due to this reason this method is used here.

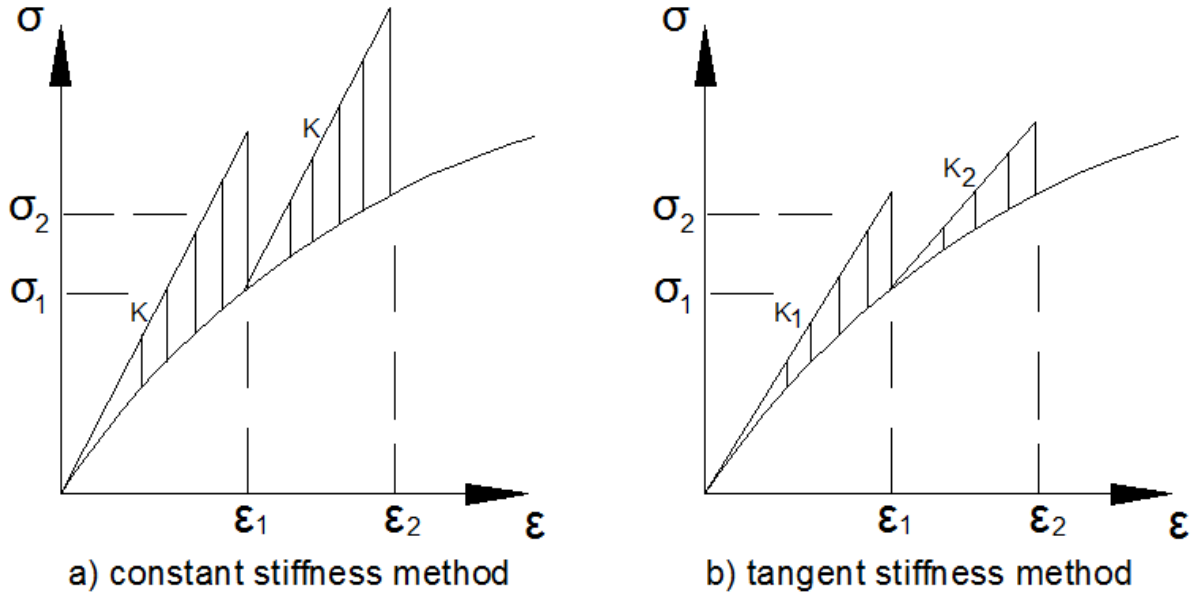


Figure 6.1 Nonlinear solution strategies

Constant stiffness methods uses repeated elastic solutions to achieve convergence by iteratively varying the loads on the system. For each load increment the displacement increment must be solved, which in turn is used to find the total strain increments using the strain-displacement matrix.

$$\{\Delta u\}^i = [K]^{-1} \{\Delta f\}^i \dots\dots\dots 6.2$$

$$\{\Delta \varepsilon\}^i = [B] \{\Delta u\}^i \dots\dots\dots 6.3$$

If the material is yielding, the strains will contain both elastic and plastic parts,

$$\{\Delta \varepsilon\}^i = \{\Delta \varepsilon^e\}^i + \{\Delta \varepsilon^p\}^i \dots\dots\dots 6.4$$

But it is only the elastic strains that generate stresses through the elastic constitutive matrix C,

$$\{\Delta \sigma\}^i = [C] \{\Delta \varepsilon^e\}^i \dots\dots\dots 6.5$$

These stresses are then added to the previously existing stresses and the resulting stresses are used to check whether yield is violated or not.

$$\{\sigma\}^i = \{\sigma\}^{i-1} + \{\Delta\sigma\}^i \dots\dots\dots 6.6$$

If the stress is below the yield surface ($F < 0$) then it is in the elastic state and the whole process is repeated with an increased load in equation 6.2.

On the other hand, if the stress is outside the yield surface ($F > 0$), the stress is over estimated and it should be redistributed so that it can be brought back as near as possible to the yield surface ($F = 0$). At this stage the material is behaving plastically and the plastic strains should be calculated. This calculation requires the differentiation of the yield and plastic potential derivatives with respect to stress and the general formula is given by Vermeer (1979) as:-

$$\{\Delta\varepsilon^p\} = \Delta\lambda \left[(1-\omega) \left(\frac{\partial G}{\partial \sigma} \right)^{i-1} + \omega \left(\frac{\partial G}{\partial \sigma} \right)^i \right] \dots\dots\dots 6.7$$

Where $\Delta\lambda$ is the increment of the plastic multiplier and ω is a parameter which depends on the type of time integration used. For $\omega = 0$ the integration is called explicit and for $\omega = 1$ it is called implicit. During implicit integration the derivative of the plastic potential function has to be calculated at the current stress, which is unknown, this complicates the coding process and is not used here though it is advantageous. Hence using the explicit integration for which differentials are calculated at the previous known stress, equation 6.7 will be reduced to:-

$$\{\Delta\varepsilon^p\} = \Delta\lambda \left(\frac{\partial G}{\partial \sigma} \right)^{i-1} \dots\dots\dots 6.8$$

These plastic strains are used to calculate the ‘excess’ load at the nodes and these loads are used in addition to the external load increment for the next iteration. These load can be generated using two methods, the Visco-plastic method and the initial stress method. The Visco-plastic method, which is used here, is discussed next.

6.2 Explicit integration with visco-plastic method

As outlined before, if the stress state violates yield, the stress state can only be sustained momentarily and plastic straining (referred to as visco-plastic strains in this method) occurs. The magnitude of the visco-plastic strain rate is determined by the value of the yield function, which is a measure of the degree by which the current stress state exceeds the yield condition. The visco-plastic strains increase with time or iteration, causing the material to relax with a reduction in the yield function and hence the visco-plastic strain rate.

For time independent Elasto-plastic materials, the visco-plastic strain rate is given by,

$$\{\delta\epsilon^{vp}\}^i = \Delta t * F * \left\{ \frac{\partial G}{\partial \sigma} \right\} \dots\dots\dots 6.9$$

Where

Δt = a pseudo time step and for a Mohr Coulomb material it is given by,

$$\Delta t = \frac{4(1+\nu)(1-2\nu)}{E(1-2\nu + \sin^2 \phi)} \dots\dots\dots 6.10$$

F = the value of the yield function at the current stress.

$\{\partial G/\partial \sigma\}$ = the plastic potential derivative at the current stress state.

This strain is accumulated from one iteration to the next as

$$\{\Delta\epsilon^{vp}\}^i = \{\Delta\epsilon^{vp}\}^{i-1} + \{\delta\epsilon^{vp}\}^i \dots\dots\dots 6.11$$

When the visco-plastic strain rate in equation 6.8 becomes very small (at convergence), the accumulated visco-plastic strain and the associated stress change are equal to the incremental plastic strain and stress change respectively.

The stress increment associated with the visco-plastic strain change is,

$$\{\Delta\sigma^{vp}\}^i = [C]\{\delta\epsilon^{vp}\}^i \dots\dots\dots 6.12$$

And the associated nodal force vectors would be

$$\{\Delta r\}^i = \int_{vol} B^T C \{\delta\epsilon^{vp}\}^i dV \dots\dots\dots 6.13$$

These nodal reactions are accumulated at each iteration by summing equation 6.13 for all elements with yielding gauss points. Finally the applied load vector would take the form,

$$\{\Delta f\}^i = \{\Delta f\}^{i-1} + \sum_{\substack{\text{all} \\ \text{elements}}} \int_{vol} B^T C \{\delta \varepsilon^{vp}\}^i dV \dots\dots\dots 6.14$$

This load is then used in equation 6.2 and the process is repeated until convergence is achieved. If convergence is achieved, the stresses, strains and displacements from the last iteration are recorded for use in the next load step.

6.3 Convergence

During the iteration process from equation 6.2 through 6.14 the process is repeated until the stresses are close enough to the yield surface within a certain error which should be less than a predefined tolerance which is set to be 0.01. When this criterion is met the solution is said to be converged. Convergence is indicated by either of the following

- The incremental displacements from equation 6.2 is nearly the same from one iteration to the next
- The yield function, F is too small
- The visco-plastic strain is too small
- The stress increment from equation 6.5 is nearly the same from one iteration to the next

To calculate the error the displacement condition is used. In this condition the difference between the absolute maximum of the incremental displacements in iteration i and iteration i-1 should be less than the predefined tolerance. The whole load stepping is shown in the flow chart in the next page.

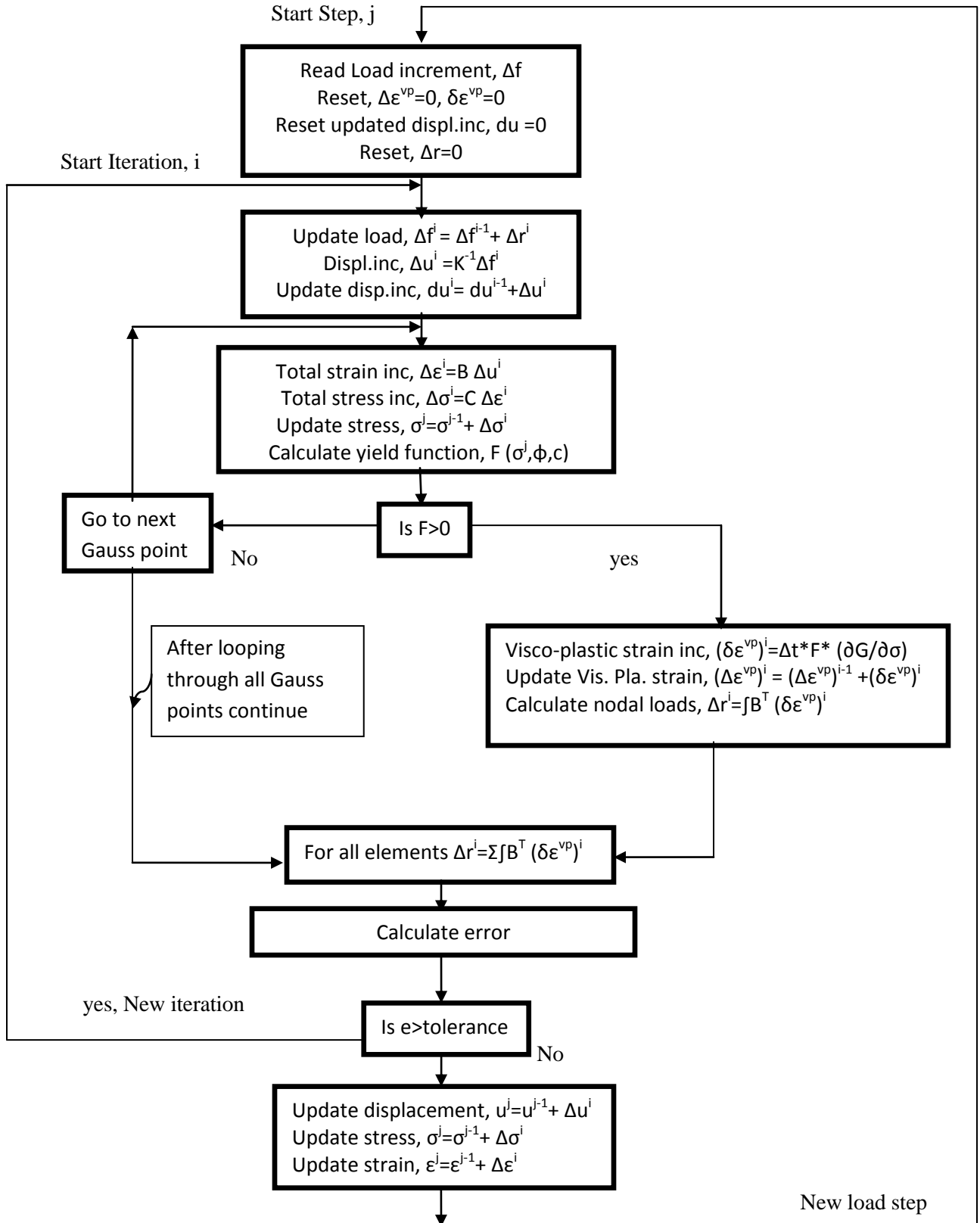


Figure 6.2:- A single load step

CHAPTER 7

USING THE MATLAB CODE

7.1 Introduction

The functions that have been developed are arranged in four groups depending on what they primarily perform.

1. Meshing:- contains functions used for discretization of the domain.
2. Integrations:- contains functions used for the Gauss numerical integration.
3. Subfunctions:- contains the main functions for finite element analysis.
4. Plotting:- contains functions that are used for plotting the output results.

The description of each function and the functions themselves are provided in appendix B and C respectively. Additionally the descriptions of frequently used variables in these functions are also given in appendix A. Two files are presented depending on the analysis type to be performed.

1. `biaxialtest.m` :- Performs biaxial test for a Mohr Coulomb soil.
2. `bearingcapacity.m` :- calculates bearing capacity of a strip footing on a Mohr Coulomb soil.

7.2 Input parameters

The input parameters for these analyses are shown in the following tables.

parameter name	Variable name	Unit of input values	Remark
Length of region	L	Meter	
Depth of region	D	Meter	
No. of Element in x-dir.	numx		
No. of Element in y-dir.	numy		
Young's modulus	E	KPa	
Poissons ratio	nu		
cohesion	c	KPa	
Angle of friction	phi	°	
Dilatancy angle	tsi	°	
Initial vertical stress	s_v	KPa	Negative
Initial horizontal stress	s_h	KPa	Negative
Load increment	df	KN/m	Negative
Number of steps	nsteps		
Type of element	elemType	'Q9'/'Q8'/'Q4' 'T6'/'T3'	Capitalized And quoted
Order of integration	normal_order		

Table 7.1:- Input parameters for `biaxialtest.m`

parameter name	Variable name	Unit of input values	Remark
Length of region	L	Meter	
Depth of region	D	Meter	
Number of Elements in x-direction	numx		
Number of Elements in y-direction	numy		
Young's modulus	E	KPa	
Poissons ratio	nu		
cohesion	c	KPa	
Angle of friction	phi	°	
Dilatancy angle	tsi	°	
Unit weight	gamma	KN/m ³	
Load increment	df	KN/m	Negative
Number of steps	nsteps		
x-coordinate of load starting point	load_edge1		Usually at 0
x-coordinate of load ending point	load_edge2		known node coordinate
Type of element	elemType	'Q9'/'Q8'/'Q4' 'T6'/'T3'	Capitalized And quoted
Order of integration	normal_order		

Table 7.2:- Input parameters for bearingcapacity.m

The minimum and maximum values for the order of integration depends on the type of element selected. For T6, Q8 and Q9 elements the minimum value is 3 while for Q4 and T3 elements it is 2. For Q4, Q8 and Q9 elements the maximum value is 7 while for T3 and T6 elements it is 6. Additionally the number of elements in the x- direction for T3 and T6 elements is twice the value inserted for numx. For example, to get a region with 30 elements in x direction and 40 in y direction the input for triangular elements should be , numx=15 and numy=40 which gives a total of 1200 elements.

Care should be taken when inserting the ending location of the applied load increment (load_edge2) for bearingcapacity.m. This is because the user has to provide the coordinate by dividing the length (L) with the number of elements in x direction which gives the spacing between each node from which the value of the load ending point coordinate can be calculated.

In both cases the user have to open the file and insert the input parameters directly which are located at the beginning of the script files. After giving the values and saving, it can be executing using the Run (F5) button. Eventually calculation progress will show up and then the outputs will be displayed.

7.3 Outputs

The outputs of the analysis are displayed automatically when the calculation is finished. The default way of displaying the calculation results is in picture form these are:-

- The finite element mesh.
- The deformed finite element mesh
- Displacement shadings (u_x and u_y)
- Stress shadings (σ_{xx} , σ_{yy} , τ_{xy} and σ_{zz})
- Deviatoric stress vs. mean stress plot.
- Vertical stress vs. vertical strain plot.
- Deviatoric stress vs. volumetric strain plot.
- Load vs. vertical displacement plot.

The mesh and the shading diagrams show the overall property of the domain while the curves are plotted for a stress point in the element at the upper left corner. In this element the first gauss point is selected as a stress point for rectangular elements and the third gauss point in triangular elements. Examples of outputs from a bearing capacity problem are shown in chapter 8, pages 49-53.

Another way of accessing the outputs is to see the numerical results to several decimal places at each node on the Command window by using the following commands.

