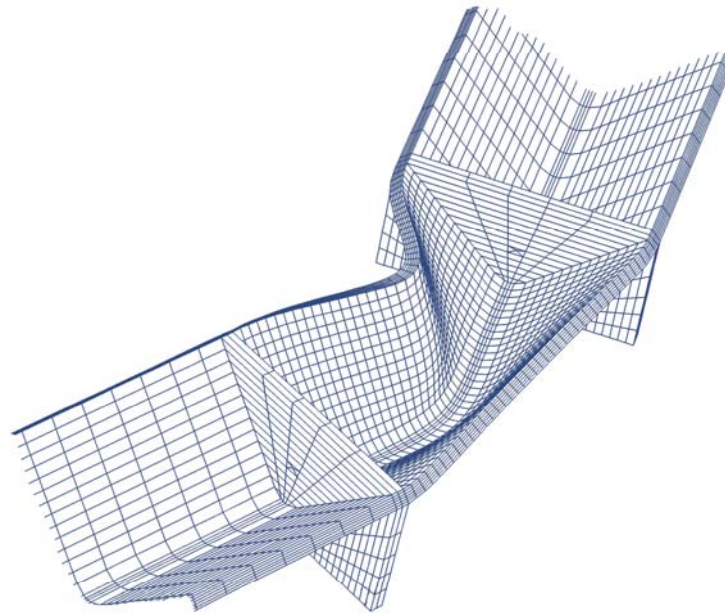
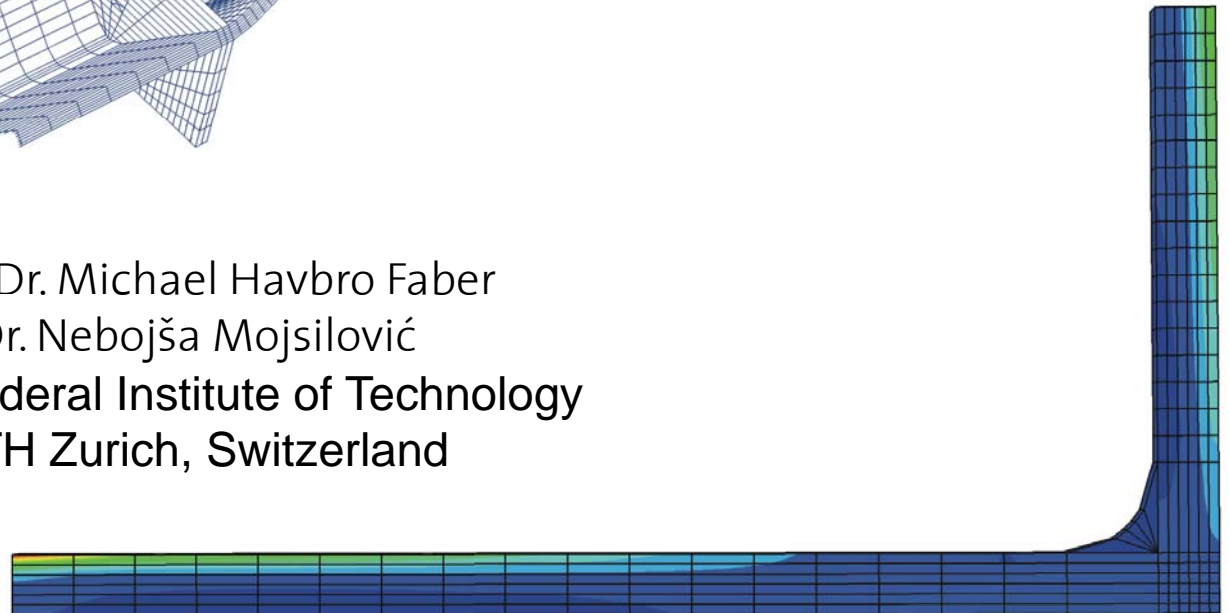


