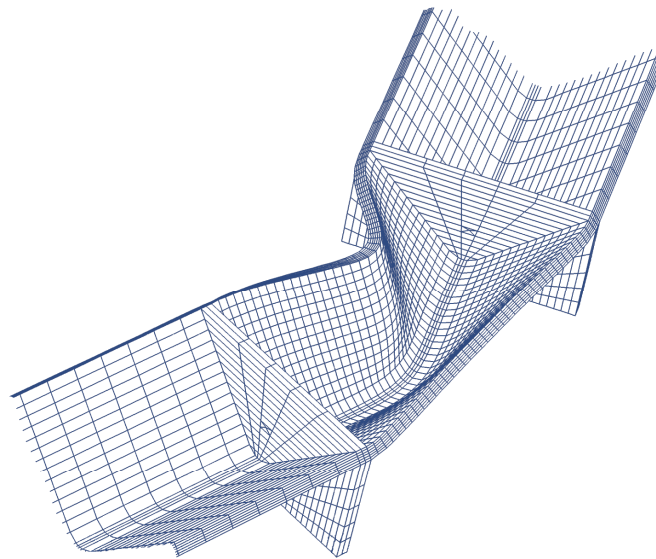
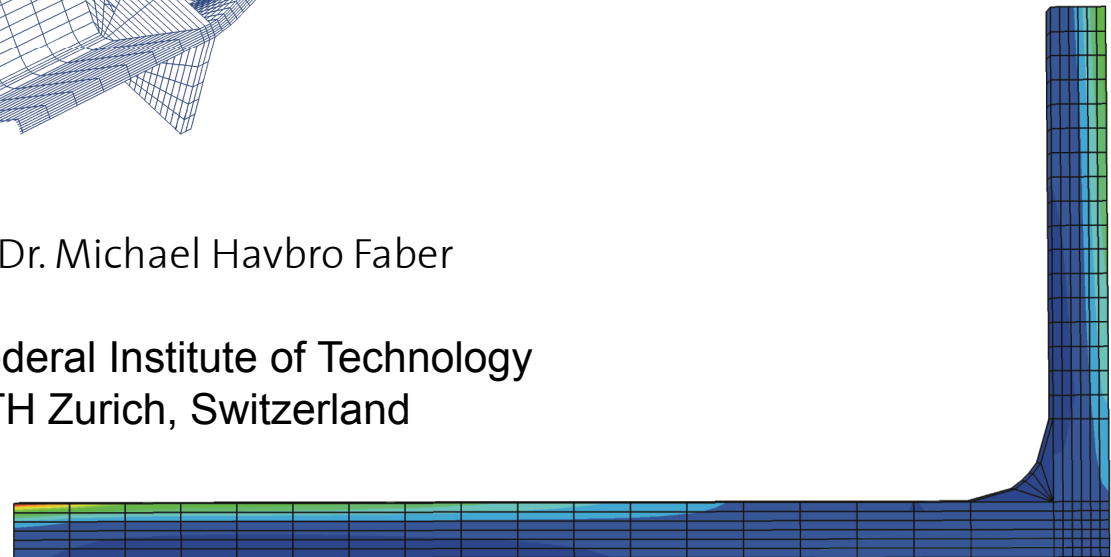


The Finite Element Method for the Analysis of Linear Systems



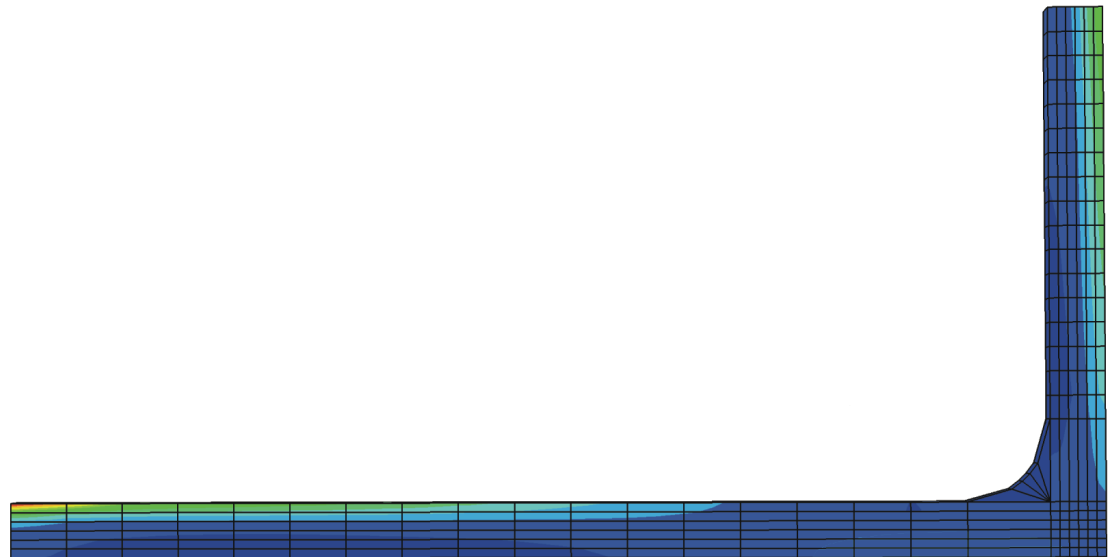
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Swiss Federal Institute of Technology
ETH Zurich, Switzerland



Contents of Today's Lecture

- **The Method of Finite Elements (principle of virtual displacements)**
- **Properties of FEM solutions**
- **On the choice of shape functions**
 - Lagrange
 - Hermitian
 - Serendipity
- **Natural coordinates**



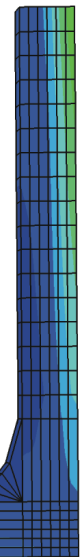
The Method of Finite Elements

The principle of virtual displacements:

We may now express the principle of virtual displacements in a more general form:

$$\underbrace{\int_V \bar{\boldsymbol{\varepsilon}}^T \boldsymbol{\tau} dV}_{\text{Internal virtual work}} = \underbrace{\int_V \bar{\mathbf{U}}^T \mathbf{f}^B dV + \int_{S_f} \bar{\mathbf{U}}^{S_f T} \mathbf{f}^{S_f} dS + \sum_i \bar{\mathbf{U}}^{iT} \mathbf{R}_C^i}_{\text{External virtual work}}$$

\uparrow \uparrow \uparrow \uparrow \uparrow \uparrow \uparrow
 Virtual strains corresponding to virtual displacements Stresses in equilibrium with applied loads



The Method of Finite Elements

Finite Element Equations:

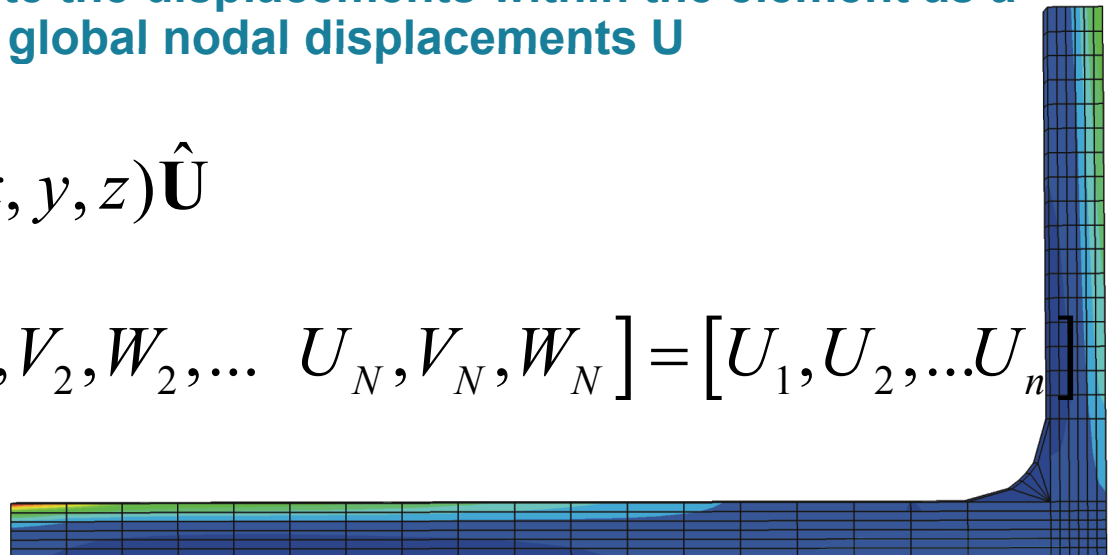
We now consider the volume modeled as an assemblage of N elements connected in the nodal points on the element boundaries

The displacements within the individual elements are measured in a convenient local coordinate system x, y, z

For element m we now write the displacements within the element as a function of the total set of global nodal displacements \mathbf{U}

$$\mathbf{u}^{(m)}(x, y, z) = \mathbf{H}^{(m)}(x, y, z) \hat{\mathbf{U}}$$

$$\hat{\mathbf{U}}^T = [U_1, V_1, W_1, U_2, V_2, W_2, \dots, U_N, V_N, W_N] = [U_1, U_2, \dots, U_n]$$



