FINITE ELEMENT METHOD II

Autumn 2015

Lectures (15h):

1. Accuracy, error estimation and adaptive remeshing
2. Heat flow and thermal stresses in FEM
3. Introduction to structural dynamics, free vibrations
4. Nonlinear problems in mechanics of structures - basic numerical techniques
5. Orthotropic materials and composite structures
6. Parametric modeling and design optimization

Computer lab (15h):

Modeling simple problems of: thermal stresses, contact mechanics, plasticity and residual stresses, free vibrations, buckling, parametric modeling

References:

[1] Lecture notes from the web site: http://meil.pw.edu.pl/zwmik/ZWMiK/Dla-studentow2/Finite-Element-Method-II

[2] Moaveni S.: *Finite element analysis. Theory and applications with ANSYS*. Paerson Education, 2015.

[3] Kleiber M. (red.): *Komputerowe metody mechaniki ciał stałych*, seria Mechanika Techniczna XI, Warszawa PWN 1995.

[4] Xiaolin Chen, Yijun Liuv: *Finite Element Modeling and Simulation with ANSYS.* Workbench, CRC Press 2014

[5] Huebner K. H., Dewhirst D. L., Smith D.E., Byrom T. G.: *The finite element method for engineers*, J. Wiley & Sons, Inc., 2001.

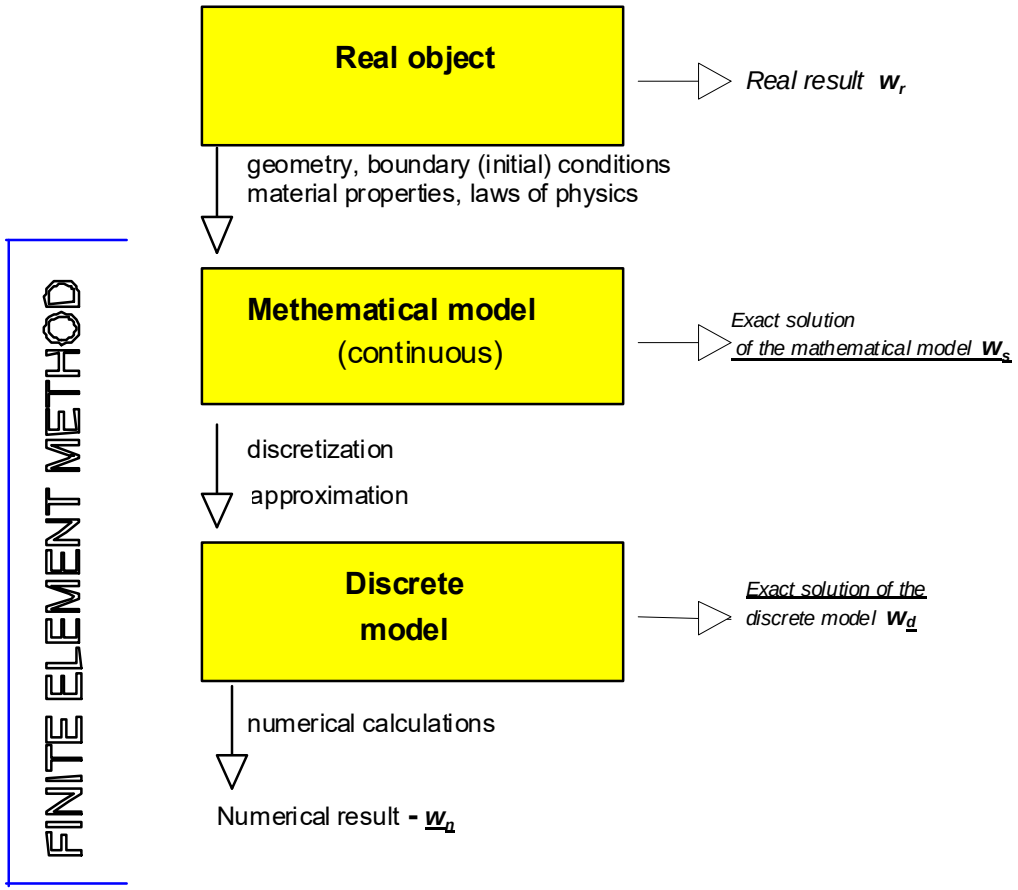
[6] Zienkiewicz O.C., Taylor R.: *The Finite Element Method* - different publishers and editions

[7] Krzesiński G., Zagrajek T., Marek P., Borkowski P.*: MES w mechanice materiałów i konstrukcji. Rozwiązywanie wybranych zagadnień za pomocą programu ANSYS*, Of. Wyd.PW 2015

[8] Bijak-Żochowski M., Jaworski A., Krzesiński G., Zagrajek T.: *Mechanika Materiałów i Konstrukcji*, Tom 2, Warszawa, Of. Wyd. PW, 2014

Assessment based on the final test and the results of computer lab work

1. ACCURACY OF FE ANALYSIS. ERROR ESTIMATION AND ADAPTIVE REMESHING



$\varepsilon_s = w_s - w_r$ – **modeling error**,

$\varepsilon_d = w_d - w_s$ – **discretization error**

$\varepsilon_n = w_n - w_d$ – **numerical error**

Total error:

$\varepsilon_c = \varepsilon_s + \varepsilon_d + \varepsilon_n = w_n - w_r$

The most effective FEM analysis:

$\varepsilon_s \sim \varepsilon_d \sim \varepsilon_n$

Approximate methods – flow chart

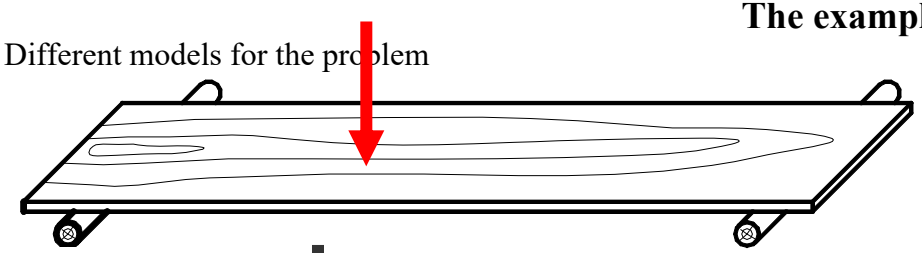
(valid also for other methods e.g. Boundary Element Method - BEM and Finite Differences Method - FDM)