Command	Result displayed
<code>sigma(:, :, 1)</code>	σ_{xx} at all gauss points
<code>sigma(:, :, 2)</code>	σ_{yy} at all gauss points
<code>sigma(:, :, 3)</code>	τ_{xy} at all gauss points
<code>sigma(:, :, 4)</code>	σ_{zz} at all gauss points
<code>stress (1, :, :)</code>	σ_{xx} at all gauss points
<code>stress (2, :, :)</code>	σ_{yy} at all gauss points
<code>stress (3, :, :)</code>	τ_{xy} at all gauss points
<code>stress (4, :, :)</code>	σ_{zz} at all gauss points
<code>u_x</code>	Horizontal displacement at each node
<code>u_y</code>	Vertical displacement at each node
<code>min(u_y)</code>	Largest downward displacement
<code>max(u_y)</code>	Largest upward displacement
<code>min(u_x)</code>	Largest left ward displacement
<code>max(u_x)</code>	Largest right ward displacement

Table 7.3:- commands for numerical outputs

In table 7.3 both sigma and stress shows the stresses at the gauss points but in a different manner.

`sigma(:, :, :)`

The first position is the element number which is between 1 and number of elements in the domain. The second position is the gauss point number which is between 1 and kk (the number of gauss points in the element). The third position is the stress name as shown in the table above. For example `sigma(4,1,2)` displays the value of σ_{yy} for the first gauss point in the fourth element.

`stress(:, :, :)`

The first position is the stress name as shown in the table above. The second position is the gauss point number which is between 1 and kk (the number of gauss points in the element). The third position is element number which is between 1 and number of elements in the domain. For example `stress(2,1,4)` displays the value of σ_{yy} for the first gauss point in the fourth element.

CHAPTER 8

VERIFICATION AND CONCLUSION

8.1 Introduction

Now all the required functions developed, it is time to see the results and compare these results with values from theoretical solutions and the PLAXIS software. For this purpose two types of problem sets are discussed, biaxial test and bearing capacity problem.

8.2 Biaxial test

A typical biaxial test setup used is shown in figure 8.1. The soil is initially subjected to an initial stress in the axial (σ_v) and lateral (σ_h) directions. When the test start incremental load is applied in the vertical direction until failure.

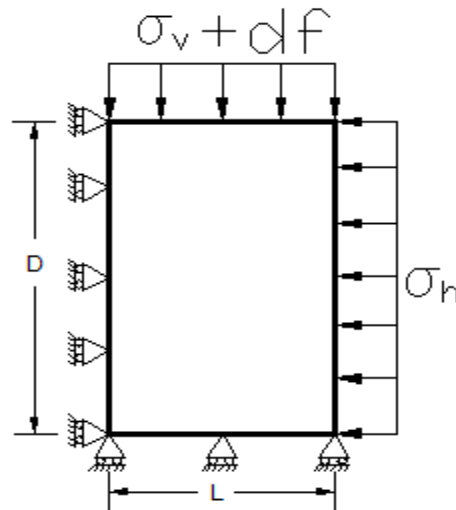


Figure 8.1:- a typical biaxial test setup

In this example the following input parameters are used

Modulus of elasticity, $E=10,000\text{KPa}$	Poisson's ratio, $\nu=0.25$,
Angle of friction, $\phi=30^\circ$	Dilatancy angle, $\psi=30^\circ$
Cohesion, $c=1\text{KPa}$	unit weight, $\gamma=0$
Initial vertical load, $\sigma_v = -1\text{KPa}$	Initial horizontal load, $\sigma_h = -1\text{KPa}$
Length of region, $L=1\text{m}$	depth of region, $D=2\text{m}$
Number of steps, $nsteps=100$	load increment, $df=-0.1\text{KN/m}$

The theoretical solution for the failure load, according to Mohr Coulomb failure criteria is given by,

$$\sigma_1 = \sigma_2 \left(\frac{1 + \sin \phi}{1 - \sin \phi} \right) - 2c \cdot \left(\frac{\cos \phi}{1 - \sin \phi} \right) \dots\dots\dots 8.1$$

$$\sigma_1 = -1 * \left(\frac{1 + \sin 30}{1 - \sin 30} \right) - 2 * 1 * \left(\frac{\cos 30}{1 - \sin 30} \right) = -6.464 \text{ KPa}$$

In the Matlab code the above problem is solved using different combinations of element types and numbers. The results are summarized in table 8.1

Element type	numx=3 and numy=4		numx=5 and numy=6	
	σ_{yy} (KPa)	u_y (mm)	σ_{yy} (KPa)	u_y (mm)
Q4, Q9, T6	-6.4	-2.7	-6.4	-2.4
Q8	-6.4	-2.8	-6.4	-2.8
T3	-6.4	-2.6	-6.4	-2.4

Table 8.1:- results from Matlab code

As can be seen the results are unchanged regardless of the number of elements involved for biaxial test. Plots of stresses and strains for this problem are shown below.

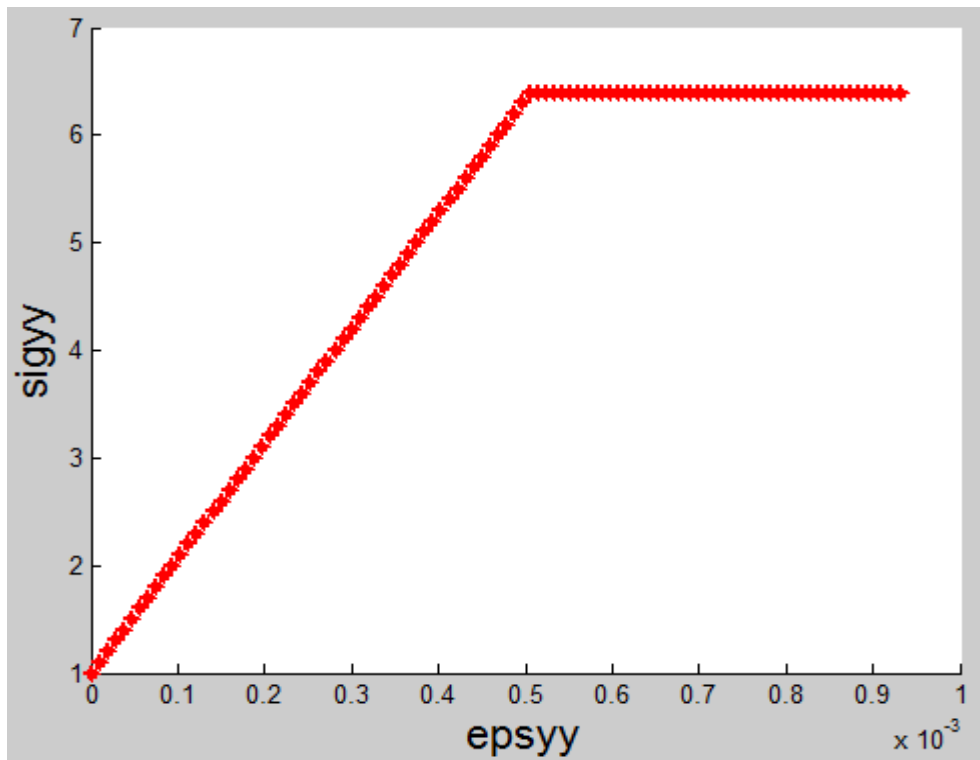


Figure 8.2:- vertical stress –strain curve

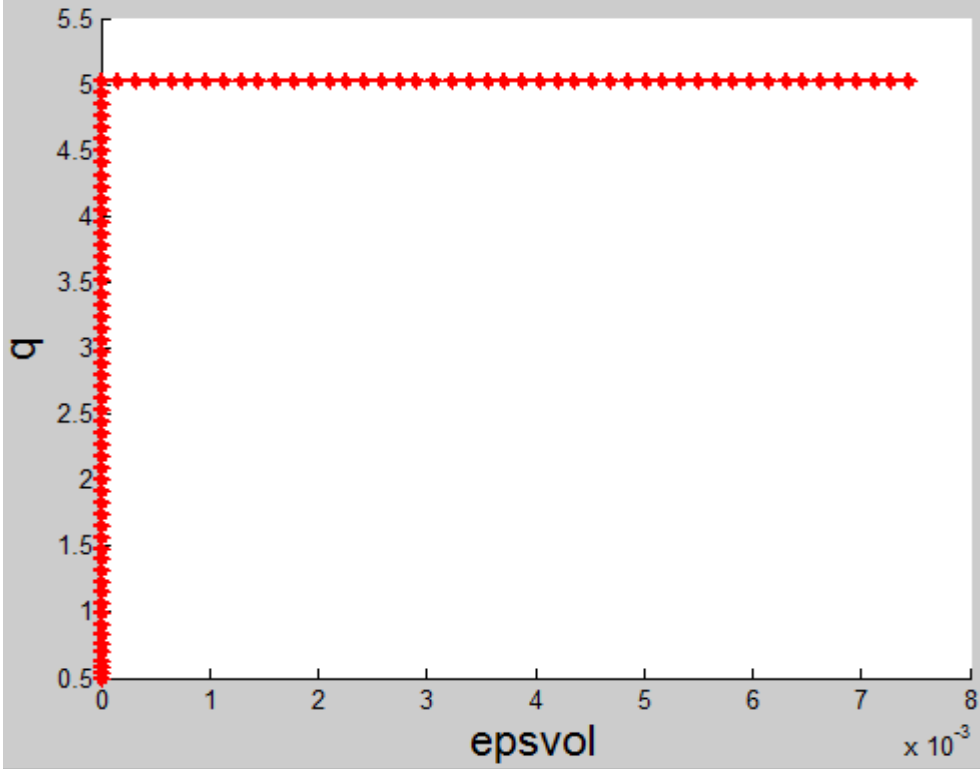


Figure 8.3:- deviatoric stress – volumetric strain curve

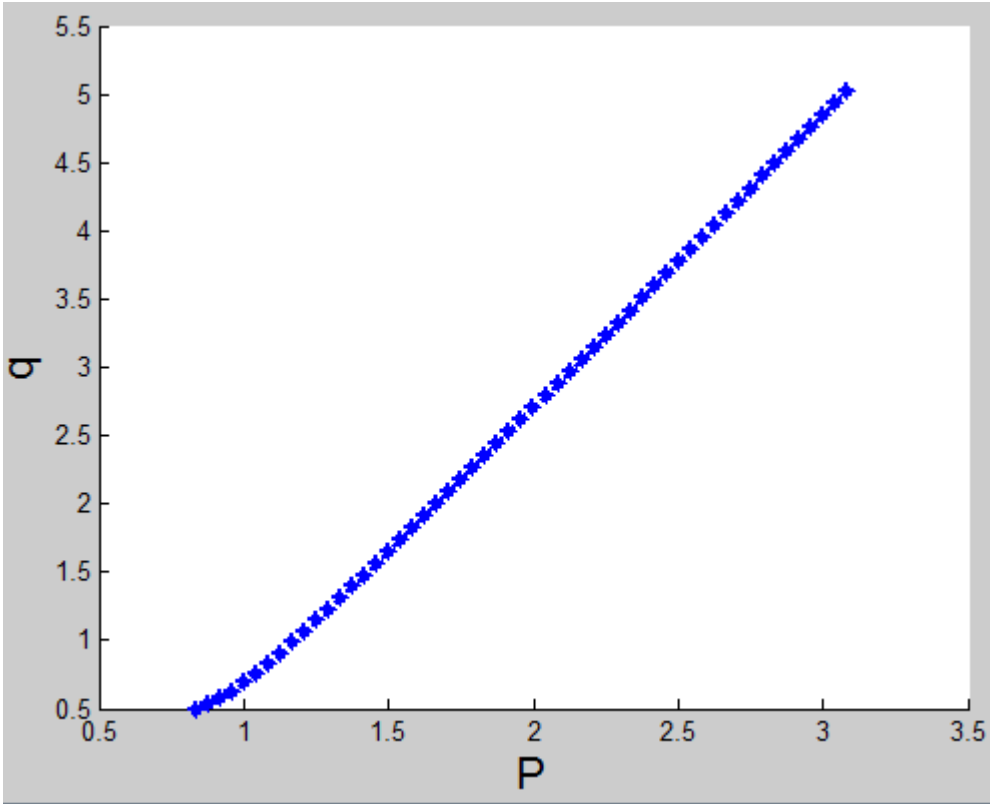


Figure 8.4:- deviatoric stress –mean stress curve

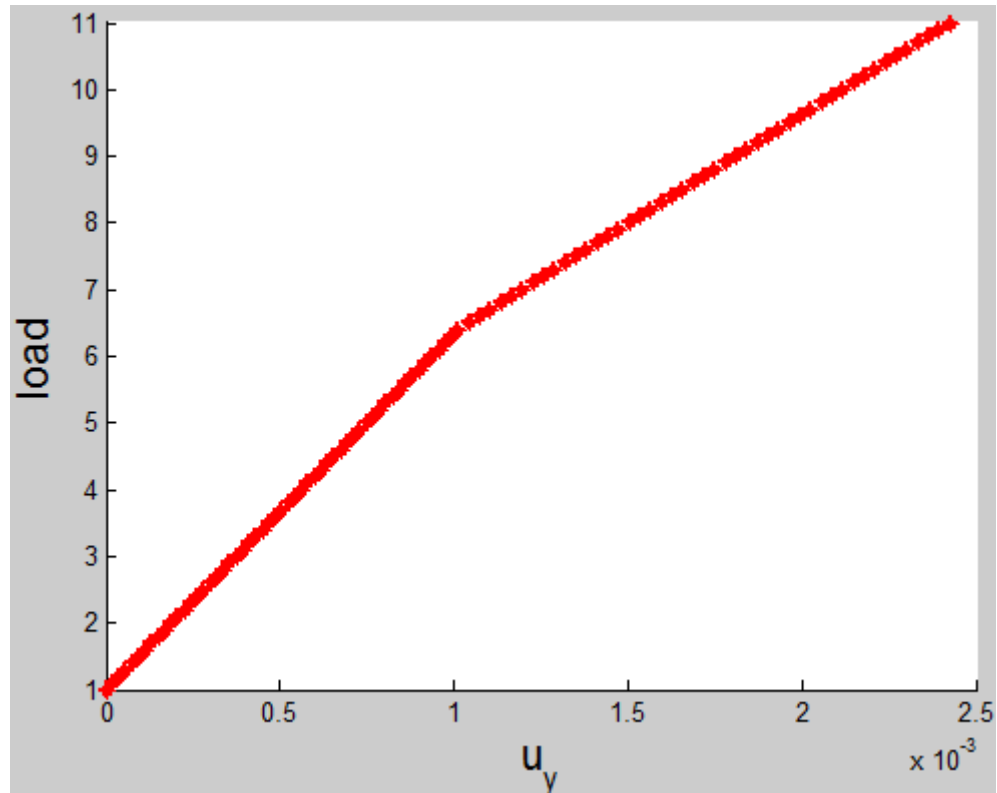


Figure 8.5:- load –displacement curve

Note that these plots the stresses and loads start from the initial confining stress value. A problem encountered here is the preparation of the load displacement curve as shown in figure 8.5. This is because there is no mechanism implemented that would reduce the load increments at the beginning of each load step when failure is approached. Hence on the load displacement curve presented here ‘load’ is just the summation of the load increments at the beginning of each step (before equilibrium). Solving this problem with the PLAXIS software gives a vertical yield stress of -6.464KPa which is close to the theoretical value calculated above and also with the Matlab code result.

8.3 Bearing capacity problem

The bearing capacity problems that can be analysed in the code are surface footings with a uniform vertical loading. A typical schematic diagram of a bearing capacity problem is shown below. For the time being only the bearing capacity of weightless soil is considered in the code but templates for initial stress generation with gravity loading and k_0 procedure are included in the code but not used here.

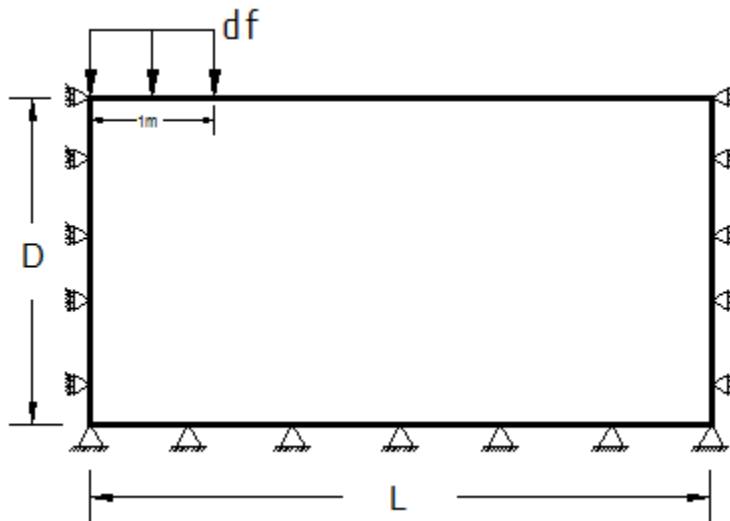


Figure 8.6:- schematic diagram for the bearing capacity problem

In this problem the following input parameters are used

Modulus of elasticity, $E=10,000\text{KPa}$	Poisson's ratio, $\nu=0.3,$
Angle of friction, $\phi=5^\circ$	Dilatancy angle, $\psi=5^\circ$
Cohesion, $c=1\text{KPa}$	unit weight, $\gamma=0$
load_edge1= 0	load_edge2= 1
Length of region, $L=5\text{m}$	Number of steps, nsteps=80
load increment, $df=-0.1\text{KN/m}$	

Note that the difference between load_edge1 and load_edge2 is the loaded area which is 1m wide as shown in figure 8.6. Also the maximum number of iteration is kept to be 20 since the load increment is small and hence the solution will converge after a few iterations. The order of integration (normal_order) used is 2 for Q4 and T3 elements and 3 for T6, Q8 and Q9 elements.