The Method of Finite Elements

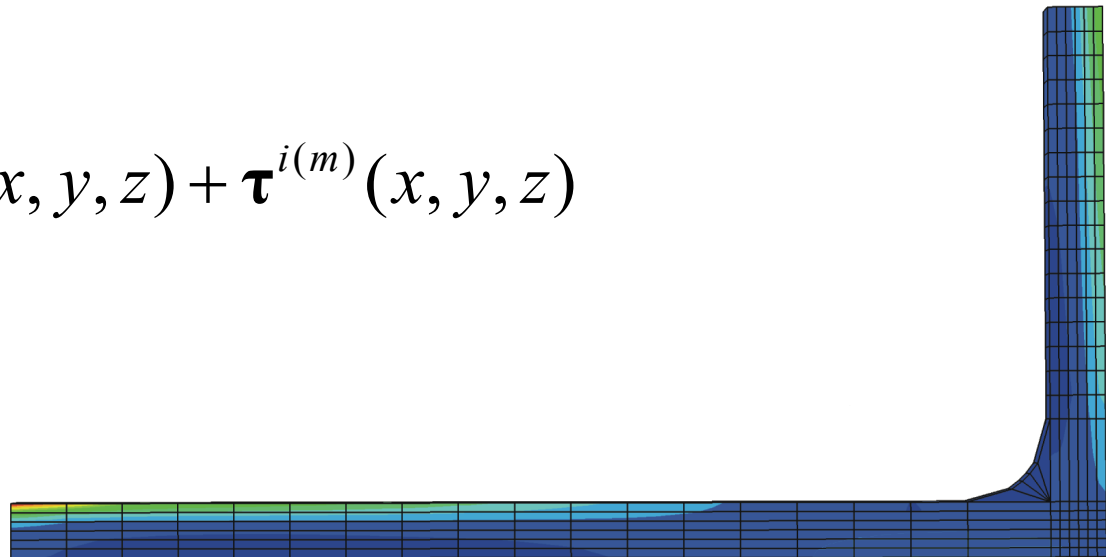
Finite Element Equations:

For element m we now write the strains within the element as a function of the total set of global nodal displacements \mathbf{U}

$$\boldsymbol{\varepsilon}^{(m)}(x, y, z) = \mathbf{B}^{(m)}(x, y, z) \hat{\mathbf{U}}$$

The stresses are then:

$$\boldsymbol{\tau}^{(m)}(x, y, z) = \mathbf{C} \boldsymbol{\varepsilon}^{(m)}(x, y, z) + \boldsymbol{\tau}^{i(m)}(x, y, z)$$



The Method of Finite Elements

Finite Element Equations:

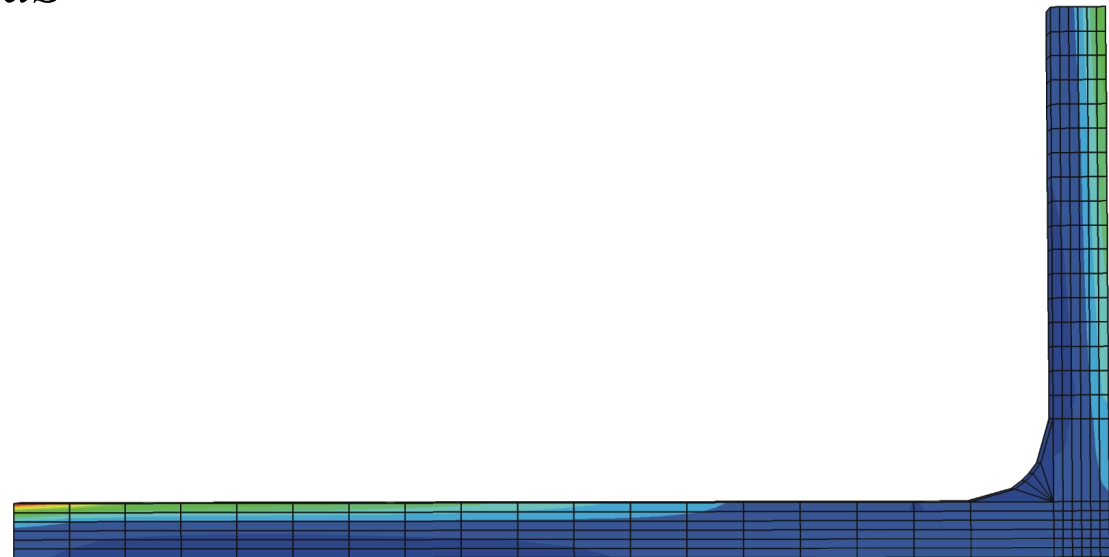
We can now write the equilibrium equations for the total volume by summing up over the N elements

$$\sum_{m=1}^N \int_{V^{(m)}} \bar{\boldsymbol{\varepsilon}}^{(m)T} \boldsymbol{\tau}^{(m)} dV^{(m)} = \sum_{m=1}^N \int_{V^{(m)}} \bar{\mathbf{U}}^{(m)T} \mathbf{f}^{B(m)} dV^{(m)} +$$

$$\sum_{m=1}^N \int_{S_{f1}^{(m)}, S_{f2}^{(m)}, \dots} \bar{\mathbf{U}}^{S_f(m)T} \mathbf{f}^{S_f(m)} dS^{(m)} -$$

$$\sum_{m=1}^N \int_{V^{(m)}} \bar{\boldsymbol{\varepsilon}}^{(m)T} \boldsymbol{\tau}^{i(m)} dV^{(m)} +$$

$$\sum_i \bar{\mathbf{U}}^{iT} \mathbf{R}_C^i$$



The Method of Finite Elements

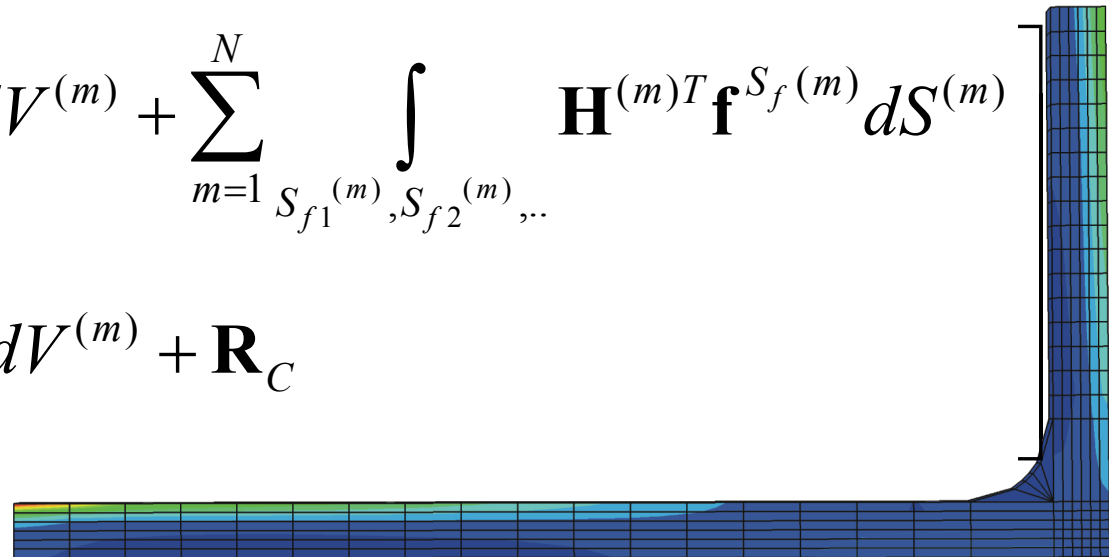
Finite Element Equations:

As a next step we represent both the real unknown displacement fields as well as the virtual displacement fields through the interpolation functions (provides symmetrical stiffness matrixes ☺)

$$\bar{\hat{\mathbf{U}}}^T \left[\sum_{m=1}^N \int_{V^{(m)}} \mathbf{B}^{(m)T} \mathbf{C}^{(m)} \mathbf{B}^{(m)} dV^{(m)} \right] \bar{\mathbf{U}} =$$

$$\bar{\hat{\mathbf{U}}}^T \left[\sum_{m=1}^N \int_{V^{(m)}} \mathbf{H}^{(m)T} \mathbf{f}^{B(m)} dV^{(m)} + \sum_{m=1}^N \int_{S_{f1}^{(m)}, S_{f2}^{(m)}, \dots} \mathbf{H}^{(m)T} \mathbf{f}^{S_f(m)} dS^{(m)} \right]$$

$$\left[- \sum_{m=1}^N \int_{V^{(m)}} \mathbf{B}^{(m)T} \boldsymbol{\tau}^{i(m)} dV^{(m)} + \mathbf{R}_C \right]$$



The Method of Finite Elements

Finite Element Equations:

Now we may finally simplify as

$$\mathbf{K}\mathbf{U} = \mathbf{R}$$

$$\mathbf{K} = \sum_{m=1}^N \int_{V^{(m)}} \mathbf{B}^{(m)T} \mathbf{C}^{(m)} \mathbf{B}^{(m)} dV^{(m)}$$