Modeling error ε_s

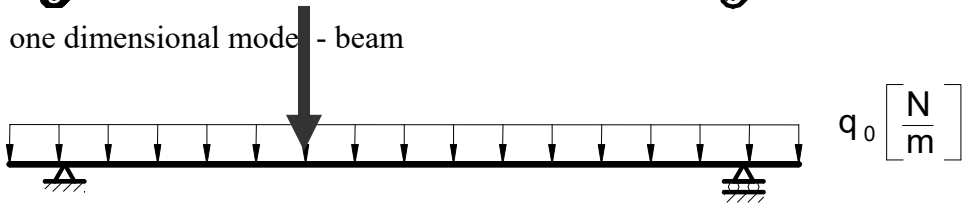
Depends on the accuracy of the available information about the problem and knowledge of the analyst (1D – 2D – 3D models, linear, nonlinear, assumed simplifications, reliable information concerning material properties, loads.)

The example – wooden board

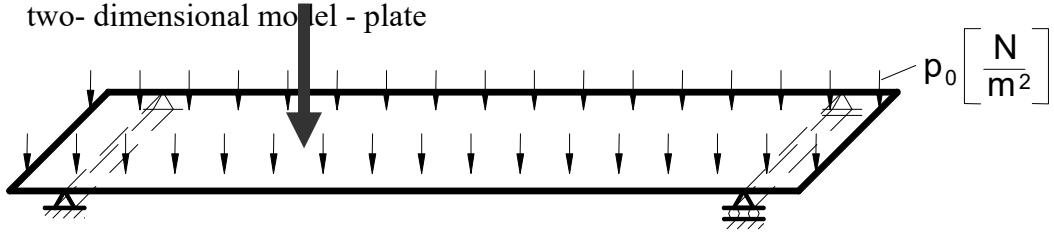
Different models for the problem



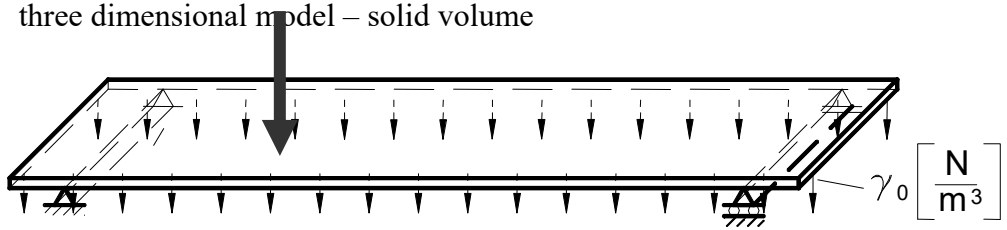
one dimensional model - beam



two- dimensional model - plate



three dimensional model – solid volume

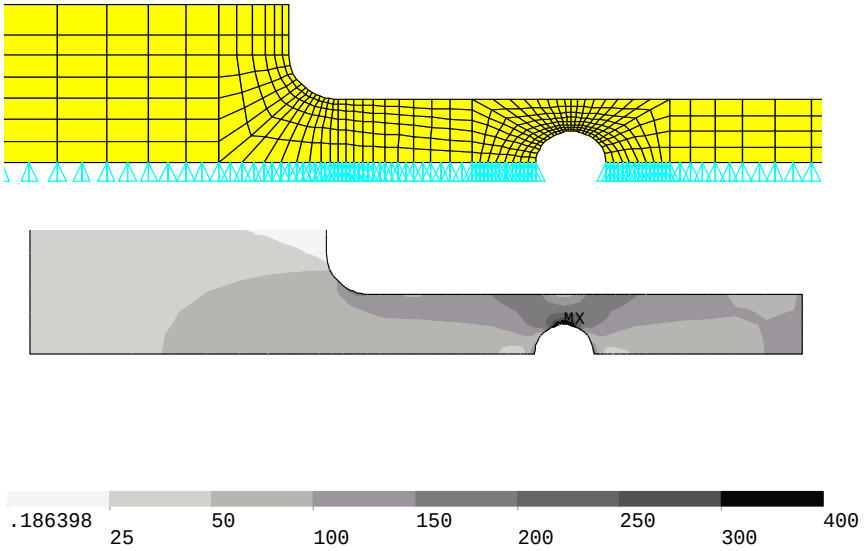


Mathematical model:

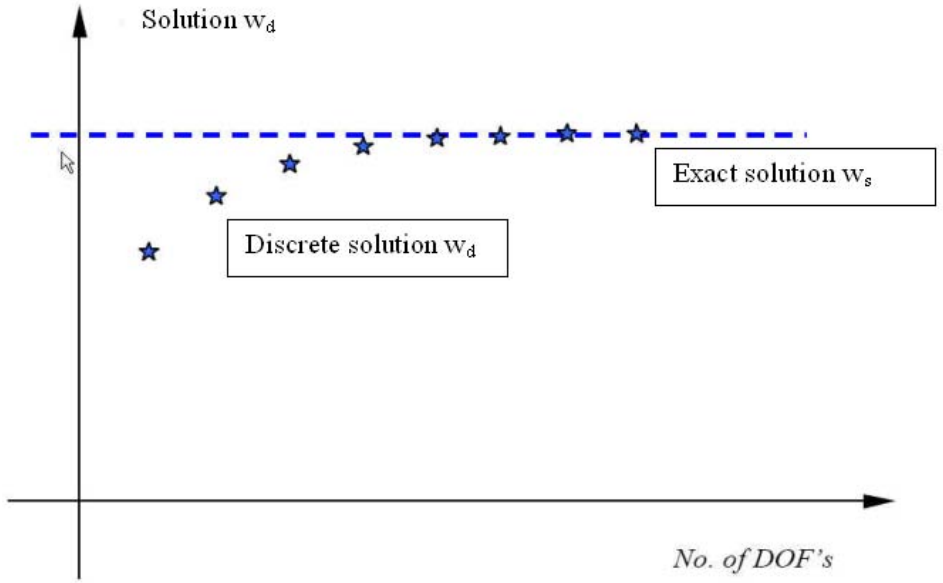
- 1,2, 3 dimensional
- properties of wood: isotropic-orthotropic, moist-dry, homogenous-inhomogeneous, rate dependent-rate-independent)
- contact interaction model (friction)
- loadings

Discretization error ε_d

Depends on the mesh density, types of the elements – shape functions, the shapes of the finite elements



FE mesh and von Mises stress distribution



Discrete solution versus exact solution of the continuous problem

There are some mathematical convergence requirements in FEA concerning the mesh, shape functions, and rules of FE model building.