The Finite Element Method for the Analysis of Non-Linear and Dynamic Systems

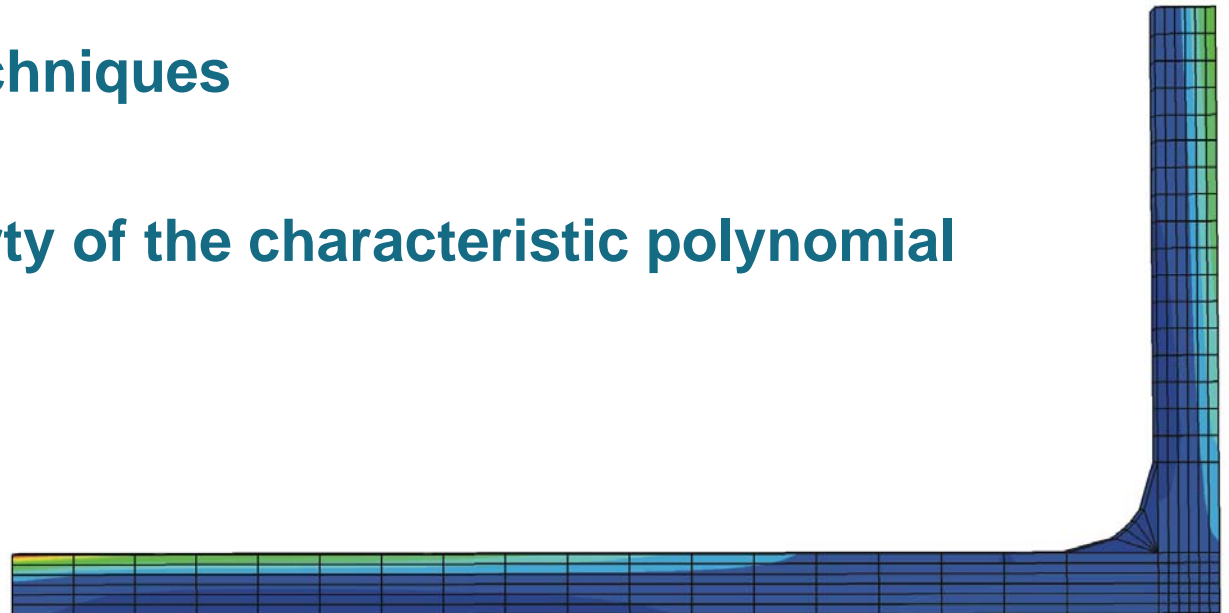


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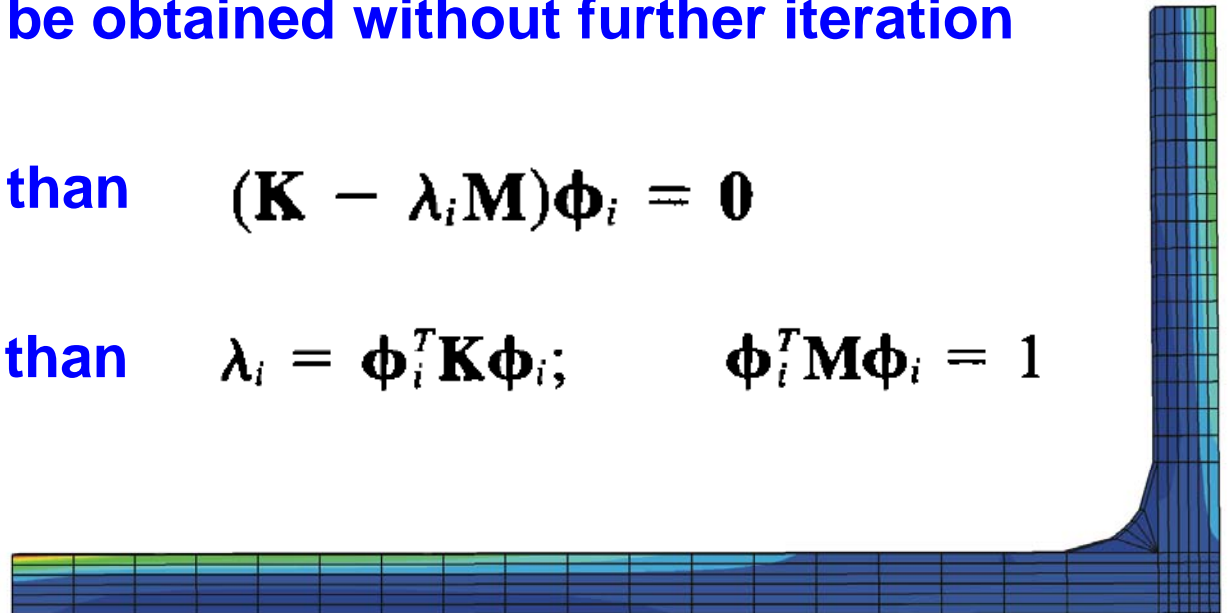
Contents of Today's Lecture

- Solution methods for eigenproblems $\mathbf{K}\boldsymbol{\phi} = \lambda\mathbf{M}\boldsymbol{\phi}$
- Vector iteration methods
- Transformation methods
- Polynomial iteration techniques
- Sturm sequence property of the characteristic polynomial