These are the finite element equations to be solved 😊

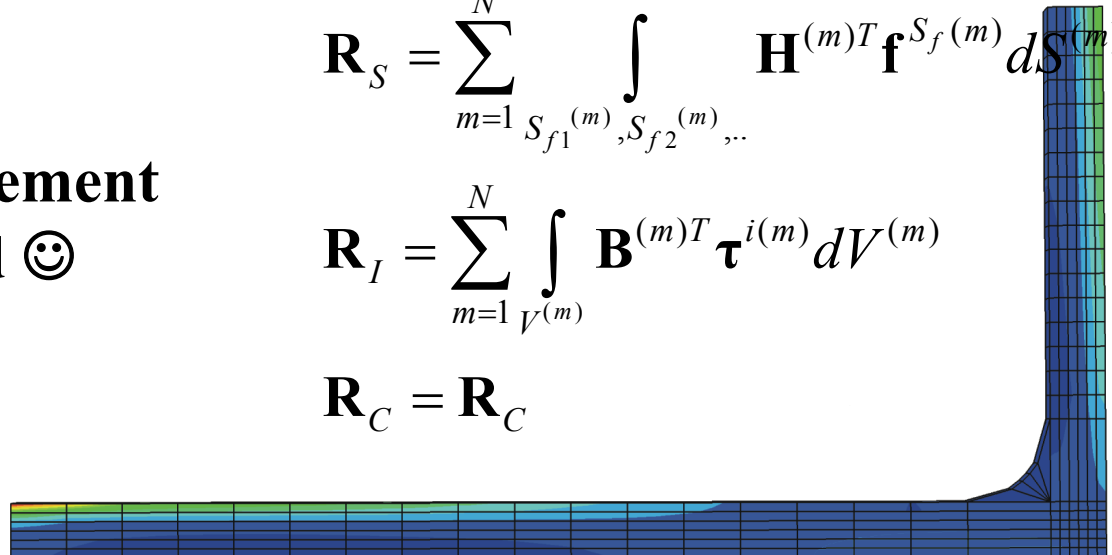
$$\mathbf{R} = \mathbf{R}_B + \mathbf{R}_S - \mathbf{R}_I + \mathbf{R}_C$$

$$\mathbf{R}_B = \sum_{m=1}^N \int_{V^{(m)}} \mathbf{H}^{(m)T} \mathbf{f}^{B(m)} dV^{(m)}$$

$$\mathbf{R}_S = \sum_{m=1}^N \int_{S_{f1}^{(m)}, S_{f2}^{(m)}, \dots} \mathbf{H}^{(m)T} \mathbf{f}^{S_f(m)} dS^{(m)}$$

$$\mathbf{R}_I = \sum_{m=1}^N \int_{V^{(m)}} \mathbf{B}^{(m)T} \boldsymbol{\tau}^{i(m)} dV^{(m)}$$

$$\mathbf{R}_C = \mathbf{R}_C$$



Generalized coordinate models

Generalized coordinate models:

The principle behind these models is:

Formulate displacement field in terms of polynomials

one-dimensional

$$u(x) = \alpha_1 + \alpha_2 x + \alpha_3 x^2 + \alpha_4 x^3 + \dots$$

two-dimensional

$$u(x, y) = \alpha_1 + \alpha_2 x + \alpha_3 xy + \alpha_4 x^2 + \dots$$

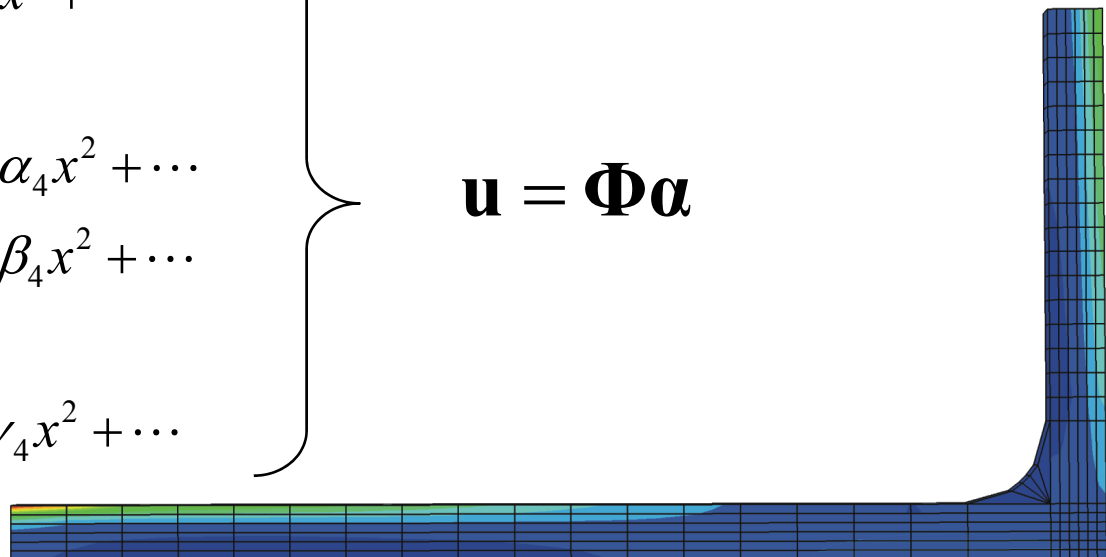
$$v(x, y) = \beta_1 + \beta_2 x + \beta_3 xy + \beta_4 x^2 + \dots$$

plate bending

$$w(x, y) = \gamma_1 + \gamma_2 x + \gamma_3 xy + \gamma_4 x^2 + \dots$$

α, β, γ : Generalized coordinates

$$\mathbf{u} = \mathbf{\Phi} \mathbf{\alpha}$$



Generalized coordinate models

Generalized coordinate models:

Next step is to relate the generalized coordinates to the nodal displacements:

We insert the nodal coordinates into

$$\mathbf{u} = \Phi \boldsymbol{\alpha}$$

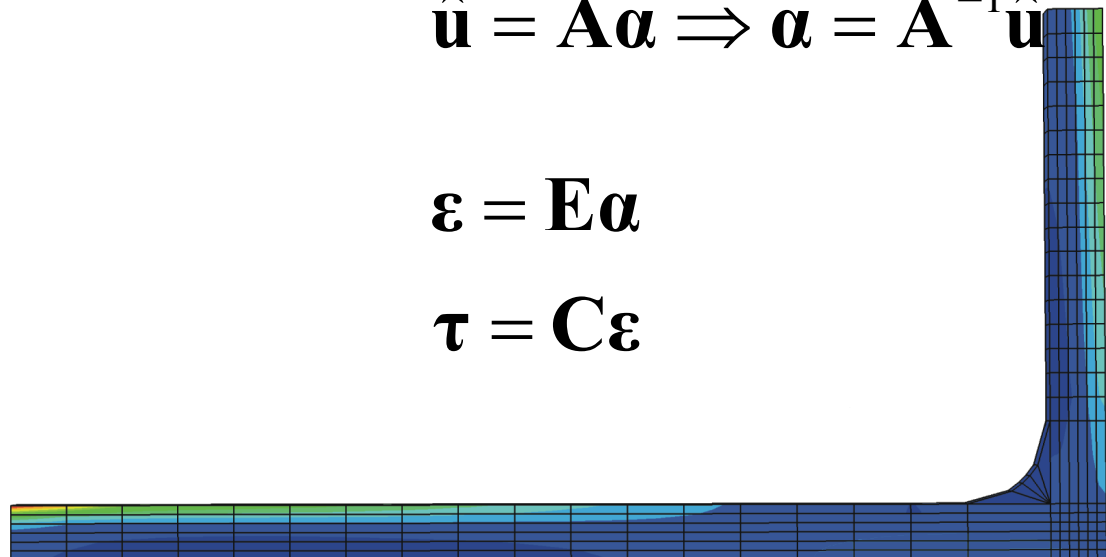
and get:

$$\hat{\mathbf{u}} = \mathbf{A} \boldsymbol{\alpha} \Rightarrow \boldsymbol{\alpha} = \mathbf{A}^{-1} \hat{\mathbf{u}}$$

Now we can obtain

$$\boldsymbol{\varepsilon} = \mathbf{E} \boldsymbol{\alpha}$$

$$\boldsymbol{\tau} = \mathbf{C} \boldsymbol{\varepsilon}$$



On the choice of shape functions

Requirements to shape functions:

First of all the functions which we want to represent (displacements and their derivatives) need to be able to represent the physics of the type of problem we are aiming to model within the individual elements.

Secondly we also need to be concerned with continuity over the borders of the elements.

We here introduce continuity requirements:

C^0 continuity

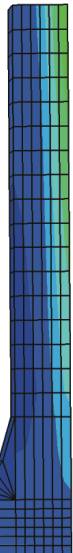
Continuity of displacement field

C^1 continuity

Continuity of the first order derivative of the displacement field

C^m continuity

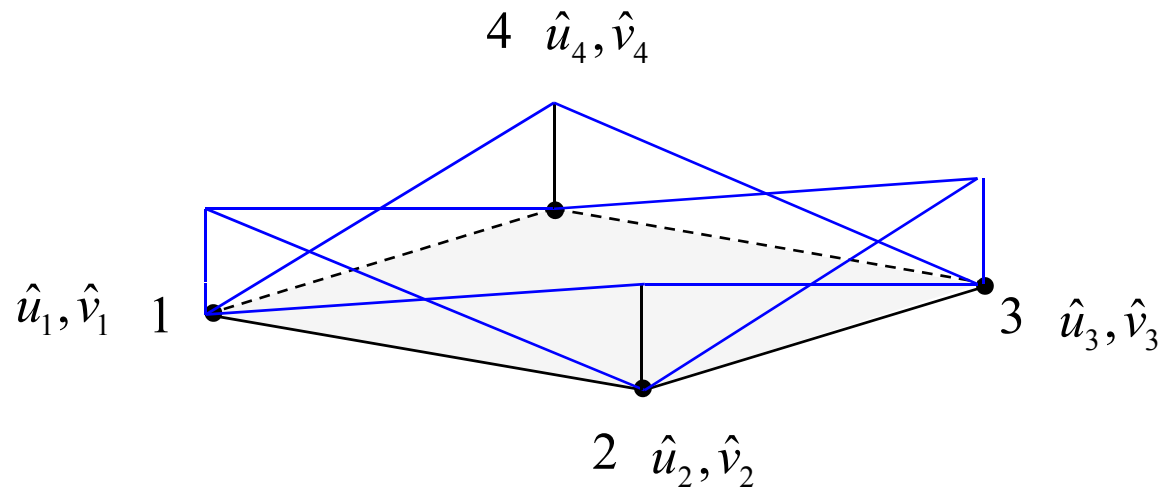
Continuity of the m^{th} order derivative of the displacement field