Numerical error ε_n

$$[K]\{q\} = \{F\},$$

$$[K+\delta K]\{q+\delta q\} = \{F+\delta F\}. \qquad \rightarrow \qquad \frac{\|\delta q\|}{\|q\|} \le \|[K]\|\,\|[K]^{-1}\|\left(\frac{\|\{\delta F\}\|}{\|\{F\}\|} + \frac{\|[\delta K]\|}{\|[K]\|}\right),$$

condition number of the matrix K

$$cond([K]) = \|[K]\|\,\|[K]^{-1}\|.$$

Norm of a matrix (vector)– a measure of magnitude

L_2- Euclidean norms

Vector norm $\|\{q\}\| = \left(\sum_i (q_i)^2\right)^{\frac{1}{2}},$

Matrix norm $\|[K]\| = \left(\sum_j \sum_i (k_{ij})^2\right)^{\frac{1}{2}}$ (matrix norm induced by the vector norm)

Max norms ($L_\sim$ norms)

$$\|\{q\}\| = \max_i |q_i|, \qquad \|[K]\| = \max_i \left(\sum_j |k_{ij}|\right).$$

A problem with a small condition number is said to be **well-conditioned**, while a problem with a high condition number is said to be **ill-conditioned**.

cond(K) $| \ge 1$ cond(K) ≈ 1 - problem well-conditioned cond(K)» 1 - problem ill-conditioned

Reasons of ill-conditioning of the problems in FE stress analysis – great differences between stiffness of FE elements, unstable boundary conditions

Wskaźnik uwarunkowania macierzy [edytuj | edytuj kod]

Zasugerowano, aby ta sekcja została **przeniesiona** do nowego artykułu nazwanego Wskaźnik uwarunkowania macierzy. (dyskusja)

Wskaźnik uwarunkowania macierzy A w równaniu $Ax = b$ jest charakterystyczną własnością macierzy informującą o tym, jakie wzmocnienie będzie miała zmiana normy macierzy A na normę rozwiązania x.

Wskaźnik uwarunkowania macierzy definiuje się bardziej precyzyjnie jako maksymalny stosunek błędu względnego wektora rozwiązania x do błędu względnego b.

Załóżmy, że e jest błędem b. Stąd błąd w rozwiązaniu $A^{-1}b$ wynosi $A^{-1}e$. Stąd stosunek relatywnego błędu rozwiązania do relatywnego błędu w b wynosi:

$$\frac{\|A^{-1}e\|/\|A^{-1}b\|}{\|e\|/\|b\|}.$$

Można to przekształcić do:

$$(\|A^{-1}e\|/\|e\|) \cdot (\|b\|/\|A^{-1}b\|).$$

Maksymalna wartość (dla niezerowych b i e) będzie iloczynem dwóch norm (definiowanych w różny sposób, np. często jako normę traktuje się maksymalną sumę wartości bezwzględnych wierszy):

$$\kappa(A) = \|A^{-1}\| \cdot \|A\|.$$

Definicja ta jest taka sama dla każdej zwartej normy. Liczba ta pojawia się tak często w algebrze liniowej, że nadano jej nazwę **wskaźnika uwarunkowania macierzy**.

Zastosowania [edytuj | edytuj kod]

Wskaźnik uwarunkowania macierzy pozwala na oszacowanie, z jaką (maksymalnie) dokładnością (do ilu miejsc po przecinku) możemy podać wynik. Dokładność jest zależna od iloczynu epsilonu maszynowego i wskaźnika uwarunkowania. Załóżmy dla przykładu, że mamy macierz $\mathbf{A}$:

$$\mathbf{A} = \begin{bmatrix} 1 & 2 \\ 3 & 4 \end{bmatrix},$$

$$\mathbf{A^{-1}} = \begin{bmatrix} -2 & 1 \\ 1{,}5 & -0{,}5 \end{bmatrix}.$$

Stosując tak zdefiniowaną normę: $\|A\| = max_{1<i<n} \sum_{j=1}^{n} |a_{ij}|$ możemy obliczyć wskaźnik uwarunkowania $\kappa(A) = \|A^{-1}\| \cdot \|A\| = 21.$

Jako inny przykład rozpatrzmy prosty układ równań typu $\mathbf{Ax} = \mathbf{b}$. Jeśli do naszych obliczeń wybierzemy macierz o wysokim wskaźniku uwarunkowania, np.:

$$\mathbf{A} = \begin{bmatrix} 1 & 2 \\ 2 & 3{,}999 \end{bmatrix}; \mathbf{b} = \begin{bmatrix} 4 \\ 7{,}999 \end{bmatrix}; cond(\mathbf{A}) = 35\ 988,$$

to otrzymane rozwiązanie jest niestabilne. Oznacza to, że mała zmiana wartości współczynników może znacząco wpłynąć na wynik.

W podanym wyżej przypadku rozwiązanie wynosi $\mathbf{x} = \begin{bmatrix} 2 \\ 1 \end{bmatrix}$. Jeśli zmodyfikujemy następująco wektor $\mathbf{b} = \begin{bmatrix} 4 \\ 8 \end{bmatrix}$, to otrzymamy rozwiązanie $\mathbf{x} = \begin{bmatrix} 4 \\ 0 \end{bmatrix}$.