For a weightless soil, $\gamma = 0$, the ultimate bearing capacity of strip foundations has closed form solutions according to Prandtl,

$$q_u = \begin{cases} c \cdot N_c & \text{for } \phi \neq 0 \\ (\pi + 2) \cdot c & \text{for } \phi = 0 \end{cases} \dots\dots\dots 8.2$$

Where

N_c is a bearing capacity factor given by,

$$N_c = (N_q - 1) \cot \phi \dots\dots\dots 8.3$$

While,
$$N_q = e^{\pi \tan \phi} \left(\frac{1 + \sin \phi}{1 - \sin \phi} \right) \dots\dots\dots 8.4$$

For the soil parameters given above the bearing capacity of the soil using equation 8.2 will be

$$q_u = c \cdot N_c = 1 * 6.489 = 6.489 \text{ KPa}$$

This problem is first solved using the Matlab code for different domain depth (D) using a Q4 element. Keeping the number of element constant (numx=30 and numy=80) and varying the depth for a 5m long region gave the following results.

Depth (m)	Bearing capacity (KPa)
3	5.63
4	6.055
6	6.289

Table 8.2:- comparison of bearing capacity values for different depth of region

From these results it can be seen that for an accurate result the dimension of the region should be sufficiently large. Hence for a footing with a width B, the length of the region should be at least 5*B and the depth should be 4*B. for the proceeding calculations a region 5m long a 6m deep is used. Since the stress shadings and plots does not show these values with such precision, they can be seen in Matlab by typing `stress(:, :, ule)` in the command window. The term *ule* refers to the element at the left top corner and this element's gauss point is the stress point selected for plotting the curves.

For comparison of results for different element types the number of elements is set to be 1200 and the dimension of the region 5m x 6m (L x D). The results are shown in table 8.3.

Element type	numx	numy	Bearing capacity (KPa)
Q4	30	40	6.167
T3	15	40	5.943
T6	15	40	6.41
Q8	30	40	6.403
Q9	30	40	6.4023

Table 8.3:- comparison of bearing capacity values for different element types

As can be seen from these results, higher order elements (T6, Q8 and Q9 with normal_order =3) give a more accurate result than those of primary elements (T3 and Q4 with normal_order = 2). In the primary elements the number of elements can be increased beyond 1200, for example, for a Q4 element increasing the number of elements to 3200 (numx = 40 and numy = 80) gives a bearing capacity of 6.388KPa which is still less accurate than those of the higher order types with 1200 elements. For the case of higher order elements increasing the number of elements is possible but restricted by the computer memory allocated for MATLAB.

Solving this same problem in the PLAXIS software using T6 elements with fine mesh gives a bearing capacity value of 6.5113KPa. Hence the value from the Matlab code is close to both the closed form solution and the PLAXIS software. A problem encountered in the bearing capacity problem is the preparation of the load displacement curve as in the case of the biaxial test. This is because there is no mechanism implemented that would reduce the load increments at the beginning of each load step when failure is approached. Hence on the load displacement curve presented here 'load' is just the summation of the load increments at the beginning of each step (before equilibrium). Out puts for the Q9 element in table 8.2 are shown below.