On the choice of shape functions

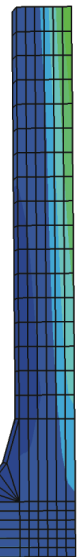
Shape functions:

The displacement e.g. u or v at any location within the element (x,y) can be represented as a function of the nodal displacements (u or v):



$$u(x, y) = f_1(\hat{\mathbf{u}})$$

$$v(x, y) = f_2(\hat{\mathbf{v}})$$



On the choice of shape functions

Shape functions:

In general we may write the approximate relation between the field representation and the nodal displacements as:

$$u(x, y) = \mathbf{H}^T(x, y) \hat{\mathbf{u}}$$

Displacements
in the u-direction

scalar

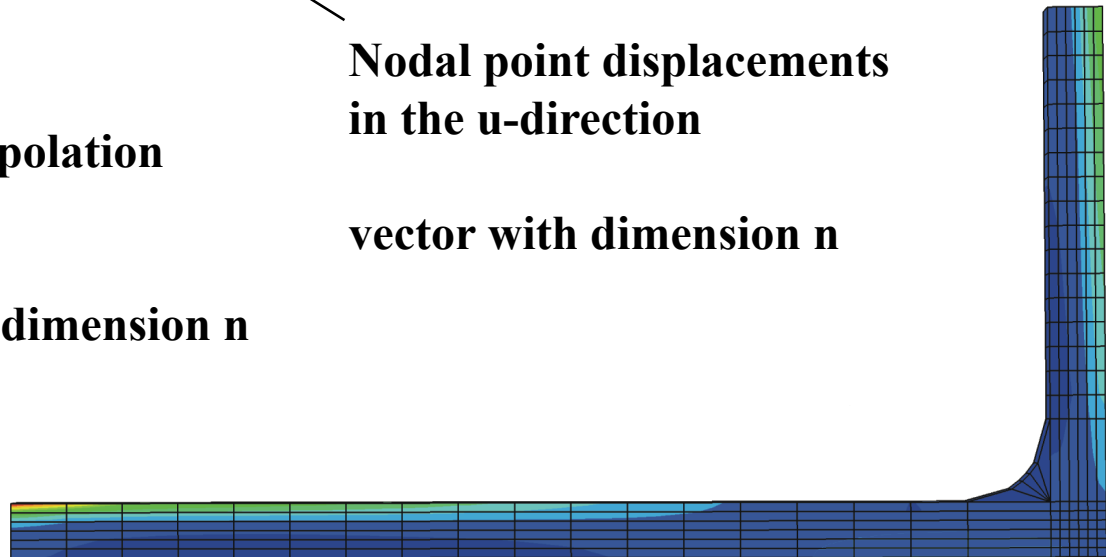
Shape/interpolation
functions

vector with dimension n

We consider an element with
n nodes

Nodal point displacements
in the u-direction

vector with dimension n

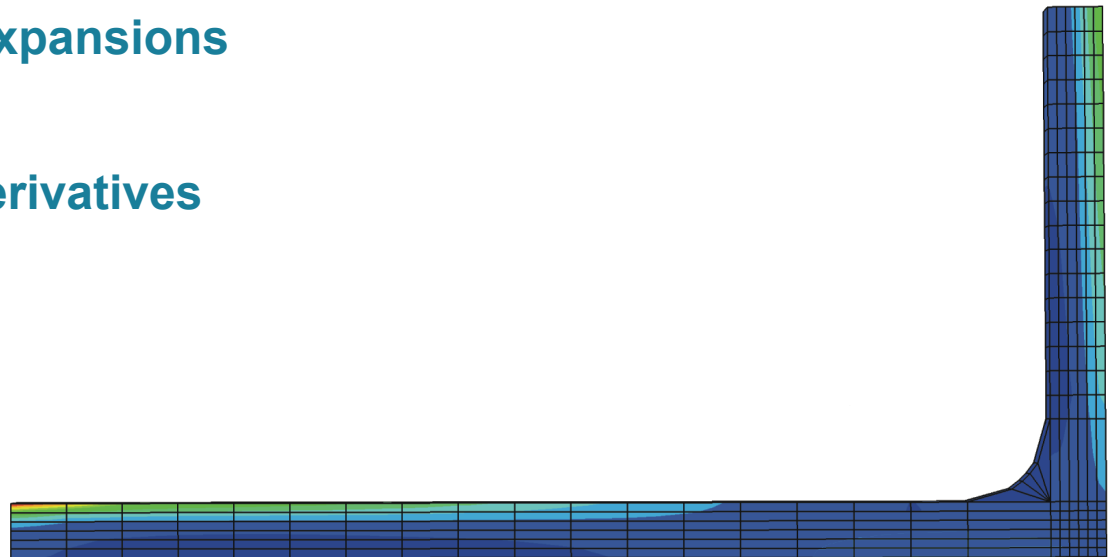


On the choice of shape functions

Shape functions:

Polynomials are usually applied for the development of shape functions (polynomials are easily differentiated analytically)

- Lagrange polynomials
complete polynomial expansions
- Serendipity polynomials
incomplete polynomial expansions
- Hermitian polynomials
polynomials including derivatives



On the choice of shape functions

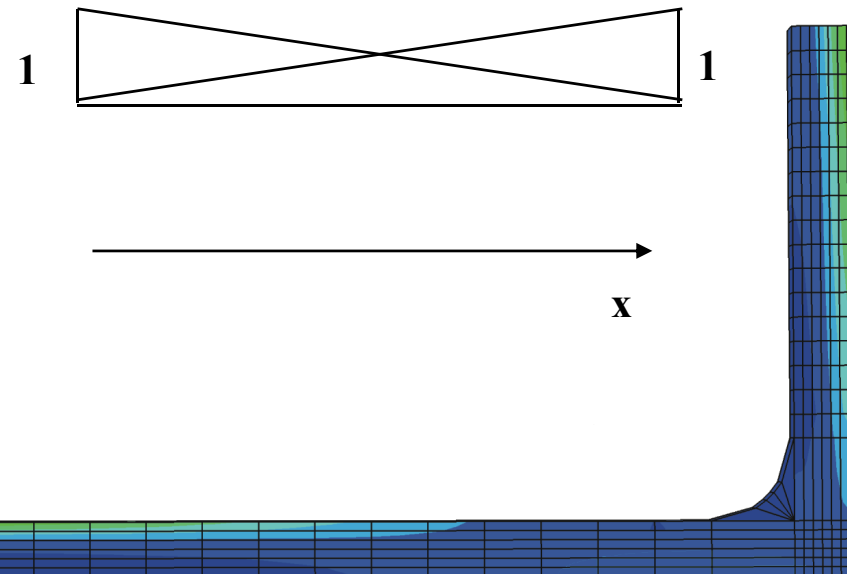
Shape functions:

Lagrange polynomials (one dimensional element):

$$H_i(x) = L_i(y) \Rightarrow u(x) = \sum_{i=1}^n L_i(x) \hat{u}_i$$

$$L_i(x) = \prod_{\substack{j=1, \\ j \neq i}}^{n+1} \frac{x - x_j}{x_i - x_j}$$

$$\begin{aligned} u(x) &= L_1(x) \hat{u}_1 + L_2(x) \hat{u}_2 \\ &= \frac{(x_2 - x)}{x_2 - x_1} \hat{u}_1 + \frac{(x - x_1)}{x_2 - x_1} \hat{u}_2 \end{aligned}$$



On the choice of shape functions

Shape functions:

Lagrange polynomials (general):

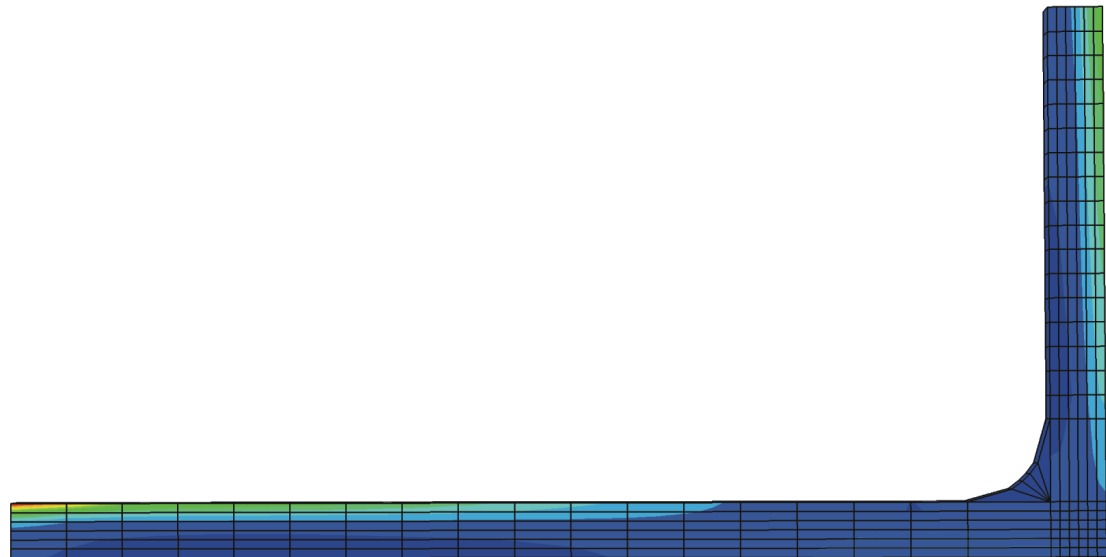
$$u(x, y, z) = \sum_{i=1}^n L_i(x, y, z) \hat{u}_i$$