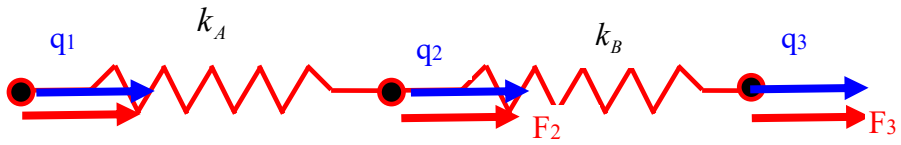
W przypadku macierzy dobrze uwarunkowanej np.:

$$\mathbf{A} = \begin{bmatrix} 1 & 2 \\ 2 & 3 \end{bmatrix}; \mathbf{b} = \begin{bmatrix} 4 \\ 7{,}999 \end{bmatrix}; cond(\mathbf{A}) = 25,$$

rozwiązanie wynosi $\mathbf{x} = \begin{bmatrix} 3{,}998 \\ 0{,}001 \end{bmatrix}$.

Jeśli zmodyfikujemy następująco wektor $\mathbf{b} = \begin{bmatrix} 4 \\ 8 \end{bmatrix}$, to otrzymamy rozwiązanie $\mathbf{x} = \begin{bmatrix} 4 \\ 0 \end{bmatrix}$, które jest zbliżone do poprzedniego.

The example (ill conditioned problem):



$$[k]_e = \underset{2\times 2}{\begin{bmatrix} k & -k \\ -k & k \end{bmatrix}} \qquad \begin{bmatrix} k_A & -k_A & \\ -k_A & k_A + k_B & -k_B \\ & -k_B & k_B \end{bmatrix} \begin{Bmatrix} q_1 \\ q_2 \\ q_3 \end{Bmatrix} = \begin{Bmatrix} F_1 \\ F_2 \\ F_3 \end{Bmatrix}$$

$$q_1 = 0 \qquad \begin{bmatrix} k_A + k_B & -k_B \\ -k_B & k_B \end{bmatrix} \begin{Bmatrix} q_2 \\ q_3 \end{Bmatrix} = \begin{Bmatrix} F_2 \\ F_3 \end{Bmatrix}$$

Let's assume: $k_A = 1 \qquad k_B = 1000$

$$\begin{Bmatrix} F_2 \\ F_3 \end{Bmatrix} = \begin{Bmatrix} 1 \\ -1 \end{Bmatrix}$$

The solution:

$$\begin{Bmatrix} q_2 \\ q_3 \end{Bmatrix} = \begin{bmatrix} \frac{1}{k_A} & \frac{1}{k_A} \\ \frac{1}{k_A} & \frac{1}{k_A} + \frac{1}{k_B} \end{bmatrix} \begin{Bmatrix} F_2 \\ F_3 \end{Bmatrix}$$

The result:

$$\begin{Bmatrix} q_2 \\ q_3 \end{Bmatrix} = \begin{Bmatrix} 0 \\ -0.001 \end{Bmatrix}$$

The example of the very small perturbation

For changed force vector : $\begin{Bmatrix} F_2 \\ F_3 \end{Bmatrix} = \begin{Bmatrix} 0.999 \\ -1.0 \end{Bmatrix}$

$$\begin{Bmatrix} q_2 \\ q_3 \end{Bmatrix} = \begin{Bmatrix} -0.001 \\ -0.002 \end{Bmatrix}$$

$$\{\delta F\} = \begin{Bmatrix} \delta F_2 \\ \delta F_3 \end{Bmatrix} = \begin{Bmatrix} -0.001 \\ 0 \end{Bmatrix} \qquad \{\delta q\} = \begin{Bmatrix} \delta q_2 \\ \delta q_3 \end{Bmatrix} = \begin{Bmatrix} -0.001 \\ -0.001 \end{Bmatrix}$$

$$\frac{\|\delta F\|}{\|F\|} = 0.707 \cdot 10^{-3} \quad \Rightarrow \quad \frac{\|\delta q\|}{\|q\|} = 1.414$$

$$cond(K) \approx 4 \bullet 10^3$$

First equation

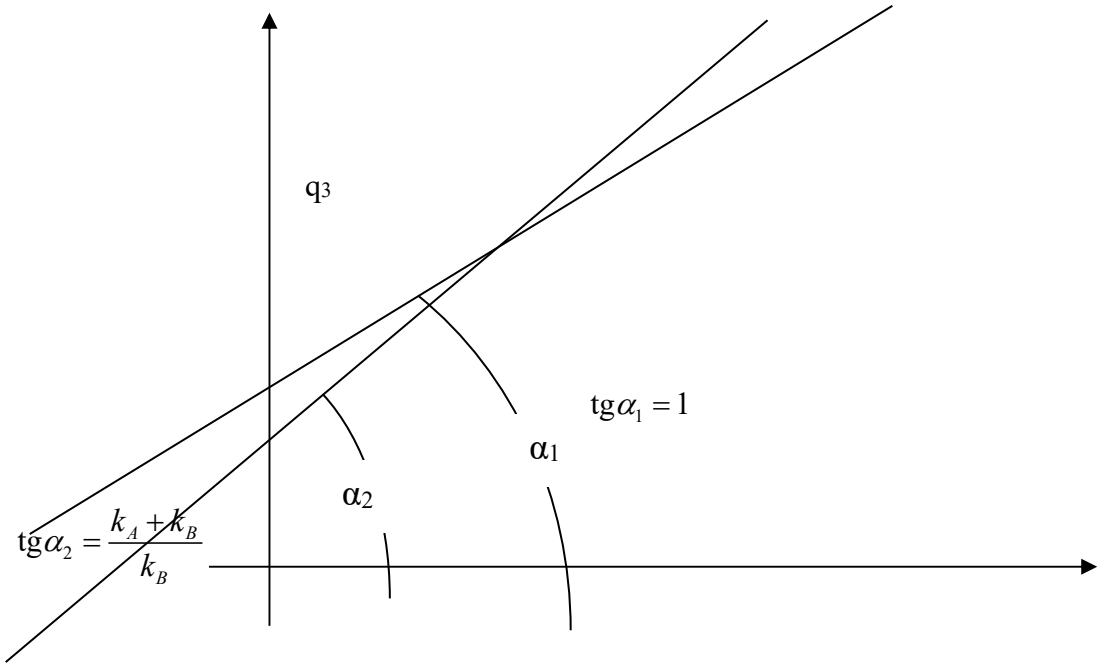
$$q_3 = \frac{k_A + k_B}{k_B} \cdot q_2 - \frac{F_2}{k_B}$$