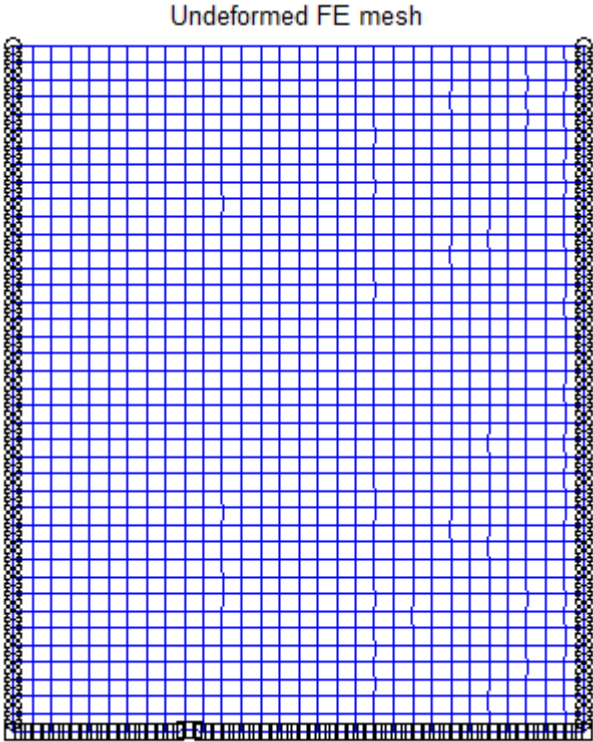


Figure 8.7:- undeformed mesh with support

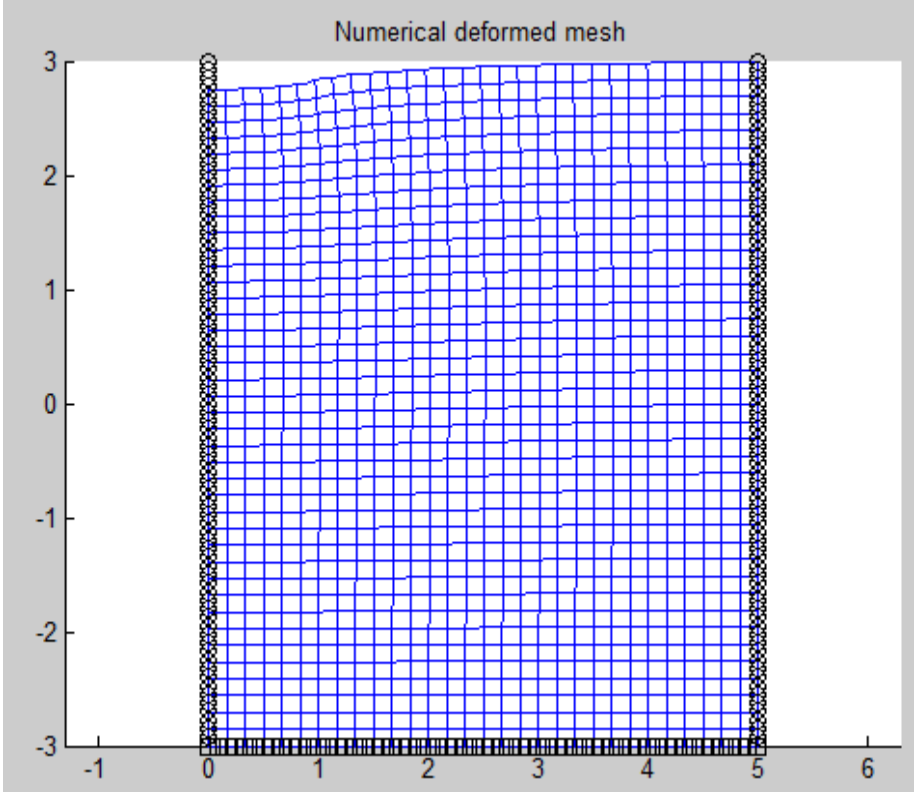


Figure 8.8:- deformed mesh with support

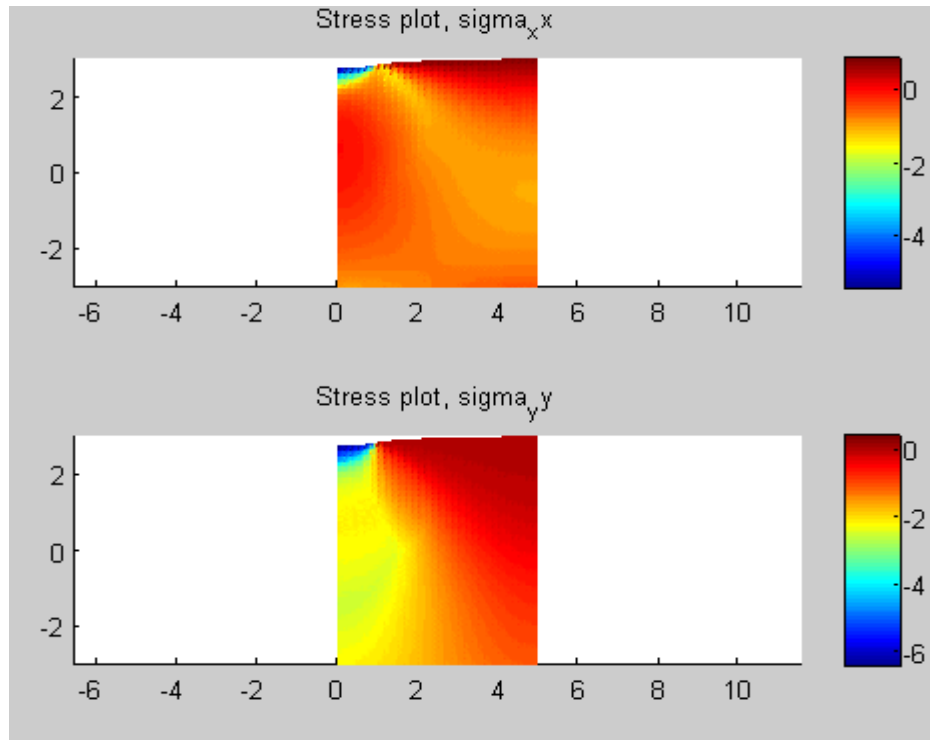


Figure 8.9:- stress shadings for σ_{xx} and σ_{yy}

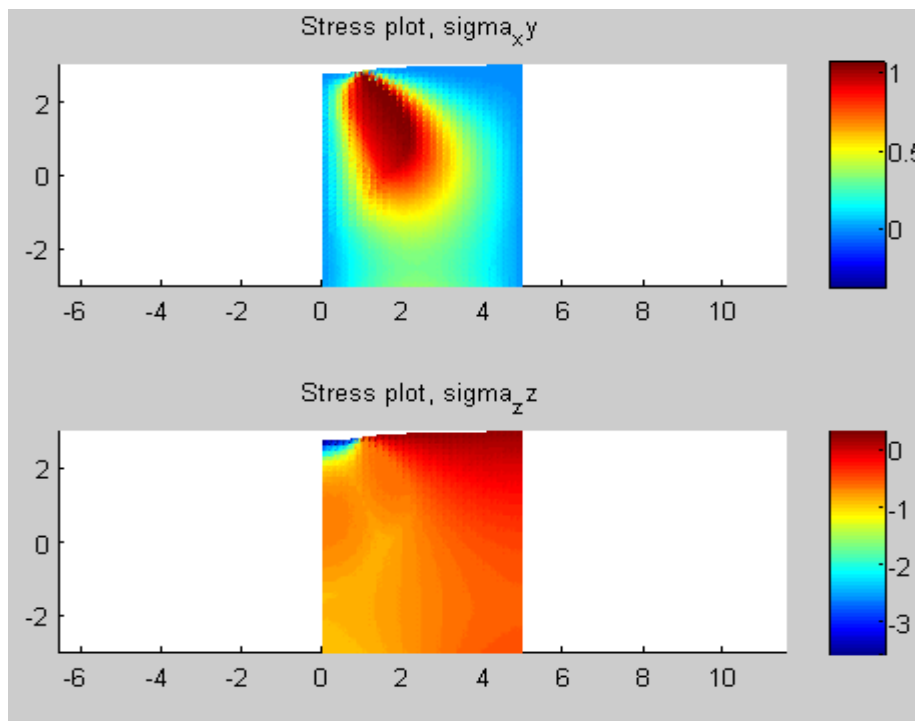


Figure 8.10:- stress shadings for σ_{xy} and σ_{zz}

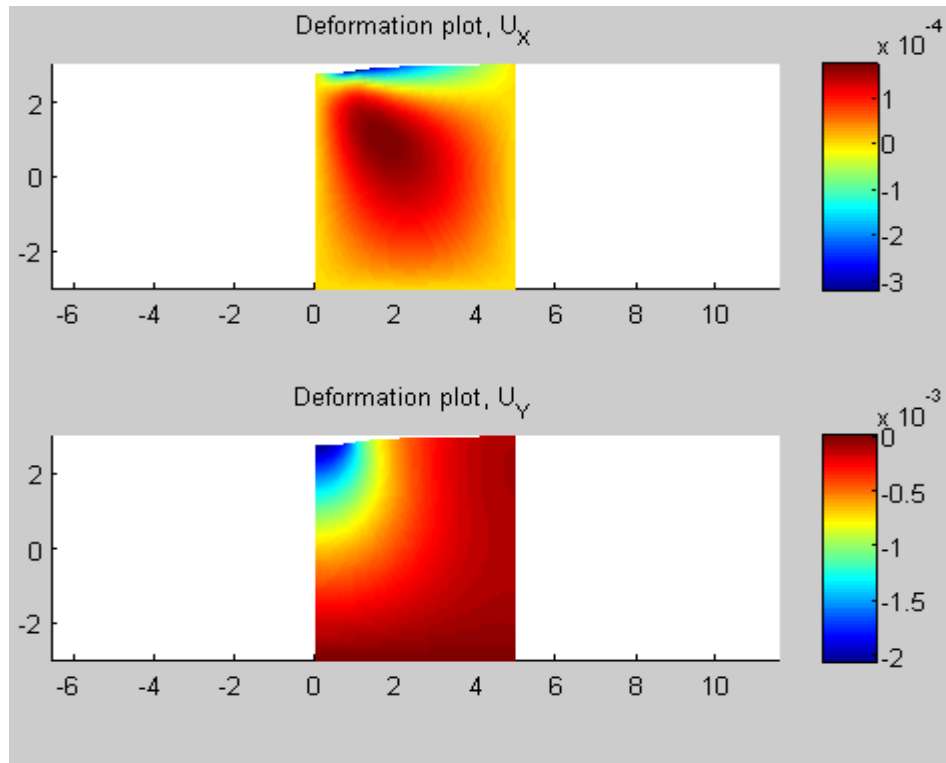


Figure 8.11:- displacement shadings

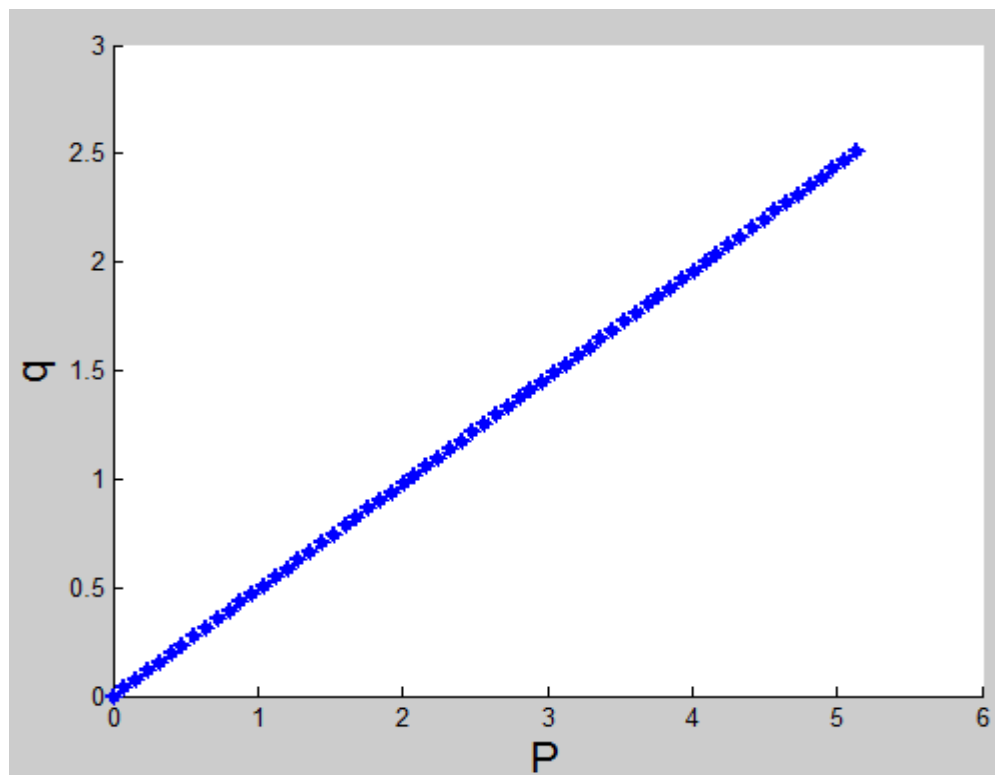


Figure 8.12:- deviatoric stress –mean stress curve

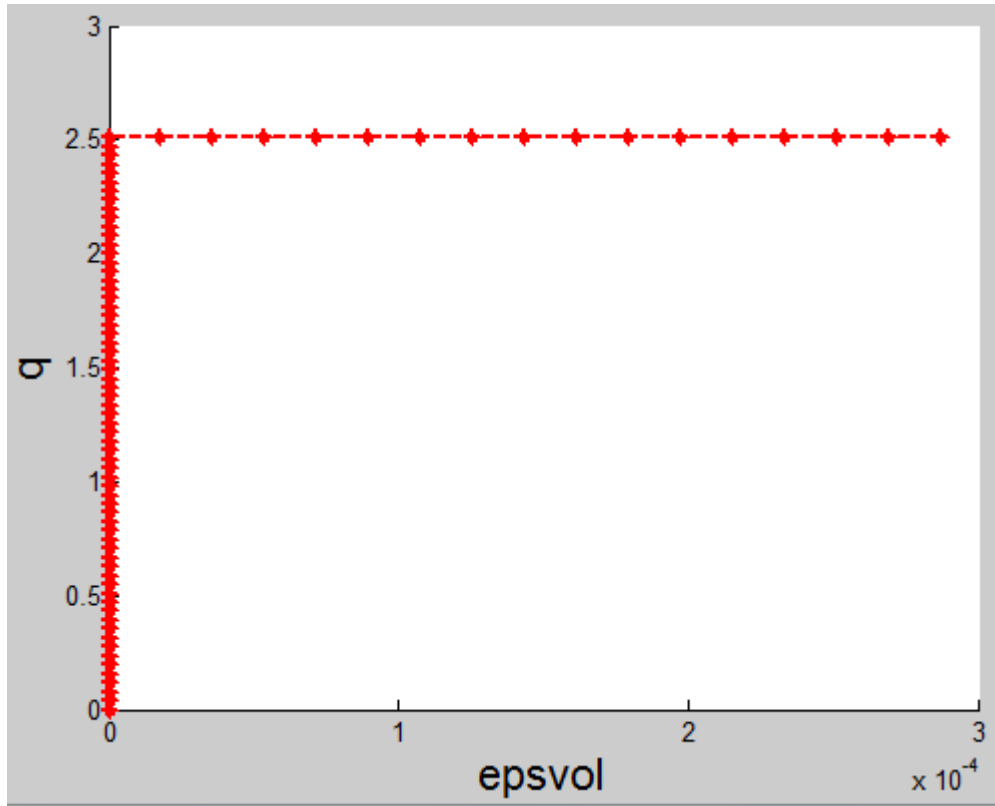


Figure 8.13:- deviatoric stress –volumetric strain curve

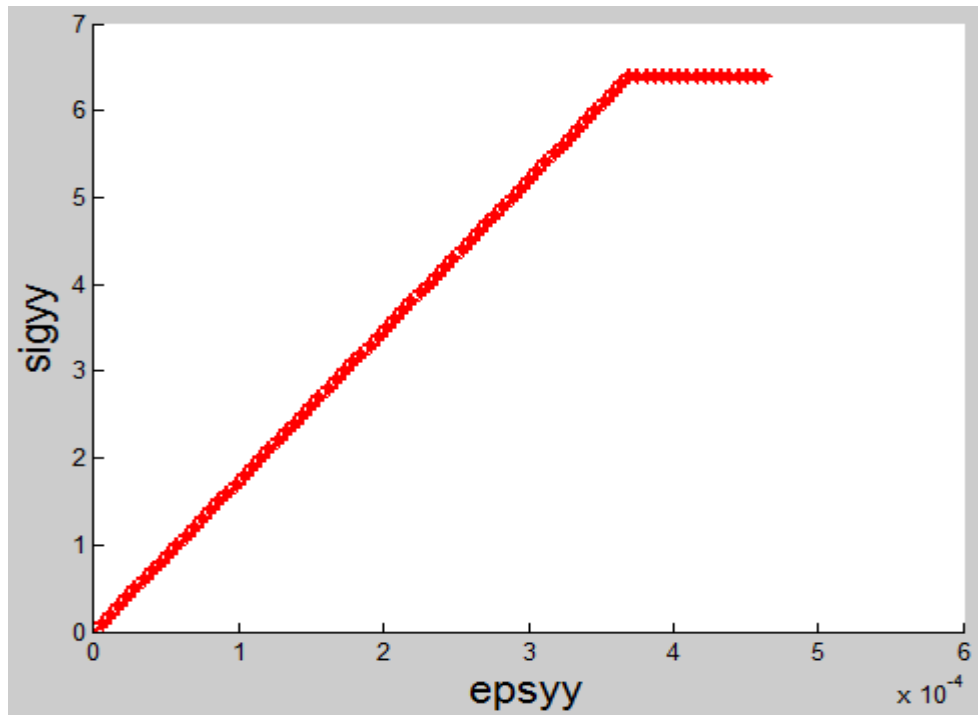


Figure 8.14:- vertical stress –vertical strain curve

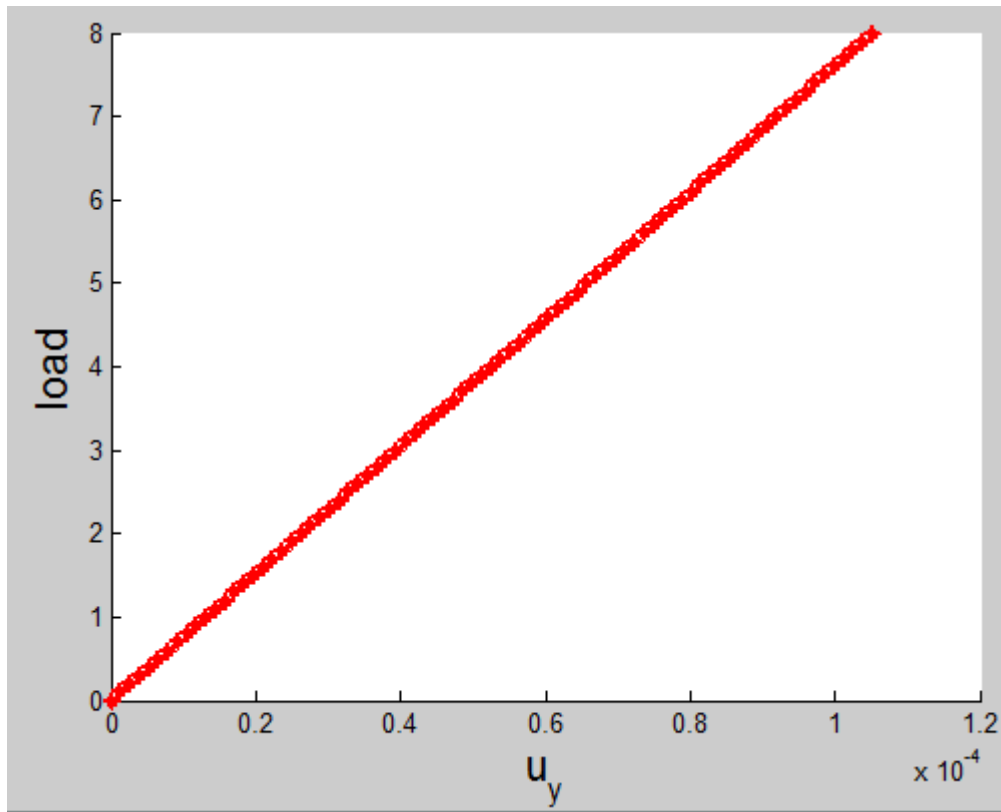


Figure 8.15:- 'load' –displacement curve

8.4 Conclusion

As can be seen from the above problems the accuracy of results is dependent on

- The type of element
- The number of elements and
- For bearing capacity problems the dimension of the domain

Elementary element types (Q4 and T3) with an order of integration of 2 give a less accurate result when compared to higher order element type (T6, Q8 and Q9) with an order of integration of 3 for a given number of elements. Increasing the number of elements in elementary element types makes the result more accurate, similarly increasing the number of elements has the same effect on higher order elements but the time it takes for analysis is too long and takes too much memory and usually results “out of memory” message from MATLAB due to the large size of the stiffness matrix. The integration order used for the above problems is only the minimum values but increasing it could make the results more accurate especially in higher order elements when great accuracy is intended. But in the present case it would not make any difference except slowing down the calculation.

In bearing capacity calculation, results are highly dependent on the dimensions of the domain compared to that of the loaded area, the larger the dimensions the more accurate is the result. Allocating a domain length, L , of five times the load span and a depth, D , of four times the load span gave a good result in the above calculations. Another issue is the number of steps and the size of load increment to be used, since both have to be inserted as an input. When the number of steps is under estimated the soil will not reach at its plastic state and the analysis will be elastic. Therefore it is advisable to make the number of steps large when the load increment is small. Due to the explicit nature of the integration used in the code, it is strongly recommended to use a small load increment. When solving a problem, it is advisable to start using the Q4 element with an integration order of 2 for a quick and rough calculation and to see if the number of steps is large enough, after which higher order elements with a larger integration order can be employed for a fine and accurate result.

CHAPTER 9

SUMMARY AND FUTURE WORK

9.1 Summary

From the calculations in the previous chapter, it is seen that higher order elements are more accurate than that of the elementary element types. A larger number of elements also improves the accuracy of results. One shortcoming in the code is there is no algorithm for the reduction of externally applied loads as failure is approached, which makes the load increment constant before and after the onset of plasticity. This problem made the plotting of the actual load displacement curve impossible. Instead the load displacement curve presented here is just the sum of the load increments at the beginning of each load step.

9.2 Future Work

In the future, this code can be expanded to include other geotechnical problems like

- Earth pressure calculations.
- Slope stability and
- Consolidation.

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Appendix A

Variable Nomenclature

Bfem:- the strain displacement matrix.

botEdge :- nodes along bottom edge of the domain in pairs

botNodes :- nodes along bottom edge of the domain

C: - The elastic compliance matrix.

coord:- matrix of the coordinates of the gauss points in the natural coordinate system.

CP: - The plastic compliance matrix.

dispNodes :- nodes which are restrained from displacement in any direction.

dispNodes1:- nodes which are restrained from displacement in x-direction.

dNdx_i:- matrix of the derivative of the shape function.

dof:- vector of the total degree of freedom.

E:-modulus of elasticity

eldof:- matrix containing the degree of freedom in an element

element :- a matrix that contains the node id in each element.

Elements: – a matrix that contains the node id in each element for Q8 or Q9 element before rearranged in the final manner.

elemType:- the type of element selected.

eta:- the natural y coordinate of the gauss points.

F: - the force matrix

f:- the externally applied force matrix.

F:- the total load (the sum of external load and self weight).

fac:- magnification factor for plotting outputs.

field:- a command for the output data to be plotted in color shading.

gamma :- self weight of soil.

iel:- a counter running from 1 up to the number of elements in the region.

inc_u:- the node counting increment in x direction (default=1)

inc_v:- the node counting increment in y direction

invJ0:- inverse of Jacobian matrix.

J0:- Jacobian matrix.

K: - the global stiffness matrix.

kk:- a counter running from 1 up to the number of gauss points in the element.

L, D: - Dimensions of the region in x and y direction respectively.

leftEdge :- nodes along left edge of the domain in pairs

leftNodes :- nodes along left edge of the domain

leftNodes1:- nodes along left edge of the domain except the lower left node.

lln :- lower left node

load_edge1:- the x-coordinate of the node at which the externally applied load starts.

load_edge2:- the x-coordinate of the node at which the externally applied load ends.

lrn :- lower right node

N:-shape function vector.

nn:- number of nodes in an element.

nnx:- number of nodes in the x direction.

nnny:- number of nodes in the y direction.

node :- matrix that contains the coordinate of the nodes.

node_pattern:- counting order of nodes arranged numerically.

Nodes :- matrix that contains the coordinate of the nodes for Q8 or Q9 element before rearranged in the final manner.

nonelem:- number of gauss points in an element.

normal_order :- number of gauss integration points (order of integration).

nu :- Poisson's ratio

numberelem:- total number of elements in the domain for a T3 or T6 element.

numelem:- number of elements in the region.

numnode:- total number of nodes in the region.

numx:- number of elements in the x direction.

numy:- number of elements in the y direction.

pt1, pt2, pt3, pt4:- coordinates of the corner points of the domain.

Q: - coordinates matrix of the gauss integration points.

qt:- type of integration to be done.

quadorder :-number of gauss integration points.

rightEdge :- nodes along right edge of the domain in pairs

rightNodes :- nodes along right edge of the domain

rightNodes1:- nodes along right edge of the domain except the lower right node.

sdim:- coordinate dimension of element for gauss integration either line element (value=1) of plane element (value=2).

se:- plotting color for the finite element mesh (default 'blue')

selfwt:- the self weight matrix.

sigmatox :- externally applied horizontal load.

sigmatoy :- externally applied vertical load.

Stresspoints:- nodes at which the stress is calculated.

topEdge :- nodes along the top edge of the domain in pairs

topEdge1:- nodes along the top edge of the domain in triples for Q9,Q8 and T6 elements.

topNodes :- nodes along the top edge of the domain

total_unknown:- the total degree of freedom in the domain.

type:-the type of element selected for shape function calculation.

U:- nodal displacements.

u_x:- nodal displacements in x-direction.

u_y:- nodal displacements in y-direction.

udof:- vector of the horizontal degree of freedom.

ule :- upper left element.

uln :- upper left node.

urn :- upper right node.

vdof:- vector of the vertical degree of freedom.

W :- vector of weight for the gauss integration.

xi:- the natural x coordinate of the gauss points.

xx:- a vector of x coordinates of nodes for a T6 element.

yy:- a vector of y coordinates of nodes for a T6 element.

Appendix B

Function Description

square_node_array:- Generates a quadrilateral array of nodes between the counterclockwise ordering of nodes pt1 - pt4.

Inputs: - pt1, pt2, pt3, pt4, nnx, nny.

Outputs: - node

q4totq8:- Forms the element and node matrices for Q8 element, from the element and node matrices of a Q4 element with the element and node matrices arranged in a counterclockwise order.

Inputs:-element, node, numx, numy.

Outputs: - Elements, Nodes

q4totq9:- Forms the element and node matrices for nine Q9 element, from the element and node matrices of a Q4 element with the element and node matrices arranged in a numerical counterclockwise order.

Inputs:-element, node, numx, numy.

Outputs: - Elements, Nodes

structured_q8_mesh:- Forms the element and node matrices for Q8 element, with the element and node matrices arranged in a nodal counterclockwise order (primary nodes then secondary nodes)

Inputs: - pt1, pt2, pt3, pt4, numx, numy.

Outputs: - node, element

structured_q9_mesh:- Forms the element and node matrices for Q9 element, with the element and node matrices arranged in a nodal counterclockwise order (primary nodes-> secondary nodes then internal nodes)

Inputs: - pt1, pt2, pt3, pt4, numx, numy.

Outputs: - node, element

make_elem:- creates a connectivity list of primary nodes in Q4 and T3 element.

Inputs: - node_pattern, numx, numy, inc_u, inc_v

Outputs: - element

mesh_region:- switching the element type selected, it forms the final nodal and element connectivity.using the above six functions for Q4, Q8, Q9 and T3 element.

Inputs: - pt1, pt2, pt3, pt4, numx, numy, elemType

Outputs: - node, element

mesh_t6_elem:- Forms the element and node matrices for T6 element, with the element and node matrices arranged in a nodal counterclockwise order (primary nodes then secondary nodes).

Inputs: - L, D, numx, numy

Outputs: - node, element

supportcond:- Forms a vector of nodes which are restrained in x direction or in both x and y direction which are the supports of the domain. It also forms a matrix of nodes for load application at the top of the domain.

Inputs: - elemType, numx, numy

Outputs: - topEdge, topEdge1, dispNodes, dispNodes1, leftNodes1

elementdof:- Forms the global degree of freedom nodes in each element from the node identification number.

Inputs: - elemType, sctr

Outputs: - eldof

gauss_pt_wt:- Returns the weights and coordinates of the gauss integration points.

Inputs: - quadorder, qt, sdim

Outputs: - W, Q

gauss_rule:- Provides the weight vector and gauss point coordinate matrix of the element based on the integration order selected.

Inputs: - iel, elemType, normal_order

Outputs: - W, Q

shape_func:- Gives the shape function and its derivatives with respect to x and y.

Inputs: - type, coord, dim

Outputs: - N, dNdx

Bmatrix:- Gives the strain displacement matrix (B matrix size 3x8) for each element.

Inputs: - pt, elemType, iel

Outputs: - Bfem

Bmatrix4:- Gives the strain displacement matrix (B matrix size 4x8) for each element.

Inputs: - pt, elemType, iel

Outputs: - Bfem4

stiffness_matrix:- Generates the element and global stiffness matrix.