$$u(x, y, z) = \sum_{i=1}^n \hat{u}_i$$

$$L_i(x_j, y_j, z_j) = 1, \quad i = j$$

$$L_i(x_j, y_j, z_j) = 0, \quad i \neq j$$

$$\sum_{i=1}^n L_i(x, y, z) = 1$$



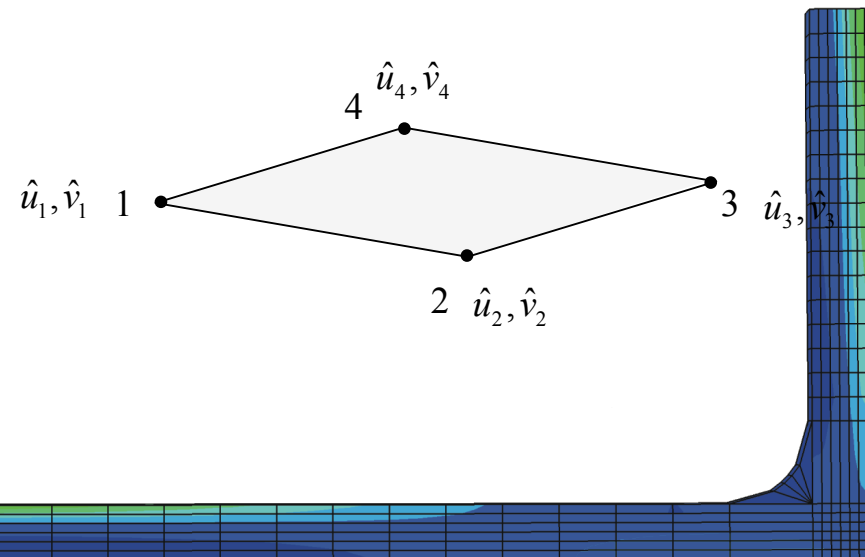
On the choice of shape functions

Shape functions:

Lagrange polynomials (four node rectangular element):

Products of two one dimensional first order Lagrange polynomials result in the **bi-linear four-node element**

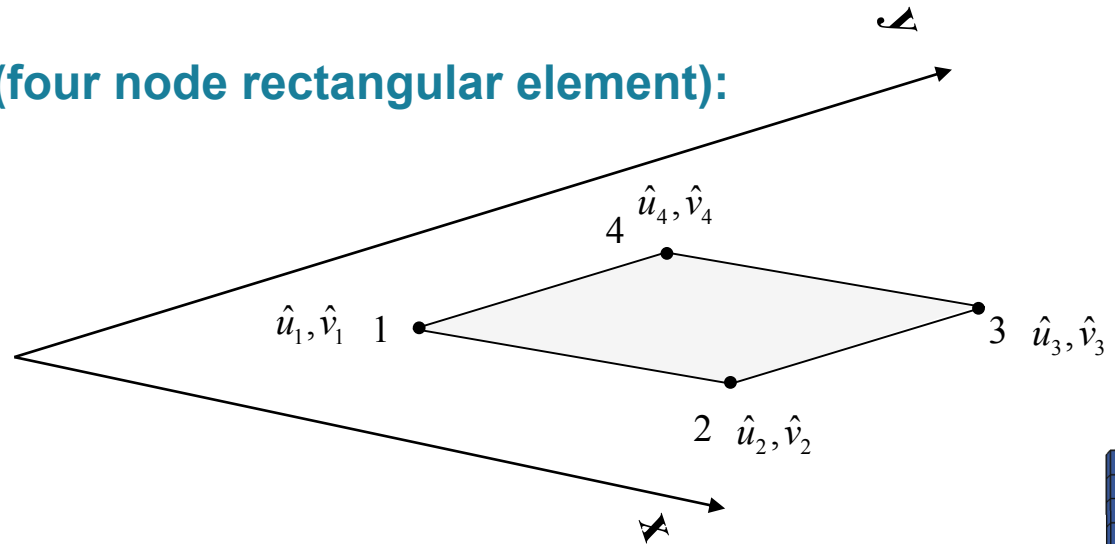
$$\begin{aligned}u(x, y) &= a_1 + a_2x + a_3y + a_4xy \\ &= (b_1 + b_2x)(b_3 + b_4y)\end{aligned}$$



On the choice of shape functions

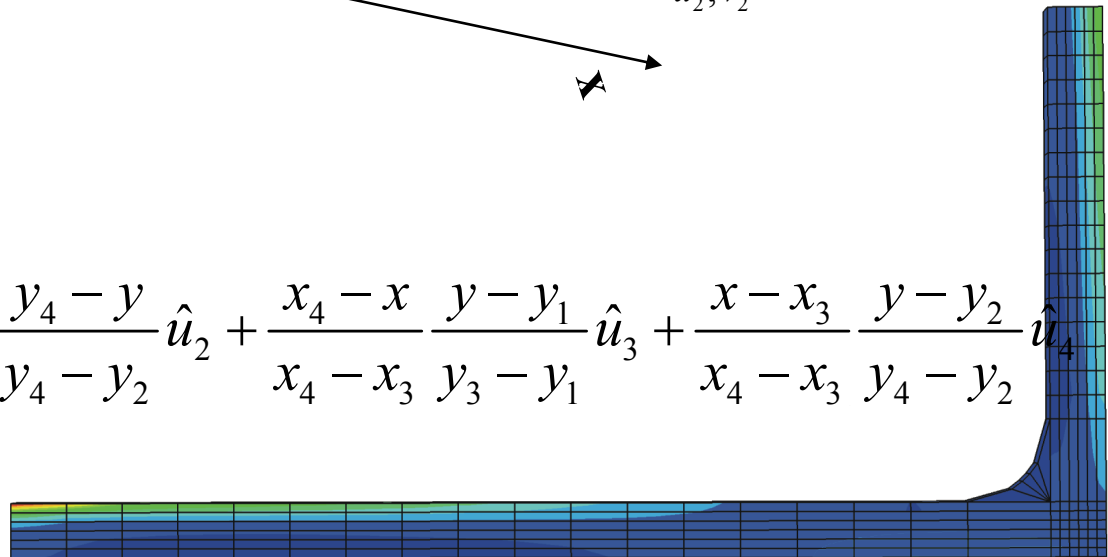
Shape functions:

Lagrange polynomials (four node rectangular element):



$$u(x, y) =$$

$$\frac{x_2 - x}{x_2 - x_1} \frac{y_3 - y}{y_3 - y_1} \hat{u}_1 + \frac{x - x_1}{x_2 - x_1} \frac{y_4 - y}{y_4 - y_2} \hat{u}_2 + \frac{x_4 - x}{x_4 - x_3} \frac{y - y_1}{y_3 - y_1} \hat{u}_3 + \frac{x - x_3}{x_4 - x_3} \frac{y - y_2}{y_4 - y_2} \hat{u}_4$$

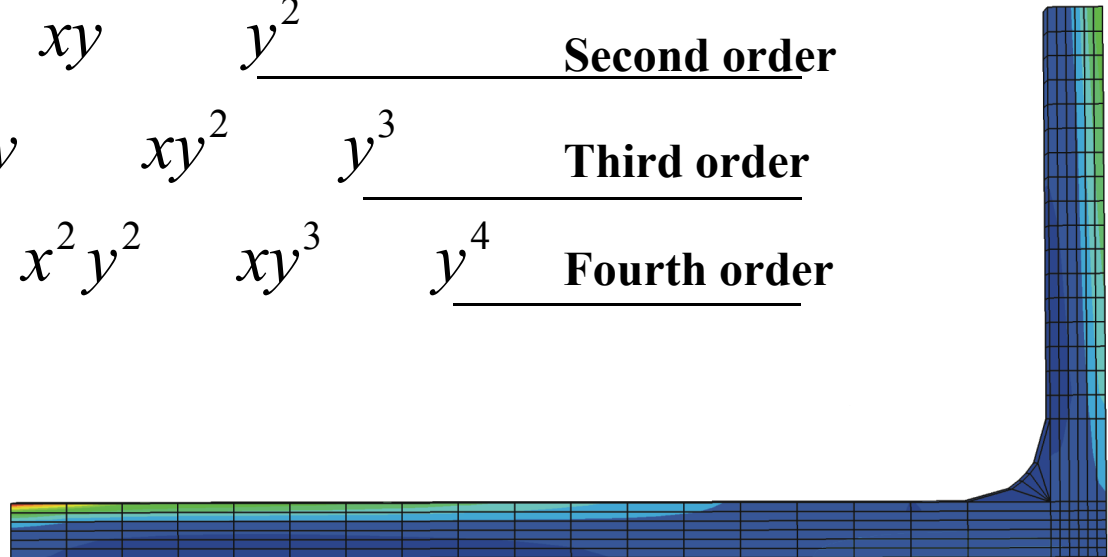


On the choice of shape functions

Shape functions:

From Pascal's triangle we can see how many nodes are required for the representation of displacement fields of any order and completeness:

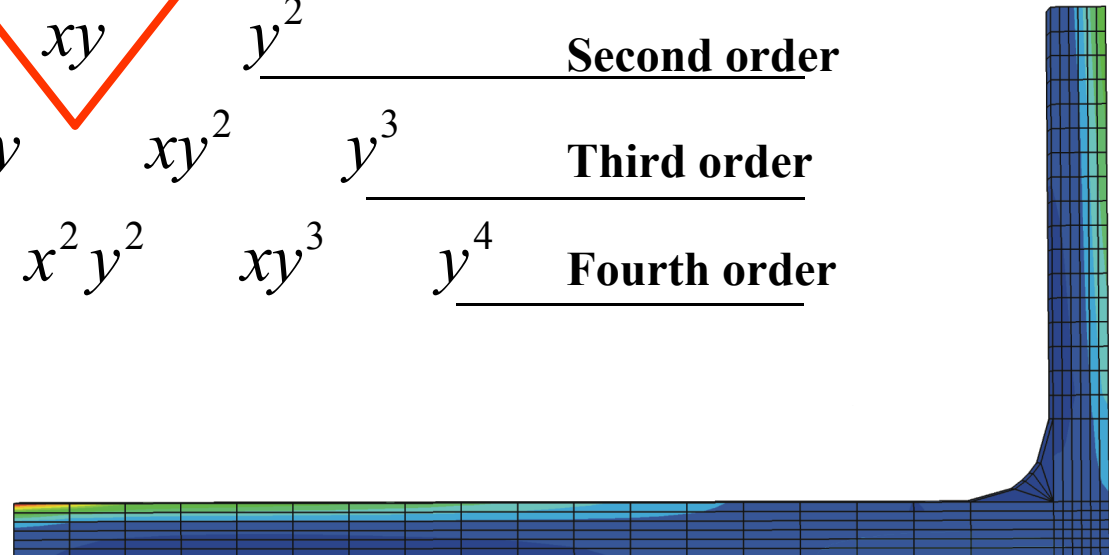
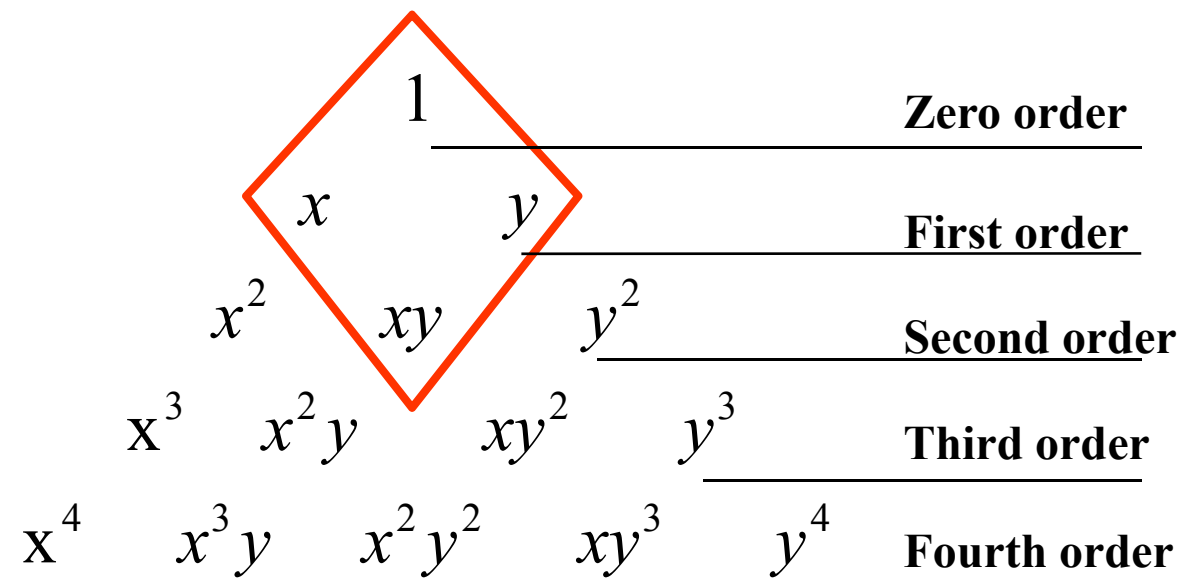
$$\begin{array}{cccccc}
 & & & & & 1 & \text{Zero order} \\
 & & & & & \hline
 & & & & x & y & \text{First order} \\
 & & & & \hline
 & & x^2 & xy & y^2 & \text{Second order} \\
 & & \hline
 x^3 & x^2y & xy^2 & y^3 & \text{Third order} \\
 & \hline
 x^4 & x^3y & x^2y^2 & xy^3 & y^4 & \text{Fourth order} \\
 & \hline
 \end{array}$$



On the choice of shape functions

Shape functions:

Products of Lagrange polynomials (bi-linear four node rectangular)

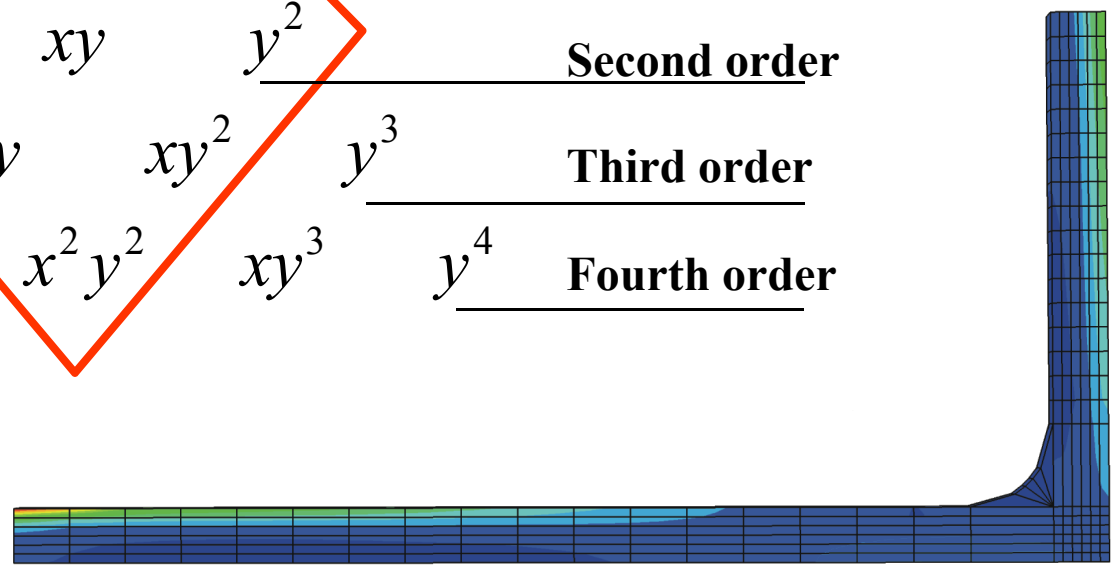
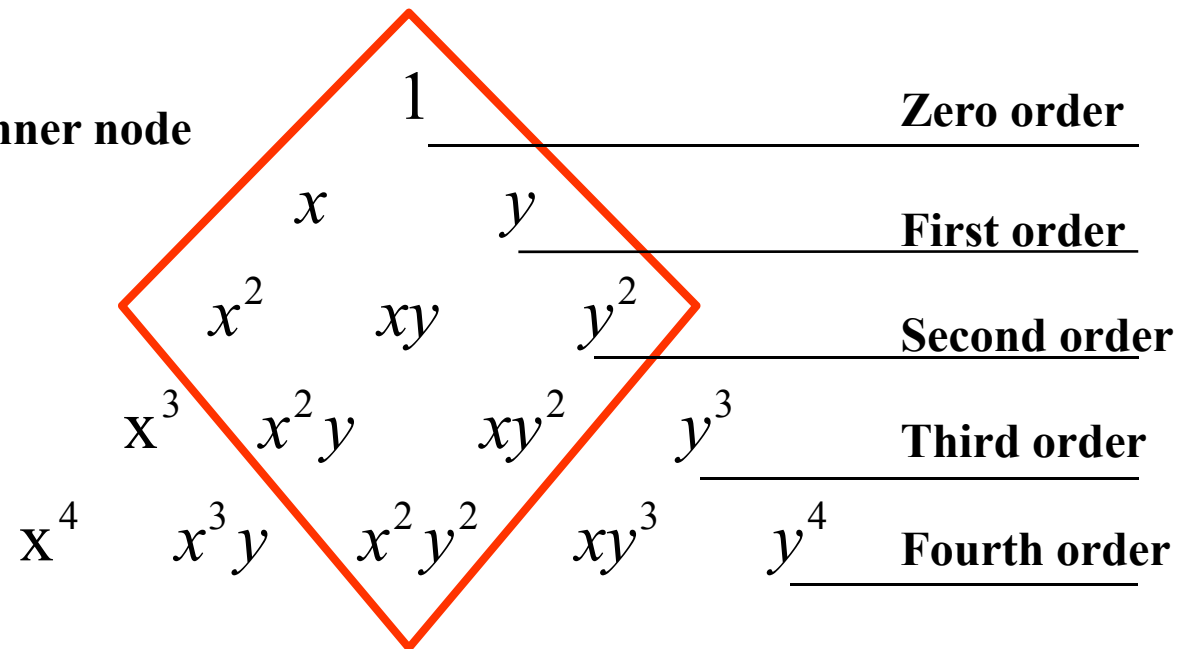