Second equation

$$q_3 = q_2 + \frac{F_3}{k_B}$$

System ill-conditioned if

$$\frac{k_A + k_B}{k_B} \to 1 \quad \text{or} \quad \frac{k_A}{k_B} \to 0$$



Sensitivity to the slop change

$K_A << K_B$ **system ill conditioned**

Round–off error

As a general rule, if the condition number cond(K) = 10^k, then you may lose up to maximum k digits of accuracy during the solution of the system of linear equations. However, the condition number does not give the exact value of the maximum inaccuracy that may occur in the algorithm.

$$r \geq p - \log_{10}\left(cond\left([K]\right)\right)$$

p – number of significant digits in the computer representation of numbers

r – number of significant digits of the result

In FE models cond(K) reaches 10^8

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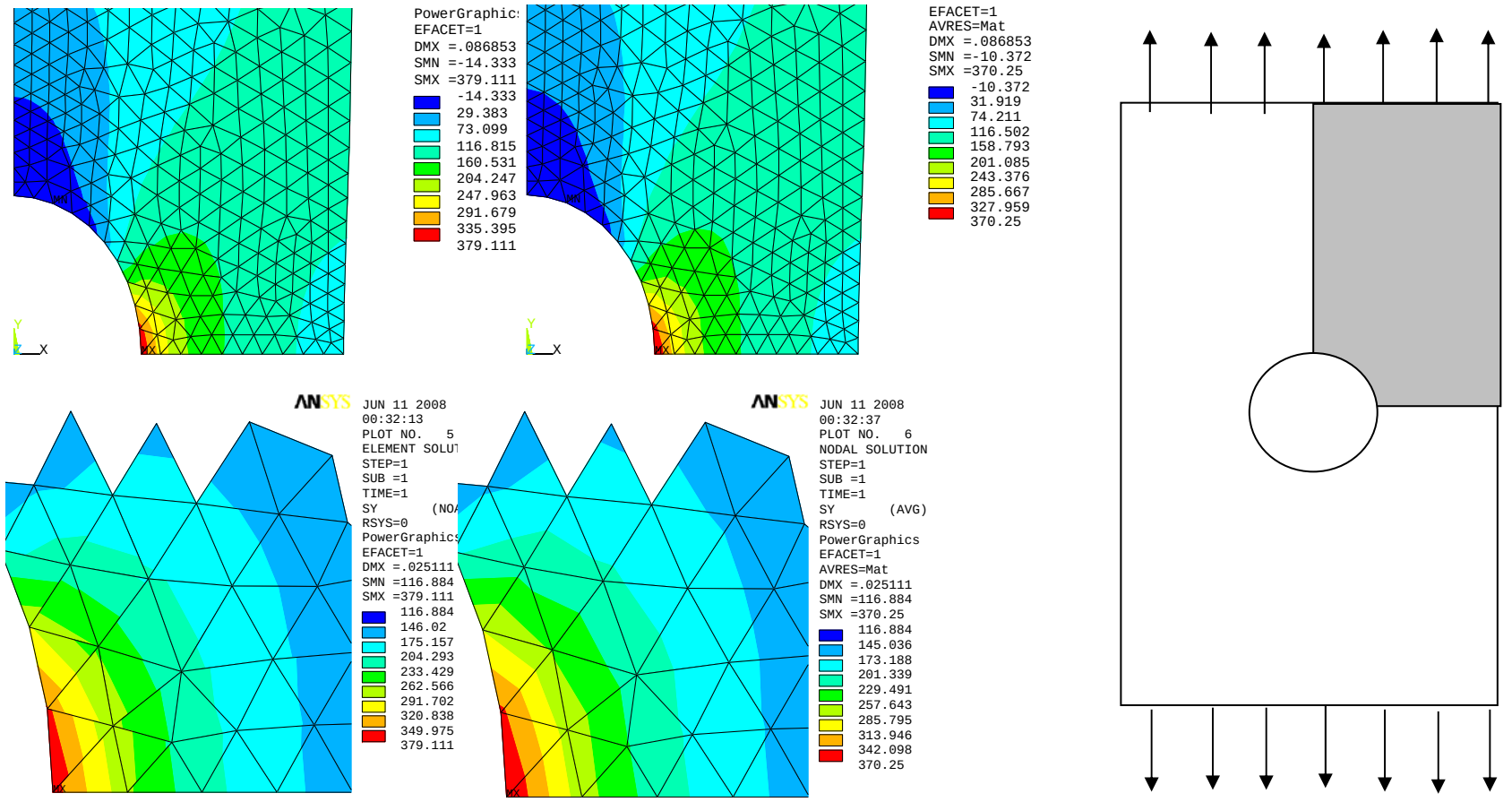
*** SYSTEM INFORMATION MESSAGE 4159 (DFMSA)
    THE DECOMPOSITION OF KLL      YIELDS A MAXIMUM MATRIX-TO-FACTOR-DIAGONAL RATIO OF    2.594055E+04
*** USER INFORMATION MESSAGE 3035 (SOLVER)
                              FOR DATA BLOCK   KLR
   SUPPORT PT.NO.               EPSILON              STRAIN   ENERGY     EPSILONS LARGER THAN 0.001 ARE FLAGGED WITH ASTERISKS
                1          1.2207031E-16          6.1035156E-05
```

A posteriori error approximation techniques

Element and nodal solution in FE program postprocessors (PLESOL, PLNSOL in ANSYS)

FE solution provides the continuous displacement field (from element to element), and the discontinuous stress field.

To obtain smooth stress distribution, the averaging of the stresses in the nodes is performed *(nodal stresses*).



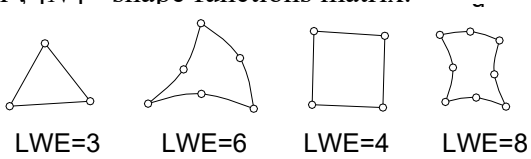
Rectangular plate with a hole under tension. The model of the quarter of the structure. The stress component σ_y
Discontinuous „element solution" (left) and averaged continuous " nodal solution" (right). Six-node triangular plane elements