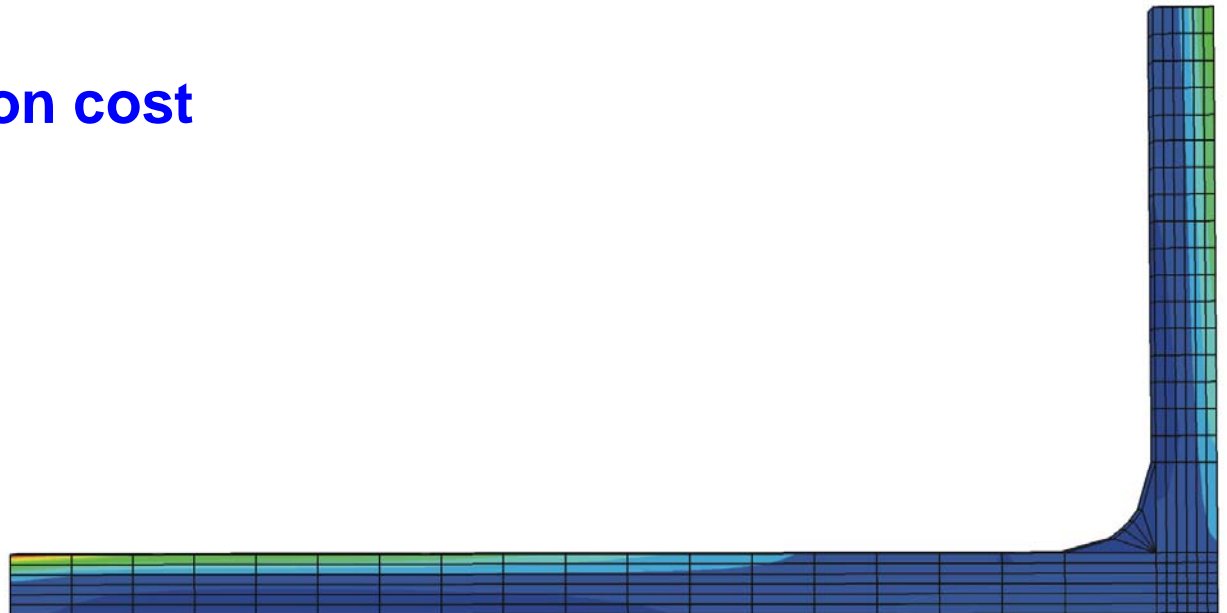
Solution methods for eigenproblems

- All solution methods are iterative since we are calculating the roots of the polynomial $p(\lambda)$, which has order of \mathbf{K} and \mathbf{M}
- There are no implicit formulas when the order of p is higher than 4
- Iteration is needed in solution of an eigenpair (λ_i, ϕ_i) ; knowing one of them the other one can be obtained without further iteration
- If λ_i solved by iteration, then $(\mathbf{K} - \lambda_i \mathbf{M})\phi_i = \mathbf{0}$
- If ϕ_i solved by iteration, then $\lambda_i = \phi_i^T \mathbf{K} \phi_i; \quad \phi_i^T \mathbf{M} \phi_i = 1$



Effectiveness of a solution method

- **Depends on two factors**
- **Firstly, on the possibility of a reliable use of the procedure, i.e. for well defined K and M matrices the solution is always obtained to the required precision without solution break-down**
- **Secondly, on the solution cost**