On the choice of shape functions

Shape functions:

Products of Lagrange polynomials (quadratic nine-node rectangular)

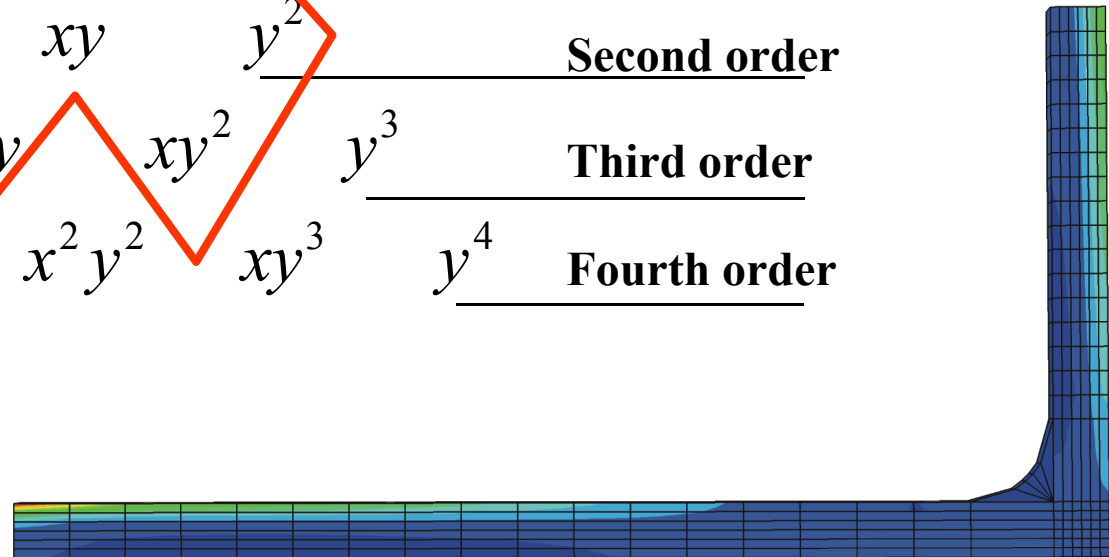
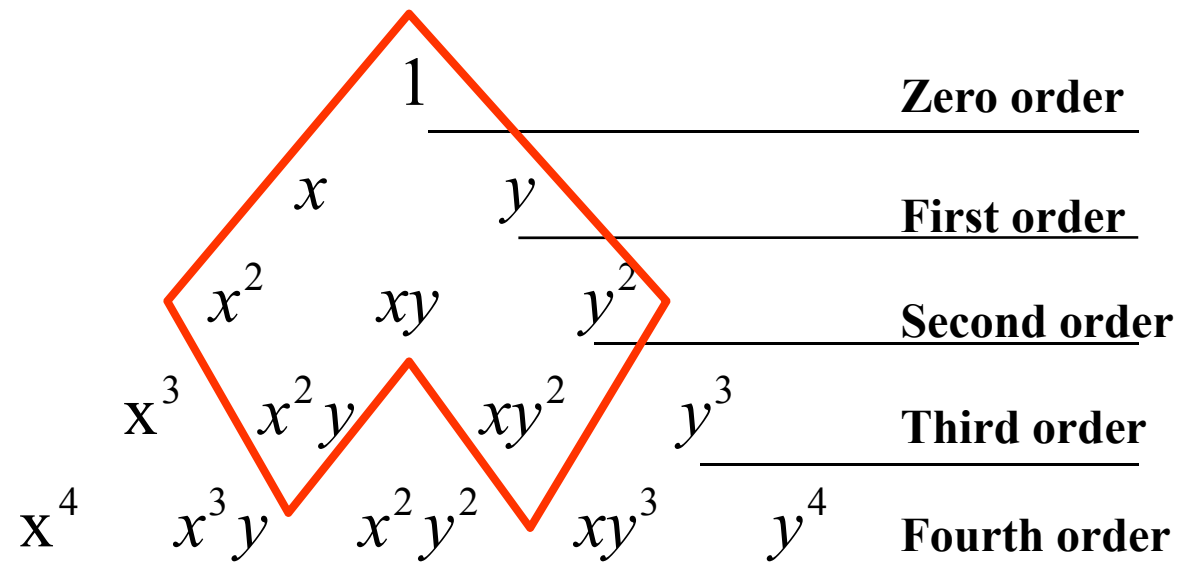
This requires an inner node
a difficulty !



On the choice of shape functions

Shape functions:

Serendipity shape functions are constructed by incomplete polynomials – avoiding inner nodes

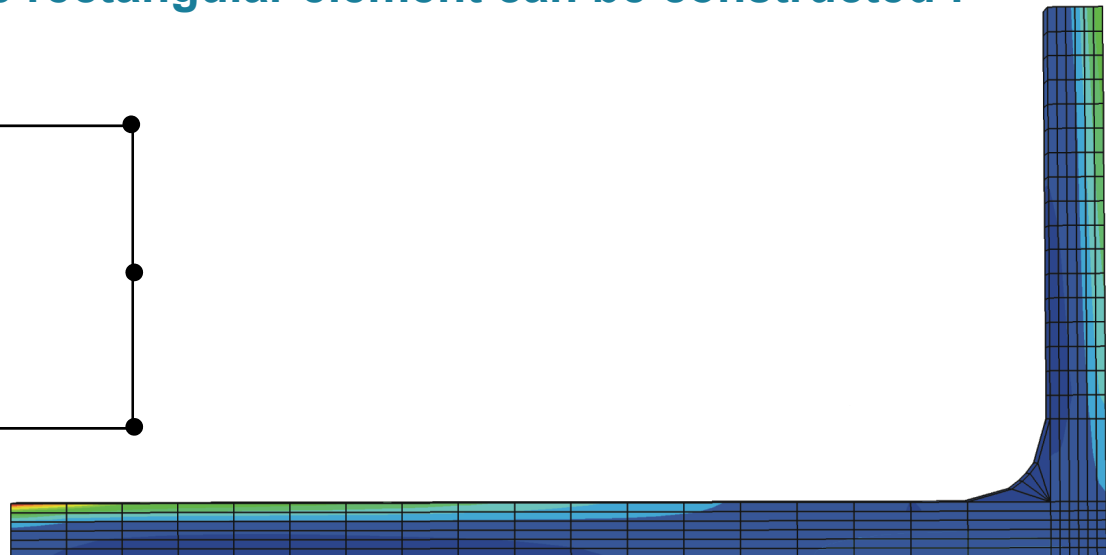
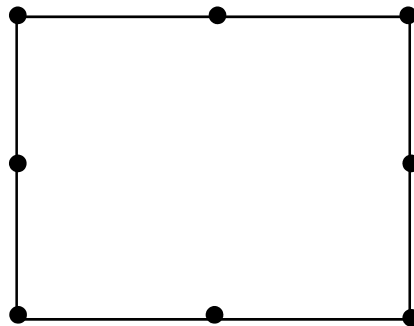


On the choice of shape functions

Shape functions:

Whereas difficulties may arise (inner nodes) when aiming to develop quadratic shape functions for rectangular elements using Lagrange polynomials the shape functions developed by incomplete polynomials (**serendipity shape functions**) – less terms necessitates less nodes !

An bi-quadratic eight node rectangular element can be constructed !



On the choice of shape functions

Shape functions:

Hermitian shape functions relate not only the displacements at nodes to displacements within the elements but also the first order derivatives

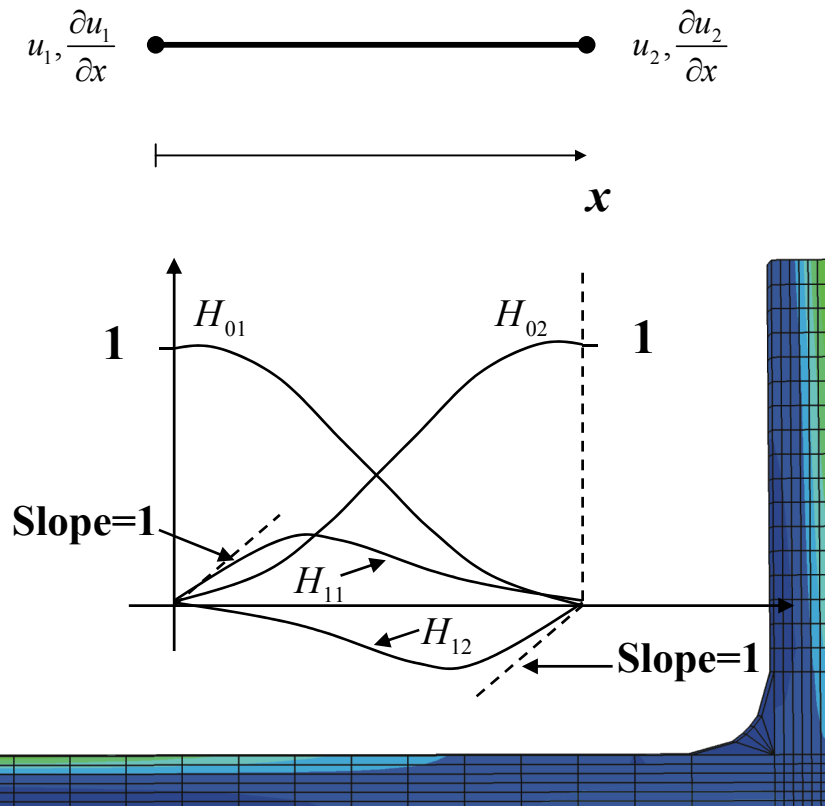
$$u(x) = \sum_{i=1}^2 \left(H_{0i}(x) \hat{u}_i + H_{1i}(x) \frac{\partial \hat{u}_i}{\partial x} \right)$$