Basic relations between displacements, strains, stresses and strain energy within finite elements (the relations discussed during FEM I lectures)

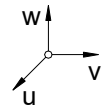
Displacement field over the element is interpolated from the nodal displacements:

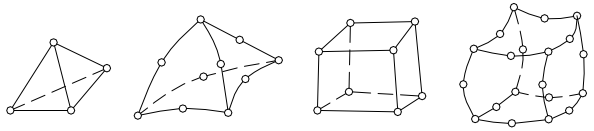
$\{u\} = [N(x,y,z)]\ \{q\}_e$, where $\{q\}_e$ - nodal displacements vector , $[N]$ - shape functions matrix.

For example for the simplest triangular element with 3 nodes and 6 DOF the relati



$$\begin{Bmatrix} u(x,y) \\ \upsilon(x,y) \end{Bmatrix} = \begin{bmatrix} N_1(x,y) & 0 & N_2(x,y) & 0 & N_3(x,y) & 0 \\ 0 & N_1(x,y) & 0 & N_2(x,y) & 0 & N_3(x,y) \end{bmatrix} \begin{Bmatrix} u_1 \\ \upsilon_1 \\ u_2 \\ \upsilon_2 \\ u_3 \\ \upsilon_3 \end{Bmatrix},$$





where N_i are the linear functions

Shape functions N_{ij} are usually polynomials defined in local (element) coordinate systems.

Displacements, strains and stresses within each element are defined as the functions of the nodal displacements

$\{u\} = [N]\{q\}_e$,

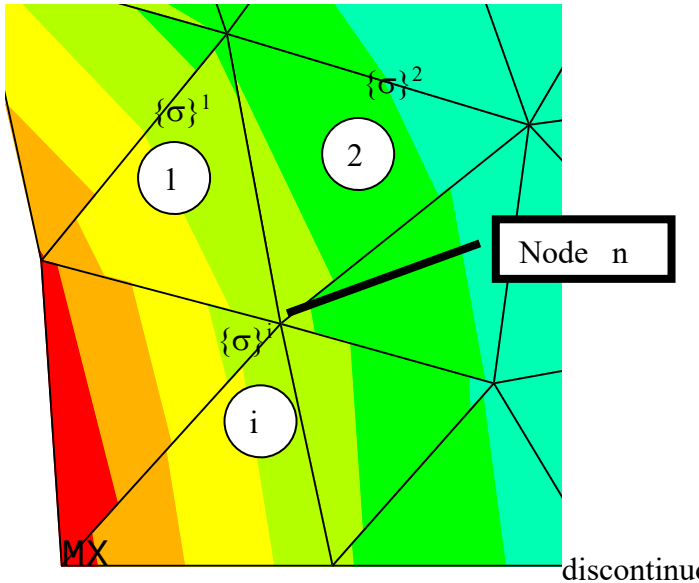
$\{\varepsilon\} = [R]\{u\} = [R][N]\{q\}_e = [B]\{q\}_e$, [B] – strain-displacement matrix, $[R]$ -gradient matrix

$\{\sigma\} = [D]\{\varepsilon\} = [D][B]\{q\}_e$.

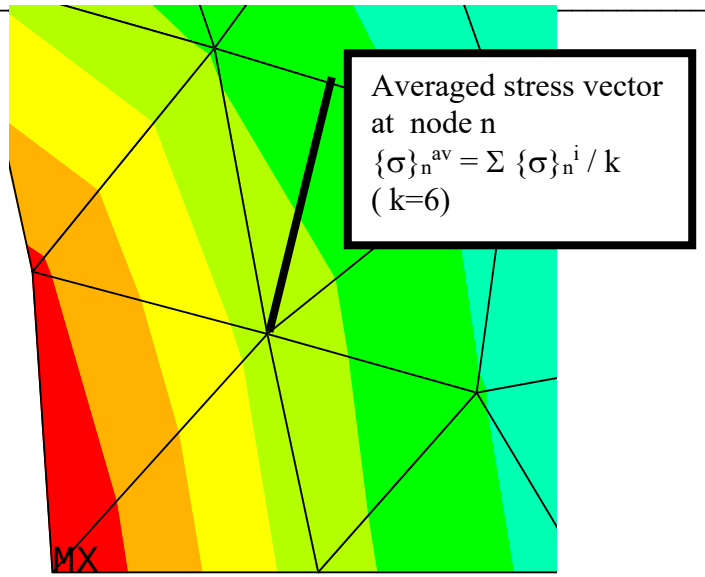
The strain energy of the element Ω_e is:

$$U_e = \frac{1}{2}\int_{\Omega_e} \lfloor \varepsilon \rfloor \{\sigma\}\, d\Omega_e, \qquad U_e = \frac{1}{2}\int_{\Omega_e} \lfloor q \rfloor_e [B]^T [D]\ [B]\ \{q\}_e\, d\Omega_e, \qquad U_e = \frac{1}{2}\lfloor q \rfloor_e [k]_e \{q\}_e.$$

where $[k]_e = \int_{\Omega_e} [B]^T [D][B] d\Omega_e$ is **the stiffness matrix of the element** (symmetrical, singular, semi-positive defined)



discontinuous



continous

stress vector at node n of element i