Inputs: - node, element, elemType, normal_order, C

Outputs: - K

force_matrix:- Generates the force matrix due to externally applied loads for Q4 and T3 elements and the position of the vertically applied load (sctry) in the global degree of freedom.

Inputs: - node, topEdge, sigmatoy, sigmatox, load_edge1, load_edge2, D

Outputs: - f, sctry

force_matrix689:- Generates the force matrix due to externally applied loads for Q9, Q8 and T6 elements and the position of the vertically applied load (sctry) in the global degree of freedom.

Inputs: - node, topEdge, topEdge1, sigmatoy, sigmatox, load_edge1, load_edge2, D

Outputs: - f, sctry

selfwt_matrix:- Generates the force matrix due to self weight.

Inputs: - elemType, normal_order, gamma, node, element

Outputs: - selfwt

displacements:- evaluates the unknown degree of freedom (displacements) at the nodes.

Inputs: - dispNodes, dispNodes1, numnode, K, f, selfwt

Outputs: - U, u_x, u_y

stress_calculation:- calculates the element strains and stresses at the nodes in x, y, xy and zz directions.

Inputs: - node, element, elemType, U, normal_order, C

Outputs: - stress, strain

invariants:- calculates the stress invariants for a single gauss point.

Inputs: - stress

Outputs: - p, q, theta

invariants1:- calculates the stress invariants at each gauss point. This can be used outside the iteration loop.

Inputs: - element, stress

Outputs: - p, q, theta

invariants2:- calculates the stress invariants at each gauss point. This can be used inside the iteration loop.

Inputs: - kk, iel, stress

Outputs: - p, q, theta

formdg:- calculates the partial derivatives of the plastic potential with respect to p , J_2 and J_3 .
Inputs: - tsi, q, theta
Outputs: - dg1, dg2, dg3

formdm:- Generates the partial derivatives of the p , J_2 and J_3 with respect to stress for use in the plastic potential derivative.
Inputs: - kk, iel, stress
Outputs: - m1, m2, m3

internalrxn:- Generates the vector of nodal force reactions due to internal stress. It can be used outside the iteration loop if the stress at each gauss point is given.
Inputs: - node, element, elemType, normal_order, stress
Outputs: - r

principal_stress:- calculates the principal stresses from the stress invariants.
Inputs: - kk, iel, p, q, theta
Outputs: - s1, s2, s3

plastic_mat:- Generates the plastic constitutive matrix CP at each gauss point.
Inputs: - E, nu, phi, tsi, iel, kk, stress
Outputs: - CP

makedgds:- Forms the partial derivatives of the plastic potential function with respect to stress at each gauss point.
Inputs: - tsi, element, iel, kk, stress
Outputs: - dgds

makedfds:- Forms the partial derivatives of the yield function with respect to stress at each gauss point.
Inputs: - phi, element, iel, kk, stress
Outputs: - dfds

plot_mesh:- Plots the finite element mesh
Inputs: - node, connect, elem_type, se
Outputs: - plot

plot_m:- Plots the finite element mesh with the support condition.
Inputs: - elemType, dispNodes, dispNodes1
Outputs: - plot

plot_def:- Plots the deformed finite element mesh with the support condition.
Inputs: - fac, u_x, u_y, elemType, dispNodes, dispNodes1
Outputs: - plot

plot_field:- Forms a color dependent finite element mesh for plotting outputs

Inputs: - node, connect, elem_type, field

Outputs: - plot

plot_sig:- plots the color coded stress distribution in the finite element region along with color bar scale.

Inputs: - fac, u_x, u_y, elemType, stress

Outputs: - plot

plot_defo:- plots the color coded displacement intensity in the finite element region along with color bar scale.

Inputs: - fac, u_x, u_y, elemType

Outputs: - plot

Appendix C

The MATLAB code

make_elem

```
function element=make_elem(node_pattern,numx,numy,inc_u,inc_v)

% creates a connectivity list of primary nodes in Q4 and T3 element
if ( nargin < 5 )
    disp('Not enough parameters specified for make_elem function')
end
inc=[zeros(1,size(node_pattern,2))];
e=1;
element=zeros(numx*numy,size(node_pattern,2));

for row=1:numy
    for col=1:numx
        element(e,:)=node_pattern+inc;
        inc=inc+inc_u;
        e=e+1;
    end
    inc=row*inc_v;
end
end % end of function
```

mesh_t6_elem

```
function [node,element] =mesh_t6_elem(L,D,numx,numy)

%Forms the element and node matrices for T6 element, with the element
%and node matrices arranged in a nodal counterclockwise order
%(primary nodes then secondary nodes).

xx= repmat((0:L/(2*numx):L)',2*numy+1,1);
yy=sort(repmat((-1*D/2:D/(2*numy):D/2)',2*numx+1,1));
node=[xx yy];
inc_u=[2 2 2 2 2 2];
inc_v=[2+(4*numx) 2+(4*numx) 2+(4*numx) 2+(4*numx) 2+(4*numx) 2+(4*numx)];
node_pattern1=[1 2 3 2*numx+3 4*numx+3 2*numx+2];
node_pattern2=[3 2*numx+4 4*numx+5 4*numx+4 4*numx+3 2*numx+3];
elements1=make_elem(node_pattern1,numx,numy,inc_u,inc_v);
elements2=make_elem(node_pattern2,numx,numy,inc_u,inc_v);

numberelem=2*numx*numy;
elements=zeros(numberelem,6);
elements(1:2:end,:)=elements1;
elements(2:2:end,:)=elements2;
    element(:,1)=elements(:,1);
    element(:,2)=elements(:,3);
    element(:,3)=elements(:,5);
    element(:,4)=elements(:,2);
    element(:,5)=elements(:,4);
    element(:,6)=elements(:,6);
end % end of function
```

mesh_region

```
function [node,element]= mesh_region(pt1, pt2, pt3, pt4,numx,numy,elemType)

    % switching the element type selected, it forms the final nodal and
    % element connectivity.using the above six functions for Q4,Q8,Q9 and
    % T3 element.

global L D
nnx = numx+1;
nny = numy+1;

switch elemType

    case 'Q4'
        node=square_node_array(pt1,pt2,pt3,pt4,nnx,nny);
        inc_u=1;
        inc_v=nnx;
        node_pattern=[ 1 2 nnx+2 nnx+1 ];
        [element]=make_elem(node_pattern,numx,numy,inc_u,inc_v);

    case 'Q9'
        [node,element]=structured_q9_mesh(pt1,pt2,pt3,pt4,numx,numy);

    case 'Q8'
        [node,element]=structured_q8_mesh(pt1,pt2,pt3,pt4,numx,numy);

    case 'T3'

        node=square_node_array(pt1,pt2,pt3,pt4,nnx,nny);
        node_pattern1=[ 1 2 nnx+1 ];
        node_pattern2=[ 2 nnx+2 nnx+1 ];
        inc_u=1;
        inc_v=nnx;
        numberelem=2*numx*numy;
        element=zeros(numberelem,3);
        element1=make_elem(node_pattern1,numx,numy,inc_u,inc_v);
        element2=make_elem(node_pattern2,numx,numy,inc_u,inc_v);
        element(1:2:end,:)=element1;
        element(2:2:end,:)=element2;

    otherwise
        error('only Q4,Q9,Q8,and T3 are supported by this function');
end

end % end of function
```

q4totq8

```
function [Elements,Nodes]=q4totq8(element,node,numx,numy)

% forms the element and node matrices for eight node rectangular element,
% from the element and node matrices of a four node rectangular element
% with the element and node matrices arranged in a counterclockwise order
% Inputs:-element,node,numx and numy.

nnx=numx+1;
num_u=numx;
num_v=numy;
inc_u=[2 2 2 1 2 2 2 1];

if numx==2
    inc_v=[8 8 8 8 8 8 8 8];

else
    inc_v=[8+3*(numx-2) 8+3*(numx-2) 8+3*(numx-2) 8+3*(numx-2) ...
          8+3*(numx-2) 8+3*(numx-2) 8+3*(numx-2) 8+3*(numx-2)];
end

node_pattern=[1 2 3 2*nnx+1 3*nnx+2 3*nnx+1 3*nnx 2*nnx];
inc=[zeros(1,size(node_pattern,2))];
e=1;
elements=zeros(num_u*num_v,size(node_pattern,2));

for row=1:num_v
    for col=1:num_u
        elements(e,:)=node_pattern+inc;
        inc=inc+inc_u;
        e=e+1;
    end
    inc=row*inc_v;
end

Elements=elements;
numNode=size(unique(Elements),1);
Nodes=zeros(numNode,2);
numElement=numx*numy;

for i=1:numElement

    Nodes(Elements(i,1),1)=node(element(i,1),1);
    Nodes(Elements(i,1),2)=node(element(i,1),2);

    Nodes(Elements(i,2),1)=(node(element(i,1),1)+node(element(i,2),1))/2;
    Nodes(Elements(i,2),2)=node(element(i,1),2);

    Nodes(Elements(i,3),1)=node(element(i,2),1);
    Nodes(Elements(i,3),2)=node(element(i,2),2);

    Nodes(Elements(i,4),1)=node(element(i,2),1);
    Nodes(Elements(i,4),2)=(node(element(i,2),2)+node(element(i,3),2))/2;
```



```

Nodes(Elements(i,5),1)=node(element(i,3),1);
Nodes(Elements(i,5),2)=node(element(i,3),2);

Nodes(Elements(i,6),1)=(node(element(i,3),1)+node(element(i,4),1))/2;
Nodes(Elements(i,6),2)=node(element(i,3),2);

Nodes(Elements(i,7),1)=node(element(i,4),1);
Nodes(Elements(i,7),2)=node(element(i,4),2);

Nodes(Elements(i,8),1)=node(element(i,4),1);
Nodes(Elements(i,8),2)=(node(element(i,4),2)+node(element(i,1),2))/2;

end

Nodes=Nodes;
Elements(:,1)=elements(:,1);
Elements(:,2)=elements(:,3);
Elements(:,3)=elements(:,5);
Elements(:,4)=elements(:,7);
Elements(:,5)=elements(:,2);
Elements(:,6)=elements(:,4);
Elements(:,7)=elements(:,6);
Elements(:,8)=elements(:,8);

end % end of function

```

q4totq9

```
function [Elements,Nodes]=q4totq9(element,node,numx,numy)

% forms the element and node matrices for nine node rectangular element
% given the element and node matrices of a four node rectangular element
% and the number of elements in x and y direction

nnx=numx+1;
num_u=numx;
num_v=numy;
inc_u=[2 2 2 2 2 2 2 2 2];

if numx==3
    inc_v=[14 14 14 14 14 14 14 14 14];
else
    inc_v=[14+4*(numx-3) 14+4*(numx-3) 14+4*(numx-3) 14+4*(numx-3)...
    14+4*(numx-3) 14+4*(numx-3) 14+4*(numx-3) 14+4*(numx-3) 14+4*(numx-3)];
end

node_pattern=[ 1 2 3 2*nnx+2 4*nnx+1 4*nnx 4*nnx-1 2*nnx 2*nnx+1];
inc=zeros(1,size(node_pattern,2));
e=1;
elements=zeros(num_u*num_v,size(node_pattern,2));

for row=1:num_v
    for col=1:num_u
        elements(e,:)=node_pattern+inc;
        inc=inc+inc_u;
        e=e+1;
    end
    inc=row*inc_v;
end

Elements=elements;
numNode=size(unique(Elements),1);
Nodes=zeros(numNode,2);
numElement=numx*numy;

for i=1:numElement

    Nodes(Elements(i,1),1)=node(element(i,1),1);
    Nodes(Elements(i,1),2)=node(element(i,1),2);

    Nodes(Elements(i,2),1)=(node(element(i,1),1)+node(element(i,2),1))/2;
    Nodes(Elements(i,2),2)=node(element(i,1),2);

    Nodes(Elements(i,3),1)=node(element(i,2),1);
    Nodes(Elements(i,3),2)=node(element(i,2),2);

    Nodes(Elements(i,4),1)=node(element(i,2),1);
    Nodes(Elements(i,4),2)=(node(element(i,2),2)+node(element(i,3),2))/2;
```

```

Nodes(Elements(i,5),1)=node(element(i,3),1);
Nodes(Elements(i,5),2)=node(element(i,3),2);

Nodes(Elements(i,6),1)=(node(element(i,3),1)+node(element(i,4),1))/2;
Nodes(Elements(i,6),2)=node(element(i,3),2);

Nodes(Elements(i,7),1)=node(element(i,4),1);
Nodes(Elements(i,7),2)=node(element(i,4),2);

Nodes(Elements(i,8),1)=node(element(i,4),1);
Nodes(Elements(i,8),2)=(node(element(i,4),2)+node(element(i,1),2))/2;

Nodes(Elements(i,9),1)=(node(element(i,1),1)+node(element(i,2),1))/2;
Nodes(Elements(i,9),2)=(node(element(i,4),2)+node(element(i,1),2))/2;
end
Nodes=Nodes;
Elements(:,1)=elements(:,1);
Elements(:,2)=elements(:,3);
Elements(:,3)=elements(:,5);
Elements(:,4)=elements(:,7);
Elements(:,5)=elements(:,2);
Elements(:,6)=elements(:,4);
Elements(:,7)=elements(:,6);
Elements(:,8)=elements(:,8);
Elements(:,9)=elements(:,9);

end % end of function

```

square_node_array

```
function node=square_node_array(pt1,pt2,pt3,pt4,nnx,ny)

% Generates a quadrilateral array of nodes between the counterclockwise
% ordering of nodes pt1 - pt4,given number of elements in x and y direction

if ( nargin < 6 )
    disp('Not enough parameters specified for square_node_array function')
end

% get node spacing along u direction
xi_pts=linspace(-1,1,nnx);

% get node spacing along v direction
eta_pts=linspace(-1,1,ny);

x_pts=[pt1(1),pt2(1),pt3(1),pt4(1)];
y_pts=[pt1(2),pt2(2),pt3(2),pt4(2)];

for r=1:ny
    eta=eta_pts(r);

    for c=1:nnx
        xi=xi_pts(c);
        % get interpolation basis at xi, eta
        N=shape_func('Q4',[xi,eta]);
        N=N(:,1);
        node((r-1)*nnx+c,:)= [x_pts*N,y_pts*N];
    end
end

end % end of function
```

structured_q8_mesh

```
function [node,element]=structured_q8_mesh(pt1,pt2,pt3,pt4,numx,numy)

%Forms the element and node matrices for Q8 element,
%with the element and node matrices arranged in a nodal
%counterclockwise order (primary nodes then secondary nodes)

nnx=numx+1;
nny=numy+1;
node=square_node_array(pt1,pt2,pt3,pt4,nnx,nny);

inc_u=1;
inc_v=nnx;
node_pattern=[ 1 2 nnx+2 nnx+1 ];

element=make_elem(node_pattern,numx,numy,inc_u,inc_v);

[element,node]=q4totq8(element,node,numx,numy);

end % end of function
```

structured_q9_mesh

```
function [node,element]=structured_q9_mesh(pt1,pt2,pt3,pt4,numx,numy)

%Forms the element and node matrices for Q9 element, with the element
%and node matrices arranged in a nodal counterclockwise order
%(primary nodes-> secondary nodes then internal nodes)

nnx=numx+1;
nny=numy+1;
node=square_node_array(pt1,pt2,pt3,pt4,nnx,nny);

inc_u=1;
inc_v=nnx;
node_pattern=[ 1 2 nnx+2 nnx+1 ];

element=make_elem(node_pattern,numx,numy,inc_u,inc_v);

[element,node]=q4totq9(element,node,numx,numy);

end % end of function
```

Bmatrix

```
function Bfem =Bmatrix(pt,elemType,iel)

% Gives the strain displacement matrix (B matrix of size 3x8)of each element

global node element

sctr = element(iel,:);
nn    = length(sctr);
[N,dNdx] = shape_func(elemType,pt); % element shape functions
J0 = node(sctr,:)'*dNdx;           % element Jacobian matrix
invJ0 = inv(J0);
dNdx  = dNdx*invJ0;               % derivatives of N w.r.t XY
Gpt = N'*node(sctr,:);           % GP in global coord, used

    Bfem = zeros(3,2*nn);
    Bfem(1,1:2:2*nn) = dNdx(:,1)';
    Bfem(2,2:2:2*nn) = dNdx(:,2)';
    Bfem(3,1:2:2*nn) = dNdx(:,2)';
    Bfem(3,2:2:2*nn) = dNdx(:,1)';

end % end of function
```