$H_{0i}(x) = 1$, and zero at the other node

$H'_{0i}(x) = 0$ at both nodes

$H_{1i}(x) = 0$, at both nodes

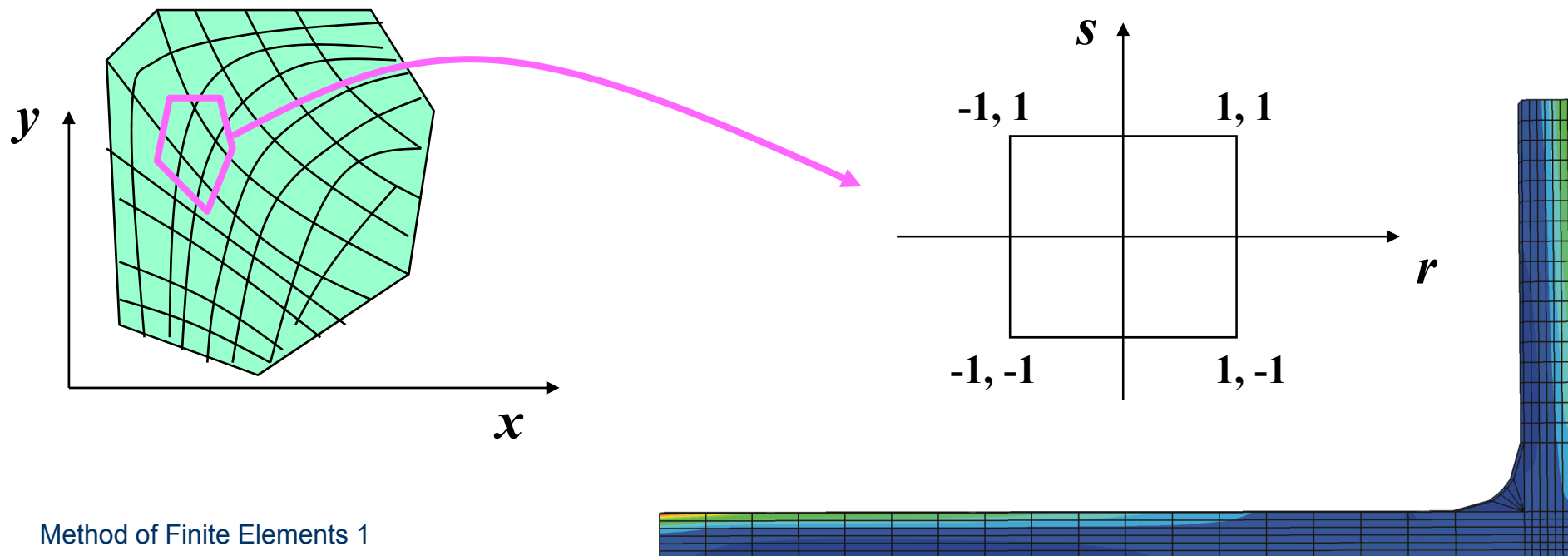
$H'_{1i}(x) = 1$, and zero at the other node



Natural coordinates

Shape functions – Natural coordinates:

As we have seen we are able to establish shape functions in global or local coordinate systems as we please. However, for the purpose of standardizing the process of developing the element matrixes it is convenient to introduce the so-called natural coordinate system.



Natural coordinates

Shape functions – Natural coordinates:

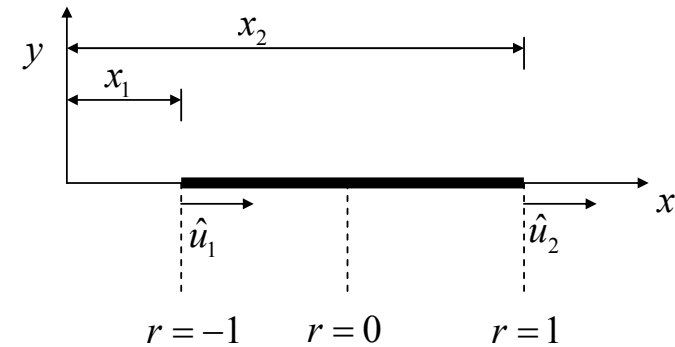
Let us consider the simple bar element

The relation between the x-coordinate and the r-coordinate is given as:

$$x = \frac{1}{2}(1-r)\hat{x}_1 + \frac{1}{2}(1+r)\hat{x}_2$$

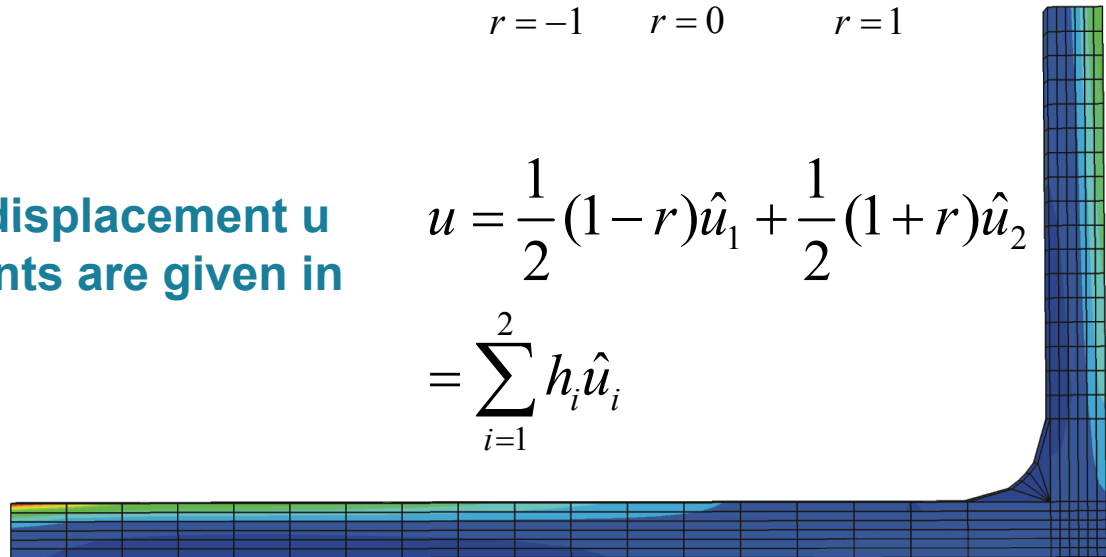
$$= \sum_{i=1}^2 h_i \hat{x}_i$$

The relation between the displacement u and the nodal displacements are given in the same way:



$$u = \frac{1}{2}(1-r)\hat{u}_1 + \frac{1}{2}(1+r)\hat{u}_2$$

$$= \sum_{i=1}^2 h_i \hat{u}_i$$



Natural coordinates

Shape functions – Natural coordinates:

Let us consider the simple bar element

We need to be able to establish the strains – meaning we need to be able to take the derivatives of the displacement field in regard to the x-coordinate

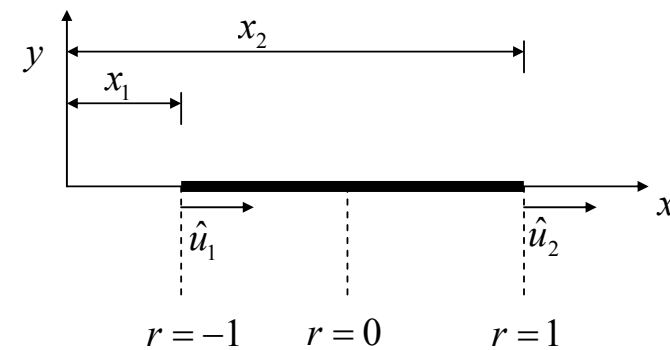
$$\varepsilon = \frac{du}{dx} = \frac{du}{dr} \frac{dr}{dx}$$

$$\frac{du}{dr} = \frac{d}{dr} \left(\frac{1}{2} (1-r) \hat{u}_1 + \frac{1}{2} (1+r) \hat{u}_2 \right) = \frac{1}{2} (\hat{u}_2 - \hat{u}_1)$$

$$\frac{dx}{dr} = \frac{d}{dr} \left(\frac{1}{2} (1-r) x_1 + \frac{1}{2} (1+r) x_2 \right) = \frac{1}{2} (x_2 - x_1)$$

$$\Downarrow$$

$$\frac{du}{dx} = \frac{(\hat{u}_2 - \hat{u}_1)}{(x_2 - x_1)} = \frac{(\hat{u}_2 - \hat{u}_1)}{L}$$



Natural coordinates

Shape functions – Natural coordinates:

Let us consider the simple bar element

The strain-displacement matrix then becomes:

$$\mathbf{B} = \frac{1}{L} \begin{bmatrix} -1 & 1 \end{bmatrix}$$

and the stiffness matrix is calculated as:

$$\mathbf{K} = \frac{AE}{L^2} \int_{-1}^1 \begin{bmatrix} -1 \\ 1 \end{bmatrix} \begin{bmatrix} -1 & 1 \end{bmatrix} \mathbf{J} dr, \quad \mathbf{J} = \frac{dx}{dr} = \frac{L}{2}$$

⇓

$$\mathbf{K} = \frac{AE}{L} \begin{bmatrix} 1 & -1 \\ -1 & 1 \end{bmatrix}$$