$$\{\sigma\}_n^i = \begin{Bmatrix} \sigma_x \\ \sigma_y \\ \sigma_z \\ \tau_{xy} \\ \tau_{yz} \\ \tau_{xz} \end{Bmatrix}_n^{\mathrm{i}} \quad \{\sigma\}_n^1 \neq \{\sigma\}_n^2 \neq \{\sigma\}_n^3 \neq \ldots$$

averaged stress vector at node n: $\{\sigma\}_n^{av} = \dfrac{\sum_{i=1}^{k} \{\sigma\}_n^i}{k}$

Stress error vector at node n of element *i* $\{\Delta\sigma\}_n^i = \{\sigma\}_n^i - \{\sigma\}_n^{av}$

The stress error vector $\{\Delta\sigma\}^i$ within the element ' i' may be determined by standard approximation using the stress error vectors at nodes of element i $\{\Delta\sigma\}_n^i$.

Strain energy of the element i

$$U_i = \frac{1}{2}\int_{\Omega_i} \lfloor\sigma\rfloor^i \{\varepsilon\}^i \, d\Omega_i$$

$\{\varepsilon\} = [D]^{-1}\{\sigma\}$, D] – stress-strain matrix

$$U_i = \frac{1}{2}\int_{\Omega_i} \lfloor\sigma\rfloor^i [D]^{-1} \{\sigma\}^i \, d\Omega_i$$

Then for each element so called energy error can be estimated

$$e_i = \frac{1}{2}\int_{\Omega_i} \lfloor\Delta\sigma\rfloor^i [D]^{-1} \{\Delta\sigma\}^i \, d\Omega_i \qquad \text{(ETABLE-SERR)}$$

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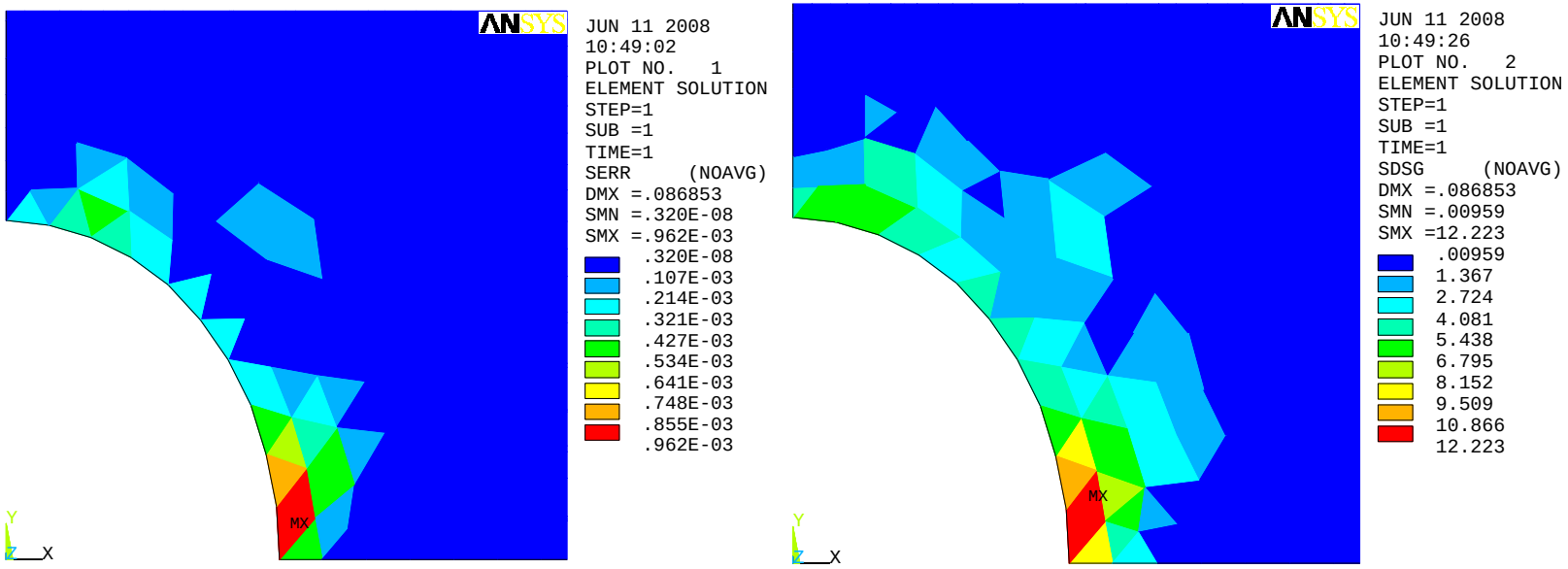


Initial mesh. SERR and SDSG error distribution

SDSG - $\Delta\sigma_i$ = maximum absolute value of any component of $\{\Delta\sigma\}_n^i = \{\sigma\}_n^i - \{\sigma\}_n^{av}$ for all nodes connected to element

The energy error over the model

$$e = \sum_{i=1}^{l.el.} e_i$$

The energy error can be normalized against the strain energy

$$\mathrm{SEPC} = 100\left(\frac{e}{U+e}\right)^{\frac{1}{2}}$$

U – total strain energy over the entire model

SEPC – percentage error in energy norm

The e_i values can be used for adaptive mesh refinement. It has been shown that if e_i is equal for all elements, then the model using the given number of elements is the most efficient one.

This concept is also referred to as "error equilibration" ($e_i = \text{const}$, SEPC< S_0).