Bmatrix4

```
function Bfem4 =Bmatrix4(pt,elemType,iel)

% Gives the strain displacement matrix (B matrix of size 4x8)of each element

global node element

sctr = element(iel,:);
nn    = length(sctr);
[N,dNdx] = shape_func(elemType,pt); % element shape functions
J0 = node(sctr,:)'*dNdx;           % element Jacobian matrix
invJ0 = inv(J0);
dNdx  = dNdx*invJ0;               % derivatives of N w.r.t XY
Gpt = N'*node(sctr,1);           % GP in global coord, used

    Bfem4 = zeros(4,2*nn);
    Bfem4(1,1:2:2*nn) = dNdx(:,1)';
    Bfem4(2,2:2:2*nn) = dNdx(:,2)';
    Bfem4(3,1:2:2*nn) = dNdx(:,2)';
    Bfem4(3,2:2:2*nn) = dNdx(:,1)';

end % end of function
```

gauss_rule

```
function [W,Q] = gauss_rule(iel,elemType,normal_order)

% Provides the weight vector and gauss point coordinate matrix of
% the element based on the integration order selected.

global node element

sctr = element(iel,:);           % element connectivity

    if ((elemType == 'Q4') & (normal_order < 8))
        [W,Q] = gauss_pt_wt(normal_order,'GAUSS',2);
    elseif ((elemType == 'Q8') & (normal_order < 8))
        [W,Q] = gauss_pt_wt(normal_order,'GAUSS',2);
    elseif ((elemType == 'Q9') & (normal_order < 8))
        [W,Q] = gauss_pt_wt(normal_order,'GAUSS',2);
    elseif elemType == 'T3'
        [W,Q] = gauss_pt_wt(normal_order,'TRIANGULAR',2);
    elseif elemType == 'T6'
        [W,Q] = gauss_pt_wt(normal_order,'TRIANGULAR',2);
    end

end % end of function
```

gauss_pt_wt

```
function [W,Q] = gauss_pt_wt(quadorder, qt, sdim )

% Returns the weights and coordinates of the gauss integration points

if ( nargin < 3 )    % set default arguments
    if ( strcmp(qt,'GAUSS') == 1 )
        dim = 1;
    else
        dim = 2;
    end
end

if ( nargin < 2 )
    type = 'GAUSS';
end

if ( strcmp(qt,'GAUSS') == 1 )
    quadpoint=zeros(quadorder^sdim ,sdim);
    quadweight=zeros(quadorder^sdim,1);
    rlpt=zeros(quadorder,1); rlwt=zeros(quadorder,1);
end
```

```

switch ( quadorder )
  case 1
    rlpt(1) = 0.0000000000000000;
    rlwt(1) = 2.0000000000000000;

  case 2
    rlpt(1) = 0.577350269189626;
    rlpt(2) = -0.577350269189626;

    rlwt(1) = 1.0000000000000000;
    rlwt(2) = 1.0000000000000000;

  case 3
    rlpt(1) = 0.774596669241483;
    rlpt(2) = -0.774596669241483;
    rlpt(3) = 0.0000000000000000;

    rlwt(1) = 0.5555555555555556;
    rlwt(2) = 0.5555555555555556;
    rlwt(3) = 0.8888888888888889;

  case 4
    rlpt(1) = 0.861134311594053;
    rlpt(2) = -0.861134311594053;
    rlpt(3) = 0.339981043584856;
    rlpt(4) = -0.339981043584856;

    rlwt(1) = 0.347854845137454;
    rlwt(2) = 0.347854845137454;
    rlwt(3) = 0.652145154862546;
    rlwt(4) = 0.652145154862546;

  case 5
    rlpt(1) = 0.906179845938664;
    rlpt(2) = -0.906179845938664;
    rlpt(3) = 0.538469310105683;
    rlpt(4) = -0.538469310105683;
    rlpt(5) = 0.0000000000000000;

    rlwt(1) = 0.236926885056189;
    rlwt(2) = 0.236926885056189;
    rlwt(3) = 0.478628670499366;
    rlwt(4) = 0.478628670499366;
    rlwt(5) = 0.5688888888888889;

  case 6
    rlpt(1) = 0.932469514203152;
    rlpt(2) = -0.932469514203152;
    rlpt(3) = 0.661209386466265;
    rlpt(4) = -0.661209386466265;
    rlpt(5) = 0.238619186003152;
    rlpt(6) = -0.238619186003152;

    rlwt(1) = 0.171324492379170;
    rlwt(2) = 0.171324492379170;
    rlwt(3) = 0.360761573048139;

```



```

    rlwt(4) = 0.360761573048139;
    rlwt(5) = 0.467913934572691;
    rlwt(6) = 0.467913934572691;

case 7
    rlpt(1) = 0.949107912342759;
    rlpt(2) = -0.949107912342759;
    rlpt(3) = 0.741531185599394;
    rlpt(4) = -0.741531185599394;
    rlpt(5) = 0.405845151377397;
    rlpt(6) = -0.405845151377397;
    rlpt(7) = 0.000000000000000;

    rlwt(1) = 0.129484966168870;
    rlwt(2) = 0.129484966168870;
    rlwt(3) = 0.279705391489277;
    rlwt(4) = 0.279705391489277;
    rlwt(5) = 0.381830050505119;
    rlwt(6) = 0.381830050505119;
    rlwt(7) = 0.417959183673469;

otherwise
    disp('unsupported integration order')

end % end of quadorder switch

n=1;

if ( sdim == 1 )
    for i = 1:quadorder
        quadpoint(n,:) = [ rlpt(i) ];
        quadweight(n) = rlwt(i);
        n = n+1;
    end

elseif ( sdim == 2 )
    for i = 1:quadorder
        for j = 1:quadorder
            quadpoint(n,:) = [ rlpt(i), rlpt(j)];
            quadweight(n) = rlwt(i)*rlwt(j);
            n = n+1;
        end
    end

end

Q=quadpoint;
W=quadweight;

% END OF GAUSSIAN QUADRATURE DEFINITION FOR RECTANGULAR ELEMENTS

```

```

elseif ( strcmp(qt,'TRIANGULAR') == 1 )

    if ( quadorder > 7 ) % check for valid quadrature order
        disp('Quadrature order too high for triangular quadrature');
        quadorder = 1;
    end

    if ( quadorder == 1 ) % set quad points and quadweights
        quadpoint = [ 0.33333333333333, 0.33333333333333 ];
        quadweight = 1;

    elseif ( quadorder == 2 )
        quadpoint = zeros( 3, 2 );
        quadweight = zeros( 3, 1 );

        quadpoint(1,:) = [ 0.16666666666667, 0.16666666666667 ];
        quadpoint(2,:) = [ 0.66666666666667, 0.16666666666667 ];
        quadpoint(3,:) = [ 0.16666666666667, 0.66666666666667 ];

        quadweight(1) = 0.33333333333333;
        quadweight(2) = 0.33333333333333;
        quadweight(3) = 0.33333333333333;

    elseif ( quadorder <= 5 )
        quadpoint = zeros( 7, 2 );
        quadweight = zeros( 7, 1 );

        quadpoint(1,:) = [ 0.1012865073235, 0.1012865073235 ];
        quadpoint(2,:) = [ 0.7974269853531, 0.1012865073235 ];
        quadpoint(3,:) = [ 0.1012865073235, 0.7974269853531 ];
        quadpoint(4,:) = [ 0.4701420641051, 0.0597158717898 ];
        quadpoint(5,:) = [ 0.4701420641051, 0.4701420641051 ];
        quadpoint(6,:) = [ 0.0597158717898, 0.4701420641051 ];
        quadpoint(7,:) = [ 0.33333333333333, 0.33333333333333 ];

        quadweight(1) = 0.1259391805448;
        quadweight(2) = 0.1259391805448;
        quadweight(3) = 0.1259391805448;
        quadweight(4) = 0.1323941527885;
        quadweight(5) = 0.1323941527885;
        quadweight(6) = 0.1323941527885;
        quadweight(7) = 0.2250000000000;

    else
        quadpoint = zeros( 13, 2 );
        quadweight = zeros( 13, 1 );

        quadpoint(1,:) = [ 0.0651301029022, 0.0651301029022 ];
        quadpoint(2,:) = [ 0.8697397941956, 0.0651301029022 ];
        quadpoint(3,:) = [ 0.0651301029022, 0.8697397941956 ];
        quadpoint(4,:) = [ 0.3128654960049, 0.0486903154253 ];
        quadpoint(5,:) = [ 0.6384441885698, 0.3128654960049 ];
        quadpoint(6,:) = [ 0.0486903154253, 0.6384441885698 ];
        quadpoint(7,:) = [ 0.6384441885698, 0.0486903154253 ];
        quadpoint(8,:) = [ 0.3128654960049, 0.6384441885698 ];

```

```

quadpoint(9,:) = [ 0.0486903154253, 0.3128654960049 ];
quadpoint(10,:) = [ 0.2603459660790, 0.2603459660790 ];
quadpoint(11,:) = [ 0.4793080678419, 0.2603459660790 ];
quadpoint(12,:) = [ 0.2603459660790, 0.4793080678419 ];
quadpoint(13,:) = [ 0.33333333333333, 0.33333333333333 ];

quadweight(1) = 0.0533472356088;
quadweight(2) = 0.0533472356088;
quadweight(3) = 0.0533472356088;
quadweight(4) = 0.0771137608903;
quadweight(5) = 0.0771137608903;
quadweight(6) = 0.0771137608903;
quadweight(7) = 0.0771137608903;
quadweight(8) = 0.0771137608903;
quadweight(9) = 0.0771137608903;
quadweight(10) = 0.1756152576332;
quadweight(11) = 0.1756152576332;
quadweight(12) = 0.1756152576332;
quadweight(13) = -0.1495700444677;

end

Q=quadpoint;
W=quadweight/2;
end

end % end of function

```



```

if size(coord,2) < 2
    disp('Error two coordinates needed for the Q4 element')
else

    xi=coord(1); eta=coord(2);

    N=1/4*[ (1-xi)*(1-eta);
            (1+xi)*(1-eta);
            (1+xi)*(1+eta);
            (1-xi)*(1+eta)];

    dNdx=1/4*[-(1-eta), -(1-xi);
               1-eta,    -(1+xi);
               1+eta,    1+xi;
               -(1+eta), 1-xi];
end

case 'Q8'
%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%% Q9 NINE NODE QUADRILATERAL ELEMENT %%%%%%%%%
%
%      4-----7-----3
%      |                               |
%      |                               |
%      |                               |
%      |                               |
%      8                               6
%      |                               |
%      |                               |
%      |                               |
%      1-----5-----2
%
if size(coord,2) < 2
    disp('Error two coordinates needed for the Q8 element')
else
    xi=coord(1); eta=coord(2);
    N=1/4*[-1*(1-xi)*(1-eta)*(1+xi+eta);
            -1*(1+xi)*(1-eta)*(1-xi+eta);
            -1*(1+xi)*(1+eta)*(1-xi-eta);
            -1*(1-xi)*(1+eta)*(1+xi-eta);
            2*(1-xi^2)*(1-eta);
            2*(1+xi)*(1-eta^2);
            2*(1-xi^2)*(1+eta);
            2*(1-xi)*(1-eta^2)];

    dNdx=1/4*[(1-eta)*(2*xi+eta), (1-xi)*(2*eta+xi);
               (1-eta)*(2*xi-eta), (1+xi)*(2*eta-xi);
               (1+eta)*(2*xi+eta), (1+xi)*(2*eta+xi);
               (1+eta)*(2*xi-eta), (1-xi)*(2*eta-xi);
               -4*xi*(1-eta),      -2*(1-xi^2);
               2*(1-eta^2),        -4*eta*(1+xi);
               -4*xi*(1+eta),      2*(1-xi^2);
               -2*(1-eta^2),      -4*eta*(1-xi)];
end

```

```

case 'Q9'
%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%% Q9 NINE NODE QUADRILATERAL ELEMENT %%%%%%%%%%
%
%      4-----7-----3
%      |                   |
%      |                   |
%      |                   |
%      8           9       6
%      |                   |
%      |                   |
%      |                   |
%      1-----5-----2
%
if size(coord,2) < 2
    disp('Error two coordinates needed for the Q9 element')
else
    xi=coord(1); eta=coord(2);
    N=1/4*[xi*eta*(xi-1)*(eta-1);
           xi*eta*(xi+1)*(eta-1);
           xi*eta*(xi+1)*(eta+1);
           xi*eta*(xi-1)*(eta+1);
          -2*eta*(xi+1)*(xi-1)*(eta-1);
          -2*xi*(xi+1)*(eta+1)*(eta-1);
          -2*eta*(xi+1)*(xi-1)*(eta+1);
          -2*xi*(xi-1)*(eta+1)*(eta-1);
           4*(xi+1)*(xi-1)*(eta+1)*(eta-1)];

    dNdxi=1/4*[eta*(2*xi-1)*(eta-1),    xi*(xi-1)*(2*eta-1);
               eta*(2*xi+1)*(eta-1),    xi*(xi+1)*(2*eta-1);
               eta*(2*xi+1)*(eta+1),    xi*(xi+1)*(2*eta+1);
               eta*(2*xi-1)*(eta+1),    xi*(xi-1)*(2*eta+1);
               -4*xi*eta*(eta-1),       -2*(xi+1)*(xi-1)*(2*eta-1);
               -2*(2*xi+1)*(eta+1)*(eta-1), -4*xi*eta*(xi+1);
               -4*xi*eta*(eta+1),       -2*(xi+1)*(xi-1)*(2*eta+1);
               -2*(2*xi-1)*(eta+1)*(eta-1), -4*xi*eta*(xi-1);
               8*xi*(eta^2-1),          8*eta*(xi^2-1)];

    end

otherwise
    disp(['Element ',type,' not yet supported'])
    N=[]; dNdxi=[];
end

I=eye(dim);
Nv=[];
for i=1:size(N,1)
    Nv=[Nv;I*N(i)];
end
end

```

```
if ( dim == 1 )
    B=dNdx1;
elseif ( dim == 2 )
    B=zeros(dim*size(N,1),3);

    B(1:dim:dim*size(N,1)-1,1) = dNdx1(:,1);
    B(2:dim:dim*size(N,1),2)   = dNdx1(:,2);

    B(1:dim:dim*size(N,1)-1,3) = dNdx1(:,2);
    B(2:dim:dim*size(N,1),3)   = dNdx1(:,1);

end
end % end of function
```


displacements

```
function [U,u_x,u_y] =displacements(dispNodes,dispNodes1,numnode,K,f,selfwt)

% evaluates the unknown degree of freedom (displacements) at the nodes

total_unknown=2*numnode;
udofs = [(dispNodes.*2)-1;(dispNodes1.*2)-1]; %prescribed disp.in x-dir
vdofs = dispNodes.*2; %prescribed disp. in y-dir
dofs=union(udofs(:),vdofs(:)); %overall prescribed disp.
unknowndof=setdiff((1:total_unknown)',dofs);

F=f(unknowndof)+selfwt(unknowndof);
u=K(unknowndof,unknowndof)\F;
U=zeros(total_unknown,1);
U(unknowndof)=u;
u_x = U(1:2:2*numnode-1) ; % 1 3 5 7 ...
u_y = U(2:2:2*numnode) ; % 2 4 6 8 ...

end % end of function
```

force_matrix

```
function f=force_matrix(node,topEdge,sigmatoy,sigmatox,load_edge1,load_edge2)

% Generates the force matrix due to externally applied loads
% for Q4 and T3 elements.

numnode = size(node,1);
total_unknown = numnode*2;
f = zeros(total_unknown,1);
[W,Q]=gauss_pt_wt(1,'GAUSS',1);

ii1=intersect(find(node(:,1)==load_edge1),unique(topEdge));
jj1=intersect(find(node(:,1)==load_edge2),unique(topEdge));
ii2=find(topEdge(:,1)==ii1);
jj2=find(topEdge(:,2)==jj1);

for e =ii2:jj2
    sctr = topEdge(e,:);
    sctry = sctr.*2 ;
    sctrx = sctr.*2-1;
    for q=1:size(W,1)
        pt = Q(q,:);
        wt = W(q);
        N = shape_func('L2',pt);
        J0 = abs( node(sctr(2))-node(sctr(1)) )/2;
        f(sctry) = f(sctry) + N*sigmatoy*det(J0)*wt;
        f(sctrx) = f(sctrx) + N*sigmatox*det(J0)*wt;
    end % end of quadrature loop
end % end of element loop

end % end of function
```