Z^2 – method

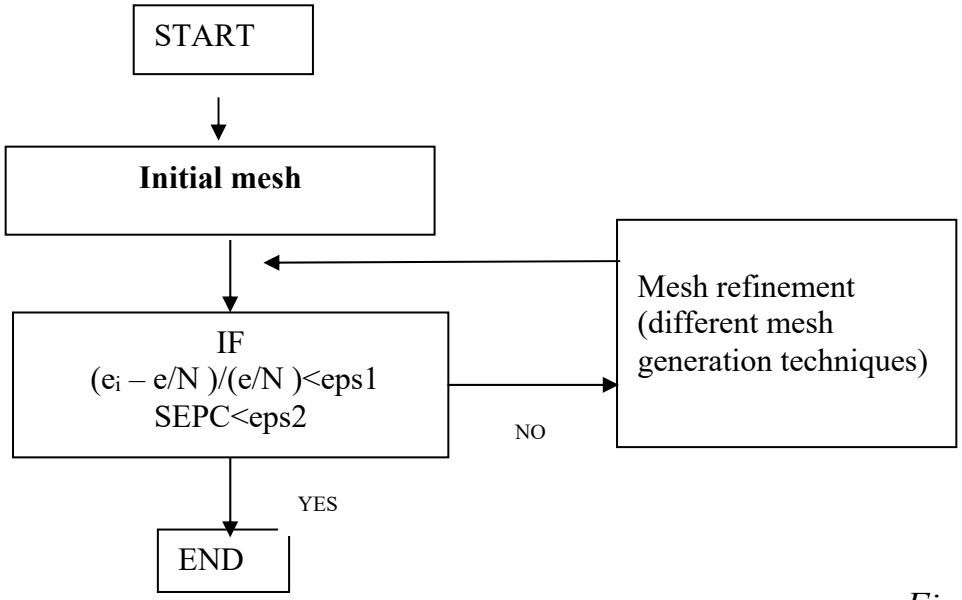
Z-Z Estimator (ZZ_Estimator) — Error estimation by Zienkiewicz and Zhu's superconvergent patch recovery.

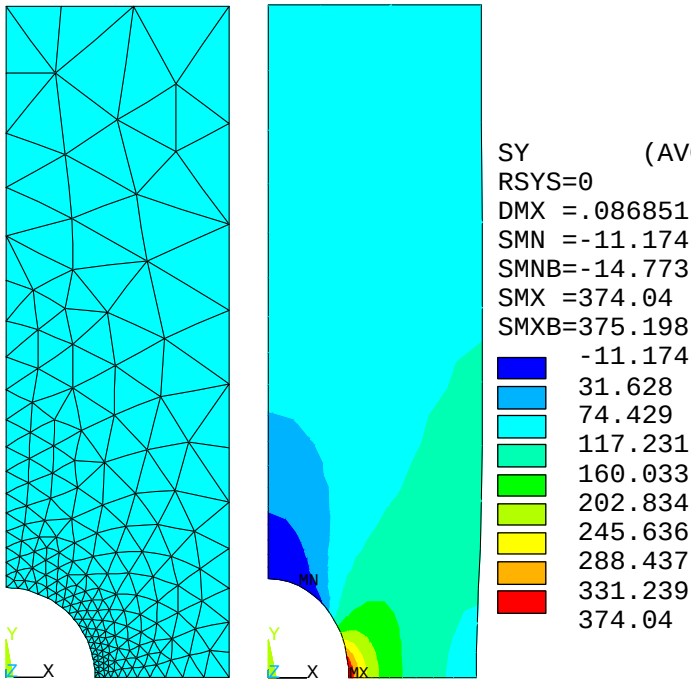
J. Z. Zhu and O. C. Zienkiewicz, "Adaptive Techniques in the Finite Element Method," *Commun. Appl. Numer. Methods*, Vol. 4, 1988, pp. 197–204.

M. Ainsworth, J. Z. Zhu, A. W. Craig, and O. C. Zienkiewicz, "Analysis of the Zienkiewicz–Zhu *a Posteriori* Error Estimator in the Finite Element Method," *Int. J. Numer. Methods Eng.*, vol. 28, 1989, pp. 2161–2174.

Adaptive Meshing Techniques

- Automatic refinement of FE meshes until converged results are obtained
- User's responsibility reduced to generation a good initial mesh

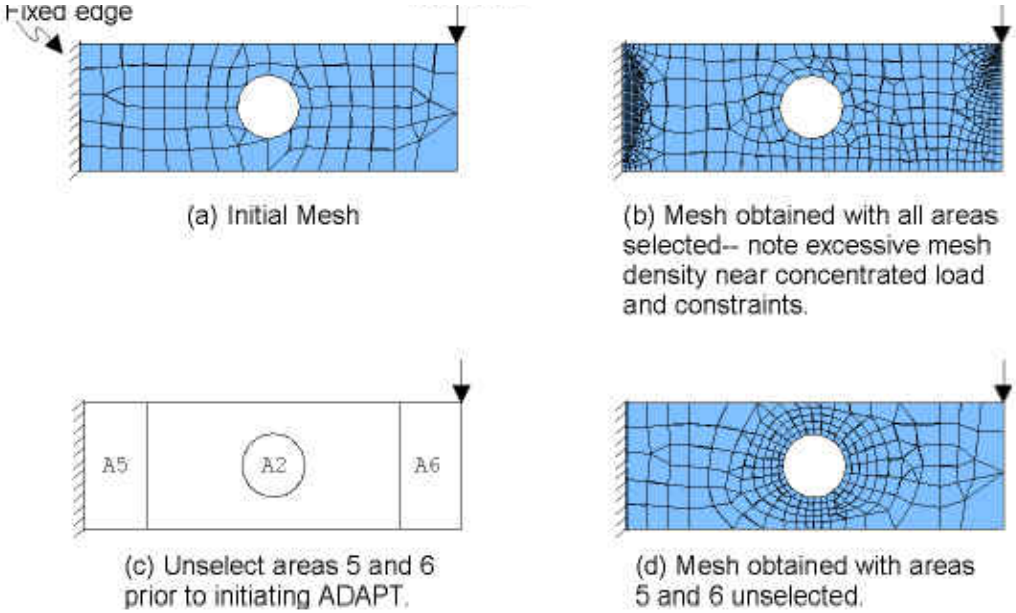




Final FE mesh and the results - Sy distribution (percentage error in energy norm *SEPC=0.811%, uniform error distribution* e_i*= const)*

Selective adaptive meshing

If mesh discretization error (measured as a percentage) is relatively unimportant in some regions of the model, the procedure may be speed up by excluding such regions from the adaptive meshing operations. Also - near singularities caused by concentrated loads and at boundaries between different materials.



(a) Initial Mesh

(b) Mesh obtained with all areas selected-- note excessive mesh density near concentrated load and constraints.

(c) Unselect areas 5 and 6 prior to initiating ADAPT.

(d) Mesh obtained with areas 5 and 6 unselected.

Selective adaptive meshing in ANSYS

Types of refinement in adaptive meshing :

h-refinement: reduction of the size of the element ("*h*" refers to the typical size of the elements)

p-refinement: increase of the order of the polynomials on an element (shape functions from linear to quadratic, etc.)

r-refinement: re-arrangement of the nodes in the mesh

hp-refinement: combination of the h- and p-refinements.