elementdof

```
function eldof =elementdof(elemType,sctr)

% forms the global degree of freedom nodes in each element from the node
% identification number.

switch (elemType)
case 'Q9'
    eldof=[sctr(1)*2-1;sctr(1)*2;...
           sctr(2)*2-1;sctr(2)*2;...
           sctr(3)*2-1;sctr(3)*2;...
           sctr(4)*2-1;sctr(4)*2;...
           sctr(5)*2-1;sctr(5)*2;...
           sctr(6)*2-1;sctr(6)*2;...
           sctr(7)*2-1;sctr(7)*2;...
           sctr(8)*2-1;sctr(8)*2;...
           sctr(9)*2-1;sctr(9)*2];
case 'Q8'
    eldof=[sctr(1)*2-1;sctr(1)*2;...
           sctr(2)*2-1;sctr(2)*2;...
           sctr(3)*2-1;sctr(3)*2;...
           sctr(4)*2-1;sctr(4)*2;...
           sctr(5)*2-1;sctr(5)*2;...
           sctr(6)*2-1;sctr(6)*2;...
           sctr(7)*2-1;sctr(7)*2;...
           sctr(8)*2-1;sctr(8)*2];
case 'Q4'
    eldof=[sctr(1)*2-1;sctr(1)*2;...
           sctr(2)*2-1;sctr(2)*2;...
           sctr(3)*2-1;sctr(3)*2;...
           sctr(4)*2-1;sctr(4)*2];
case 'T6'
    eldof=[sctr(1)*2-1;sctr(1)*2;...
           sctr(2)*2-1;sctr(2)*2;...
           sctr(3)*2-1;sctr(3)*2;...
           sctr(4)*2-1;sctr(4)*2;...
           sctr(5)*2-1;sctr(5)*2;...
           sctr(6)*2-1;sctr(6)*2];
case 'T3'
    eldof=[sctr(1)*2-1;sctr(1)*2;...
           sctr(2)*2-1;sctr(2)*2;...
           sctr(3)*2-1;sctr(3)*2];
case 'L3'
    eldof=[sctr(1)*2-1;sctr(1)*2;...
           sctr(2)*2-1;sctr(2)*2;...
           sctr(3)*2-1;sctr(3)*2];
case 'L2'
    eldof=[sctr(1)*2-1;sctr(1)*2;...
           sctr(2)*2-1;sctr(2)*2];

end

end % end of function
```

force_matrix689

```
function f=force_matrix689(node,topEdge,topEdge1,sigmatoy,sigmatox,...
                        load_edge1,load_edge2)

% Generates the force matrix due to externally applied loads
% for Q9,Q8 and T6 elements.

numnode = size(node,1);
total_unknown = numnode*2;
f = zeros(total_unknown,1);

[W,Q]=gauss_pt_wt(2,'GAUSS',1);

ii1=intersect(find(node(:,1)==load_edge1),unique(topEdge));
jj1=intersect(find(node(:,1)==load_edge2),unique(topEdge));
ii2=find(topEdge(:,1)==ii1);
jj2=find(topEdge(:,2)==jj1);
ee2=topEdge1(1:end,2).*2;ee4=ee2-1;
ee1=unique(topEdge);
ee3=ee1(1:2:end).*2;ee6=ee3-1;

for e =ii2:jj2
    sctr = topEdge(e,:);
    sctry = sctr.*2 ;
    sctrx = sctr.*2-1;

    for q=1:size(W,1)
        pt = Q(q,:);
        wt = W(q);
        [N,dNdx1] = shape_func('L2',pt);
        J0 = abs( node(sctr(2))-node(sctr(1)) )/2;
        f(sctry) = f(sctry) + N*sigmatoy*det(J0)*wt;
        f(sctrx) = f(sctrx) + N*sigmatox*det(J0)*wt;
    end % of quadrature loop

end % of element loop

f(ee3)=f(ee3)*(2/3);
f(ee2)=f(ee2)*(4/3);
f(ee6)=f(ee6)*(2/3);
f(ee4)=f(ee4)*(4/3);

end % end of function
```

formdg

```
function [dg1,dg2,dg3] =formdg(tsi,q,theta)

% calculates the partial derivatives of the plastic potential with respect
% to p, J2 and J3.

    dg1=sind(tsi);
if sind(theta)>0.49      % close to theta=30 corner,smoothen the curve with
    sw=1;              % triaxial compression case s1 > s2 = s3
    dg2=(0.25/q)*(3-sw*sind(tsi));
    dg3=0;

elseif -1*sind(theta)>049 % close to theta=-30 corner,smoothen the curve
    sw=-1;             % with triaxial extension case s1 = s2 > s3
    dg2=(0.25/q)*(3-sw*sind(tsi));
    dg3=0;
else                  % all other cases
    dg2=(sqrt(3)*cosd(theta)/(2*q))*(1+(tand(theta)*tand(3*theta))+...
        ((sind(tsi)/sqrt(3))*(tand(3*theta)-tand(theta))));
    dg3=1.5*((sqrt(3)*sind(theta))+(sind(tsi)*cosd(theta))/(q^2*cosd(3*theta)));

end
end % end of function
```

formm

```
function [m1,m2,m3] = formm(kk,iel,stress)

%Generates the partial derivatives of the p, J2 and J3 with respect to
%stress for use in the plastic potential derivative.

    s_xx=stress(1,kk,iel);s_yy=stress(2,kk,iel);
    t_xy=stress(3,kk,iel);s_zz=stress(4,kk,iel);
    p1=(s_xx+s_yy+s_zz)/3;

    sx=s_xx-p1;
    sy=s_yy-p1;
    sz=s_zz-p1;

    m1=1/(3*(s_xx+s_yy+s_zz))*[1 1 0 1;1 1 0 1;0 0 0 0;1 1 0 1];

    m2=1/3*[2 -1 0 -1;-1 2 0 -1;0 0 6 0;-1 -1 0 2];

    m3=1/3*[sx sz t_xy sz;sz sy t_xy sx;t_xy t_xy -3*sz -2*t_xy;sz sx -2*t_xy
sz];

end % end of function
```

internalrxn

```
function [r] = internalrxn( node,element,elemType,normal_order,stress)

%Generates the vector of nodal force reactions due to internal stress. it
%can be used out side the iteration loop if the stresses at each gauss
%point is given.

numnode = size(node,1);
numelem = size(element,1);
total_unknown = numnode*2;
r=zeros(total_unknown,1);

for iel = 1 : numelem
    sctr = element(iel,:); % element connectivity
    nn = length(sctr); % number of nodes per element
    eldof =elementdof(elemType,sctr); %element degree of freedom
    [W,Q] = gauss_rule(iel,elemType,normal_order); % determine GP

    for kk = 1 :nn
        pt =Q(kk,:);
        [N,dNdxi] = shape_func(elemType,pt);
        J0 = node(sctr,:)'*dNdxi;
        Bfem4 =Bmatrix4(pt,elemType,iel);
        r(eldof) =r(eldof)+Bfem4'*stress(:,kk,iel)*W(kk)*det(J0);
    end % end of looping on Gauss Points
end % end of looping on elements
end % end of function
```

invariants

```
function [p,q,theta] = invariants( stress )

%calculates the stress invariants for a single gauss point

s_xx=stress(1); s_yy=stress(2); p=(s_xx+s_yy+s_zz)/3;
t_xy=stress(3); s_zz=stress(4);
t=sqrt((s_xx-s_yy)^2+(s_yy-s_zz)^2+(s_zz-s_xx)^2+6*t_xy^2)/sqrt(3);
q=sqrt(1.5)*t;
sx=s_xx-p; sy=s_yy-p; sz=s_zz-p;
if q<1e-6
    theta=0;
else
    j3=sx*sy*sz-(sx*t_xy^2);
    sine=-3*j3*sqrt(6)/t^3;
    if sine>1;
        sine=1;
    end
    if sine<-1;
        sine=-1;
    end
    theta=1/3*(asind(sine));
end

end % end of function
```

invariants1

```
function [p,q,theta] =invariants1(element, stress)

% calculates the stress invariants at each gauss point.This can be used
% outside the iteration loop.

numelem=size(element,1);
nonelm=size(element,2);
p=zeros(1,nonelm,numelem);
q=zeros(1,nonelm,numelem);
theta=zeros(1,nonelm,numelem);

for iel = 1 : numelem
    sctr = element(iel,:); % element connectivity
    nn    = length(sctr);  % number of nodes per element

    for kk = 1:nn
        s_xx=stress(1, kk, iel);s_yy=stress(2, kk, iel);
        t_xy=stress(3, kk, iel);s_zz=stress(4, kk, iel);
        p1=(s_xx+s_yy+s_zz)/3;
        t=sqrt((s_xx-s_yy)^2+(s_yy-s_zz)^2+(s_zz-s_xx)^2+6*t_xy^2)/sqrt(3);
        q1=sqrt(1.5)*t;
        sx=s_xx-p1;
        sy=s_yy-p1;
        sz=s_zz-p1;
        if q1<1e-6
            theta11=0;
        else
            j3=sx*sy*sz-(sx*t_xy^2);
            sine=-3*j3*sqrt(6)/t^3;
            if sine>1;
                sine=1;
            end
            if sine<-1;
                sine=-1;
            end
            theta11=1/3*(asind(sine));
        end
        p(:,kk,iel)=p1;
        q(:,kk,iel)=q1;
        theta(:,kk,iel)=theta11;
    end
end
    % end of looping on GPs
end
    % end of looping on elements

end
    % end of function
```

invariants2

```
function [p,q,theta] =invariants2(kk,iel, stress)

% calculates the stress invariants at each gauss point.This can be used
% inside the iteration loop.

s_xx=stress(1, kk, iel); s_yy=stress(2, kk, iel);
t_xy=stress(3, kk, iel); s_zz=stress(4, kk, iel);
p=(s_xx+s_yy+s_zz)/3;
t=sqrt((s_xx-s_yy)^2+(s_yy-s_zz)^2+(s_zz-s_xx)^2+6*t_xy^2)/sqrt(3);
q=sqrt(1.5)*t;
sx=s_xx-p;
sy=s_yy-p;
sz=s_zz-p;
    if q<1e-6
        theta=0;
    else
        j3=sx*sy*sz-(sx*t_xy^2);
        sine=-3*j3*sqrt(6)/t^3;
        if sine>1;
            sine=1;
        end
        if sine<-1;
            sine=-1;
        end
        theta=1/3*(asind(sine));
    end

end % end of function
```

makedfds

```
function [dfds] =makedfds(phi, element, iel, kk, stress)

%Forms the partial derivatives of the yield function with
%respect to stress at each gauss point.

numelem=size(element,1);
nonelm=size(element,2);
dfds=zeros(4, nonelm, numelem);

s_xx=stress(1, kk, iel); s_yy=stress(2, kk, iel);
t_xy=stress(3, kk, iel); s_zz=stress(4, kk, iel);
p1=(s_xx+s_yy+s_zz)/3;

t=sqrt((s_xx-s_yy)^2+(s_yy-s_zz)^2+(s_zz-s_xx)^2+6*t_xy^2)/sqrt(3);
q1=sqrt(1.5)*t;
sx=s_xx-p1;
sy=s_yy-p1;
sz=s_zz-p1;
j2=-sx*sy-sy*sz-sz*sx+t_xy^2;
j3=sx*sy*sz-(sx*t_xy^2);
th=-3*sqrt(3)*j3/(2*j2^1.5);
```

```

        if th>1
            th=1;
        end
        if th<-1
            th=-1;
        end
        theta1=asind(th)/3;

if sind(theta1)>0.49 % close to theta=30 corner,smoothen the curve
                    % with triaxial compression case s1 > s2 = s3
    sw=-1;
    cphi=0.25*sqrt(3/j2)*(1+(sw*sind(phi)/3));

    dfds(:,kk,iel)=[(sind(phi)/3)+sx*cphi;...
                    (sind(phi)/3)+sy*cphi;...
                    t_xy*cphi;...
                    (sind(phi)/3)+sz*cphi];

elseif (-1*sind(theta1))>0.49 % close to theta=-30 corner,smoothen the curve
                                % curve with triaxial extension case s1 = s2 > s3
    sw=1;
    cphi=0.25*sqrt(3/j2)*(1+(sw*sind(phi)/3));

    dfds(:,kk,iel)=[(sind(phi)/3)+sx*cphi;...
                    (sind(phi)/3)+sy*cphi;...
                    t_xy*cphi;...
                    (sind(phi)/3)+sz*cphi];
else
    alpha=atand(abs((s_xx-s_yy)/(2*t_xy))); % all other conditions
    k1=1;k2=1;
        if abs(s_xx)>abs(s_yy)
            k1=-1;
        end
        if t_xy <0
            k2=-1;
        end

    dfds(:,kk,iel)=[sind(phi)+k1*sind(alpha);...
                    sind(phi)-k1*sind(alpha);...
                    2*k2*cosd(alpha);...
                    0];

end

end % end of function

```


makedgds

```
function [dgds] =makedgds(tsi,element,iel,kk, stress)

%Forms the partial derivatives of the plastic potential function with
%respect to stress at each gauss point.

numelem=size(element,1);
nonelm=size(element,2);
dgds=zeros(4,nonelm,numelem);

    s_xx=stress(1,kk,iel);s_yy=stress(2,kk,iel);
    t_xy=stress(3,kk,iel);s_zz=stress(4,kk,iel);
    p1=(s_xx+s_yy+s_zz)/3;

    t=sqrt((s_xx-s_yy)^2+(s_yy-s_zz)^2+(s_zz-s_xx)^2+6*t_xy^2)/sqrt(3);
    q1=sqrt(1.5)*t;
    sx=s_xx-p1;
    sy=s_yy-p1;
    sz=s_zz-p1;
    j2=-sx*sy-sy*sz-sz*sx+t_xy^2;
    j3=sx*sy*sz-(sx*t_xy^2);
    th=-3*sqrt(3)*j3/(2*j2^1.5);

        if th>1
            th=1;
        end
        if th<-1
            th=-1;
        end
        theta1=asind(th)/3;

    if sind(theta1)>0.49 % close to theta=30 corner,smoothen the curve
        % with triaxial compression case s1 > s2 = s3
        sw=-1;
        ctsi=0.25*sqrt(3/j2)*(1+(sw*sind(tsi)/3));

        dgds(:,kk,iel)=[(sind(tsi)/3)+sx*ctsi;...
            (sind(tsi)/3)+sy*ctsi;...
            t_xy*ctsi;...
            (sind(tsi)/3)+sz*ctsi];

    elseif (-1*sind(theta1))>0.49 % close to theta=-30 corner,smoothen the curve
        % curve with triaxial extension case s1 = s2 > s3
        sw=1;
        ctsi=0.25*sqrt(3/j2)*(1+(sw*sind(tsi)/3));

        dgds(:,kk,iel)=[(sind(tsi)/3)+sx*ctsi;...
            (sind(tsi)/3)+sy*ctsi;...
            t_xy*ctsi;...
            (sind(tsi)/3)+sz*ctsi];

    else
```

```

alpha=atand(abs((s_xx-s_yy)/(2*t_xy))); % all other conditions
k1=1;k2=1;
    if abs(s_xx)>abs(s_yy)
        k1=-1;
    end
    if t_xy <0
        k2=-1;
    end

dgds(:,kk,iel)=[sind(tsi)+k1*sind(alpha);...
sind(tsi)-k1*sind(alpha);...
2*k2*cosd(alpha);...
0];

end

end % end of function

plastic_mat

function [CP] = plastic_mat(E,nu,phi,tsi,iel,kk, stress)

% Generates the plastic constitutive matrix CP at each gauss point

s_xx=stress(1,kk,iel);s_yy=stress(2,kk,iel);
t_xy=stress(3,kk,iel);s_zz=stress(4,kk,iel);
p1=(s_xx+s_yy+s_zz)/3;

t=sqrt((s_xx-s_yy)^2+(s_yy-s_zz)^2+(s_zz-s_xx)^2+6*t_xy^2)/sqrt(3);
q=sqrt(1.5)*t;
sx=s_xx-p1;
sy=s_yy-p1;
sz=s_zz-p1;
j2=-sx*sy-sy*sz-sz*sx+t_xy^2;
j3=sx*sy*sz-(sx*t_xy^2);
th=-3*sqrt(3)*j3/(2*j2^1.5);
    if th>1
        th=1;
    end
    if th<-1
        th=-1;
    end
    theta1=asind(th)/3;

if sind(theta1)>0.495 % close to theta=30 corner,smoothen the curve
% with triaxial compression case s1 > s2 = s3
sw=-1;
cphi=0.25*sqrt(3/j2)*(1+(sw*sind(phi)/3));
ctsi=0.25*sqrt(3/j2)*(1+(sw*sind(tsi)/3));
kphi=sind(phi)*(1+nu)/3;
ktsi=sind(tsi)*(1+nu)/3;
c1=kphi+cphi*(sx*(1-nu)+nu*(sy+sz));
c2=kphi+cphi*(sy*(1-nu)+nu*(sx+sz));
c3=cphi*(1-2*nu)*t_xy;
c4=kphi+cphi*(sz*(1-nu)+nu*(sx+sy));

```

```

r1=ktsi+ctsi*(sx*(1-nu)+nu*(sy+sz));
r2=ktsi+ctsi*(sy*(1-nu)+nu*(sx+sz));
r3=ctsi*(1-2*nu)*t_xy;
r4=ktsi+ctsi*(sz*(1-nu)+nu*(sx+sy));
A=[r1*c1 r1*c2 r1*c3 r1*c4;...
   r2*c1 r2*c2 r2*c3 r2*c4;...
   r3*c1 r3*c2 r3*c3 r3*c4;...
   r4*c1 r4*c2 r4*c3 r4*c4];
CP=E*A/((1+nu)*(1-2*nu)*(kphi*sind(tsi)+2*cphi*ctsi*j2*(1-2*nu)));

elseif -1*sind(theta1)> 0.49
    sw=1; % close to theta=-30 corner,smoothen the curve
          % with triaxial extension case s1 = s2 > s3
    cphi=0.25*sqrt(3/j2)*(1+(sw*sind(phi)/3));
    ctsi=0.25*sqrt(3/j2)*(1+(sw*sind(tsi)/3));
    kphi=sind(phi)*(1+nu)/3;
    ktsi=sind(tsi)*(1+nu)/3;
    c1=kphi+cphi*(sx*(1-nu)+nu*(sy+sz));
    c2=kphi+cphi*(sy*(1-nu)+nu*(sx+sz));
    c3=cphi*(1-2*nu)*t_xy;
    c4=kphi+cphi*(sz*(1-nu)+nu*(sx+sy));
    r1=ktsi+ctsi*(sx*(1-nu)+nu*(sy+sz));
    r2=ktsi+ctsi*(sy*(1-nu)+nu*(sx+sz));
    r3=ctsi*(1-2*nu)*t_xy;
    r4=ktsi+ctsi*(sz*(1-nu)+nu*(sx+sy));
    A=[r1*c1 r1*c2 r1*c3 r1*c4;...
       r2*c1 r2*c2 r2*c3 r2*c4;...
       r3*c1 r3*c2 r3*c3 r3*c4;...
       r4*c1 r4*c2 r4*c3 r4*c4];
    CP=E*A/((1+nu)*(1-2*nu)*(kphi*sind(tsi)+2*cphi*ctsi*j2*(1-2*nu)));

else % all other cases
    alpha=atand(abs((s_xx-s_yy)/(2*t_xy)));
    k1=1;k2=1;
    if abs(s_xx)>abs(s_yy)
        k1=-1;
    end
    if t_xy <0
        k2=-1;
    end
    c1=sind(phi)+k1*(1-2*nu)*sind(alpha);
    c2=sind(phi)-k1*(1-2*nu)*sind(alpha);
    c3=k2*(1-2*nu)*cosd(alpha);
    c4=2*nu*sind(phi);
    r1=sind(tsi)+k1*(1-2*nu)*sind(alpha);
    r2=sind(tsi)-k1*(1-2*nu)*sind(alpha);
    r3=k2*(1-2*nu)*cosd(alpha);
    r4=2*nu*sind(tsi);
    A=[r1*c1 r1*c2 r1*c3 r1*c4;...
       r2*c1 r2*c2 r2*c3 r2*c4;...
       r3*c1 r3*c2 r3*c3 r3*c4;...
       r4*c1 r4*c2 r4*c3 r4*c4];
    CP=E*A/(2*(1+nu)*(1-2*nu)*(1-2*nu+sind(phi)*sind(tsi)));

end
end % end of function

```

principal_stress

```
function [s1,s2,s3] = principal_stress(kk,iel,p,q,theta)

%calculates the principal stresses from the stress invariants

s1=p(:,kk,iel)+q(:,kk,iel)*(2/3)*sind(theta(:,kk,iel)-120);

s2=p(:,kk,iel)+q(:,kk,iel)*(2/3)*sind(theta(:,kk,iel));

s3=p(:,kk,iel)+q(:,kk,iel)*(2/3)*sind(theta(:,kk,iel)+120);

end % end of function
```

selfwt_matrix

```
function selfwt=selfwt_matrix(elemType,normal_order,gamma,node,element)

% Generates the force matrix due to self weight

numelem=size(element,1);
numnode = size(node,1);
total_unknown = numnode*2;
selfwt=zeros(total_unknown,1);

for iel = 1 : numelem
    sctrl = element(iel,:); % element connectivity
    swpt=sctrl.*2; %element degree of freedom

    [W,Q] = gauss_rule(iel,elemType,normal_order);

    for q=1:size(W,1)
        pt = Q(q,:);
        wt = W(q); % quadrature point
        [N,dNdxl]=shape_func(elemType,pt);
        J0 = node(sctrl,：)'*dNdxl; % element Jacobian matrix
        selfwt(swpt) = selfwt(swpt)+ N*(-1*gamma)*det(J0)*wt;
    end
end
end % end of function
```

stiffness_matrix

```
function K=stiffness_matrix(node,element,elemType,normal_order,C)

% Generates the element and global stiffness matrix

numnode = size(node,1);
numelem = size(element,1);
total_unknown = numnode*2;
K = zeros(total_unknown,total_unknown);

for iel = 1 : numelem
    sctr = element(iel,:);           % element connectivity
    nn    = length(sctr);           % number of nodes per element
    eldof =elementdof(elemType,sctr); %element degree of freedom

    [W,Q] = gauss_rule(iel,elemType,normal_order);

    for kk = 1 : size(W,1)           % Loop on Gauss points
        pt = Q(kk,:);               % quadrature point
        [N,dNdx] = shape_func(elemType,pt);
        J0 = node(sctr,:)'*dNdx;     % element Jacobian matrix
        Bfem =Bmatrix(pt,elemType,iel);
        K(eldof,eldof) = K(eldof,eldof)+Bfem'*C(1:3,1:3)*Bfem*W(kk)*det(J0);
    end                               % end of looping on Gauss Points
end                                   % end of looping on elements
end                                   % end of function
```

stress_calculation

```
function [stress,strain] =stress_calculation(node,element,elemType,...
                                             U,normal_order,C)

% calculates the element strains and stresses at the nodes
% in x, y and xy directions.

numelem=size(element,1);

switch elemType
    case 'Q9'
        stresspoints=[-1 -1;1 -1;1 1;-1 1;0 -1;1 0;0 1;-1 0;0 0];
    case 'Q8'
        stresspoints=[-1 -1;1 -1;1 1;-1 1;0 -1;1 0;0 1;-1 0];
    case 'Q4'
        stresspoints=[-1 -1;1 -1;1 1;-1 1];
    case 'T3'
        stresspoints=[0 0;1 0;0 1];
    otherwise
        stresspoints=[0 0;1 0;0 1;0.5 0;0.5 0.5;0 0.5];
end

for iel = 1 : numelem
    sctr = element(iel,:); % element connectivity
    nn    = length(sctr);  % number of nodes per element
```

```

eldof =elementdof(elemType, sctr);
[W,Q] = gauss_rule(iel,elemType,normal_order);

for kk = 1:nn
    pt = stresspoints(kk,:);           % quadrature point
    [N,dNdx] = shape_func(elemType,pt); % element shape functions
    J0 = node(sctr,:)'*dNdx;          % element Jacobian matrix
    Bfem =Bmatrix(pt,elemType,iel);
    strain=Bfem*U(eldof);
    stress(iel, kk, :)=C*strain;

end                                     % end of looping on gauss points

end                                     % end of looping on elements

end % end of function

```

Supportcond

```

function [topEdge,topEdge1,dispNodes,dispNodes1]= ...
    supportcond(elemType,numx,numy)

% Forms a vector of nodes which are restrained in x direction or in both
% x and y direction which are the supports of the domain.
%It also forms a matrix of nodes for load application at the top of the
%domain.

switch elemType
    case 'Q9'
        nnx=numx*2+1;
        nny=numy*2+1;
        urn =nnx*nny;           % upper right node number
        uln =urn-(nnx-1);       % upper left node number
        lrn = nnx;              % lower right node number
        llm = 1;                % lower left node number

        topEdge = [ uln:1:(urn-1); (uln+1):1:urn ]';
        topEdge1=[uln:2:(urn-2); (uln+1):2:(urn-1);(uln+2):2:urn ]';
        botEdge = [ llm:1:(lrn-1); (llm+1):1:lrn ]';
        leftEdge=[ (llm:nnx:(uln-nnx));(llm+nnx:nnx:uln) ]';
        rightEdge=[ (nnx:nnx:(urn-nnx)); (nnx+nnx:nnx:urn) ]';
        % GET NODES ON ESSENTIAL BOUNDARY

        botNodes = unique(botEdge);
        topNodes = unique(topEdge);
        leftNodes = unique(leftEdge);
        rightNodes = unique(rightEdge);
        dispNodes = botNodes;
        rightNodes1=rightNodes(2:end);
        leftNodes1=leftNodes(2:end);
        dispNodes1=union(leftNodes1,rightNodes1);

```

```

case 'Q8'
    nnx=numx*2+1;
    nny=numy*2+1;
    urn = (nnx*nny) - (numx*numy); % upper right node number
    uln =urn-(nnx-1);             % upper left node number
    lrn = nnx;                    % lower right node number
    lln = 1;                     % lower left node number

    topEdge = [ uln:1:(urn-1); (uln+1):1:urn ]';
    topEdge1=[uln:2:(urn-2); (uln+1):2:(urn-1);(uln+2):2:urn ]';
    botEdge = [ lln:1:(lrn-1); (lln+1):1:lrn ]';
    % GET NODES ON ESSENTIAL BOUNDARY
    botNodes = unique(botEdge);
    topNodes = unique(topEdge);
    leftNodes = union((lln:nnx+numx+1:uln),...
                      (nnx+1:nnx+numx+1:uln-numx+1) )';
    rightNodes = union((nnx:nnx+numx+1:urn),...
                      (nnx+numx+1:nnx+numx+1:urn-nnx) )';
    dispNodes = botNodes;
    rightNodes1=rightNodes(2:end);
    leftNodes1=leftNodes(2:end);
    dispNodes1=union(leftNodes1,rightNodes1);
    leftEdge=[leftNodes(1:nny-1,:),leftNodes1];
    rightEdge=[rightNodes(1:nny-1,:),rightNodes1];
case {'Q4','T3'}
    nnx=numx+1;
    nny=numy+1;
    uln = nnx*(nny-1)+1;         % upper left node number
    urn = nnx*nny;              % upper right node number
    lrn = nnx;                  % lower right node number
    lln = 1;                     % lower left node number
    topEdge = [ uln:1:(urn-1); (uln+1):1:urn ]';
    topEdge1=topEdge;
    botEdge = [ lln:1:(lrn-1); (lln+1):1:lrn ]';
    rightEdge = (lrn:nnx:(urn))';
    % GET NODES ON ESSENTIAL BOUNDARY

    botNodes = unique(botEdge);
    topNodes = unique(topEdge);
    rightNodes = unique(rightEdge);
    leftNodes = rightNodes-(nnx-1);
    dispNodes = botNodes;
    rightNodes1=rightNodes(2:end);
    leftNodes1=leftNodes(2:end);
    dispNodes1=union(leftNodes1,rightNodes1);
case 'T6'
    nnx=numx*2+1;
    nny=numy*2+1;
    urn =nnx*nny;                % upper right node number
    uln =urn-(nnx-1);           % upper left node number
    lrn = nnx;                  % lower right node number
    lln = 1;                     % lower left node number

    topEdge = [ uln:1:(urn-1); (uln+1):1:urn ]';
    topEdge1=[uln:2:(urn-2); (uln+1):2:(urn-1);(uln+2):2:urn ]';
    botEdge = [ lln:1:(lrn-1); (lln+1):1:lrn ]';

```

```

    % GET NODES ON ESSENTIAL BOUNDARY
    botNodes = unique(botEdge);
    topNodes = unique(topEdge);
    leftNodes = (1:nx:nlx)';
    rightNodes = (1:rx:rlx)';
    dispNodes = botNodes;
    rightNodes1=rightNodes(2:end);
    leftNodes1=leftNodes(2:end);
    dispNodes1=union(leftNodes1,rightNodes1);
    leftEdge=[leftNodes(1:nny-1,:),leftNodes1];
    rightEdge=[rightNodes(1:nny-1,:),rightNodes1];
end
end % end of function

```

plot_def

```

function plot_def(fac,u_x,u_y,elemType,dispNodes,dispNodes1)

% Plots the deformed finite element mesh with the support condition

global node element

figure
clf
hold on
plot_mesh(node+fac*[u_x u_y],element,elemType,'b-');
title(' Numerical deformed mesh ')
plot(node(dispNodes,1),node(dispNodes,2),'ks');
plot(node(dispNodes1,1),node(dispNodes1,2),'ko');

end % end of function

```

plot_defo

```

function plot_defo(fac,u_x,u_y,elemType)

% plots the color coded displacement intensity in the finite element
% region along with color bar scale.

global node element

figure
clf
subplot(2,1,1);
plot_field(node+fac*[u_x u_y],element,elemType,u_x);
colorbar
title('Deformation plot, U_X')

subplot(2,1,2);
plot_field(node+fac*[u_x u_y],element,elemType,u_y);
colorbar
title('Deformation plot, U_Y')

end % end of function

```


plot_field

```
function plot_field(node,connect,elem_type,field)

% Forms a color dependent finite element mesh for plotting outputs

if ( size(field) == size(connect) )
    elementalField=1;
else
    elementalField=0;
end
% fill node if needed
if (size(node,2) < 3)
    for c=size(node,2)+1:3
        node(:,c)=[zeros(size(node,1),1)];
    end
end

holdState=ishold;
hold on

if ( strcmp(elem_type,'Q9') )      % Q9 element
    ord=[1,5,2,6,3,7,4,8,1];
elseif ( strcmp(elem_type,'T3') ) % T3 element
    ord=[1,2,3,1];
elseif ( strcmp(elem_type,'T6') ) % T6 element
    ord=[1,4,2,5,3,6,1];
elseif ( strcmp(elem_type,'Q4') ) % Q4 element
    ord=[1,2,3,4,1];
elseif ( strcmp(elem_type,'Q8') ) % Q8 element
    ord=[1,5,2,6,3,7,4,8,1];
elseif ( strcmp(elem_type,'L2') ) % L2 element
    ord=[1,2];
end

for e=1:size(connect,1)
    xpt=node(connect(e,ord),1);
    ypt=node(connect(e,ord),2);
    zpt=node(connect(e,ord),3);

    if ( elementalField )
        fpt=field(e,ord);
    else
        fpt=field(connect(e,ord));
    end

    fill3(xpt,ypt,zpt,fpt)
end

shading interp
axis equal
if ( ~holdState )
    hold off
end
end % end of fuction
```

plot_m

```
function plot_m(elemType, dispNodes, dispNodes1)

% Plots the finite element mesh with the support condition

global node element

v=get(0, 'ScreenSize');
figure('Color', [1 1 1])
hold on
plot_mesh(node, element, elemType, 'b-');
plot(node(dispNodes, 1), node(dispNodes, 2), 'ks');
plot(node(dispNodes1, 1), node(dispNodes1, 2), 'ko');
axis off

end % end of function
```

plot_sig

```
function plot_sig(fac, u_x, u_y, elemType, stress)

% plots the color coded stress distribution in the finite element
% region along with color bar scale.

global node element
figure
clf
subplot(3,1,1);
plot_field(node+fac*[u_x u_y], element, elemType, stress(:, :, 1));
colorbar
title('Stress plot, sigma_xx')

subplot(3,1,2);
plot_field(node+fac*[u_x u_y], element, elemType, stress(:, :, 2));
colorbar
title('Stress plot, sigma_yy')

subplot(3,1,3);
plot_field(node+fac*[u_x u_y], element, elemType, stress(:, :, 3));
colorbar
title('Stress plot, sigma_xy')

end % end of function
```

plot_mesh

```
function plot_mesh(node,connect,elem_type,se)

% Plots the finite element mesh

if ( nargin < 4 )
    se='w-';
end

holdState=ishold;
hold on

% fill node if needed
if (size(node,2) < 3)
    for c=size(node,2)+1:3
        node(:,c)=[zeros(size(node,1),1)];
    end
end

for e=1:size(connect,1)

    if ( strcmp(elem_type,'Q9') )           % 9-node quad element
        ord=[1,5,2,6,3,7,4,8,1];
    elseif ( strcmp(elem_type,'Q8') )      % 8-node quad element
        ord=[1,5,2,6,3,7,4,8,1];
    elseif ( strcmp(elem_type,'T3') )      % 3-node triangle element
        ord=[1,2,3,1];
    elseif ( strcmp(elem_type,'T6') )      % 6-node triangle element
        ord=[1,4,2,5,3,6,1];
    elseif ( strcmp(elem_type,'Q4') )      % 4-node quadrilateral element
        ord=[1,2,3,4,1];
    elseif ( strcmp(elem_type,'L2') )      % 2-node line element
        ord=[1,2];
    end

    for n=1:size(ord,2)
        xpt(n)=node(connect(e,ord(n)),1);
        ypt(n)=node(connect(e,ord(n)),2);
        zpt(n)=node(connect(e,ord(n)),3);
    end
    plot3(xpt,ypt,zpt,se)

end
    axis equal

if ( ~holdState )
    hold off
end
end % end of function
```