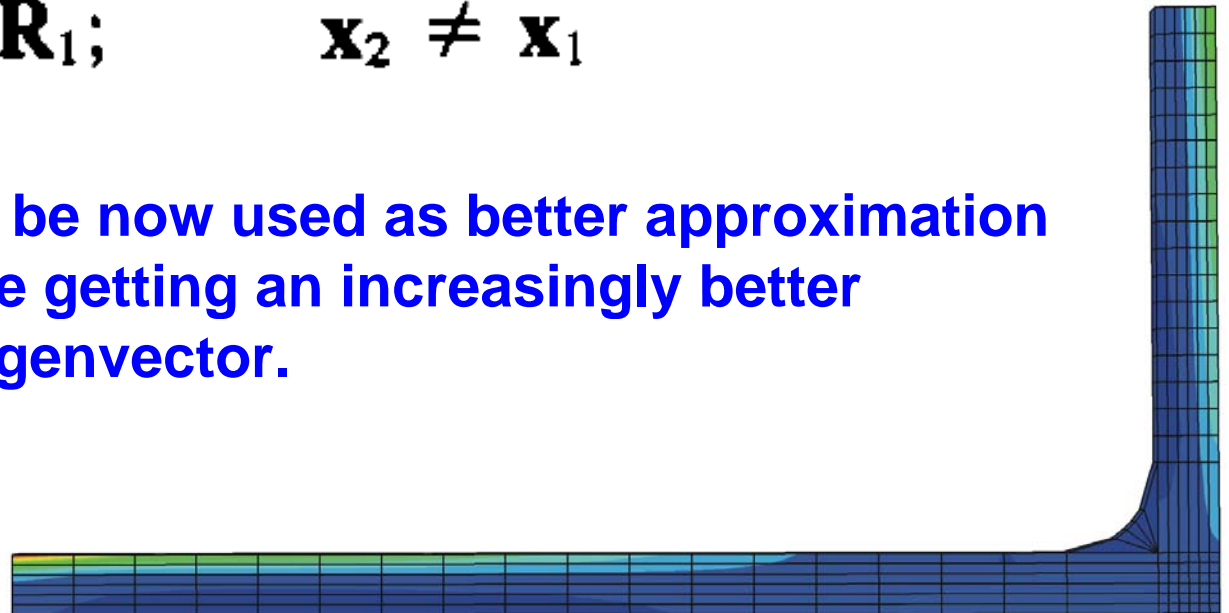


Vector iteration methods

- Problem to solve: $\mathbf{K}\boldsymbol{\phi} = \lambda\mathbf{M}\boldsymbol{\phi}$
- Assuming \mathbf{x}_1 for $\boldsymbol{\phi}$ and setting $\lambda=1$ we obtain $\mathbf{R}_1 = (\mathbf{1})\mathbf{M}\mathbf{x}_1$
- Now we can use the equilibrium equation

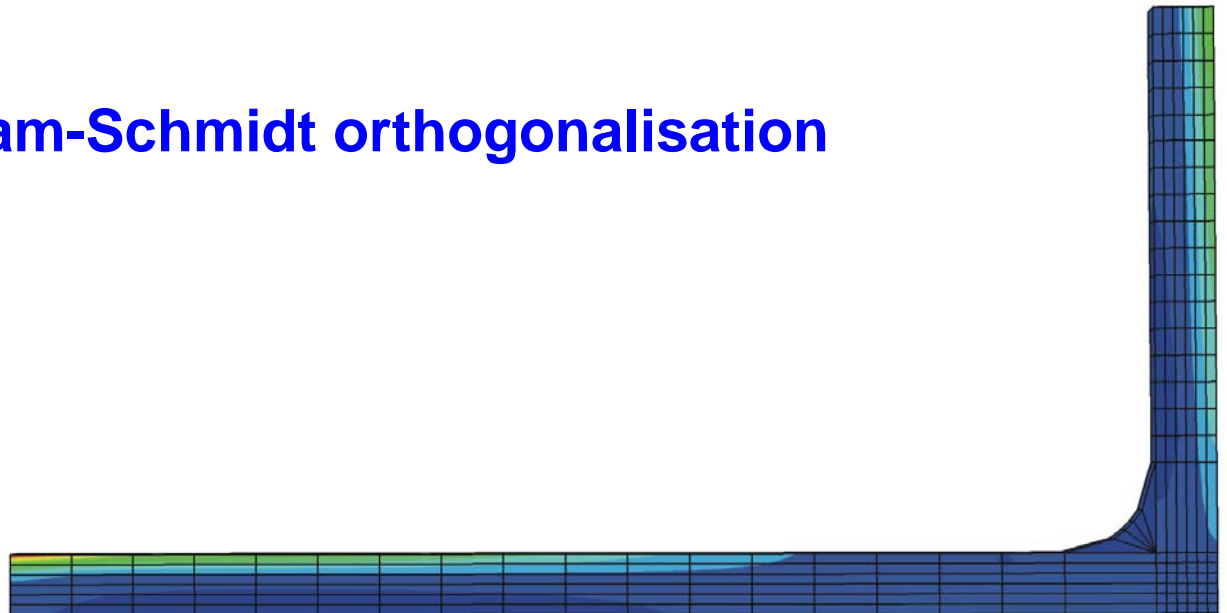
$$\mathbf{K}\mathbf{x}_2 = \mathbf{R}_1; \quad \mathbf{x}_2 \neq \mathbf{x}_1$$

- We obtain \mathbf{x}_2 which can be now used as better approximation for \mathbf{x}_1 ; in this way we are getting an increasingly better approximation for an eigenvector.



Vector iteration methods

- Inverse iteration
- Forward iteration
- Rayleigh quotient iteration
- Matrix deflation and Gram-Schmidt orthogonalisation



Inverse iteration

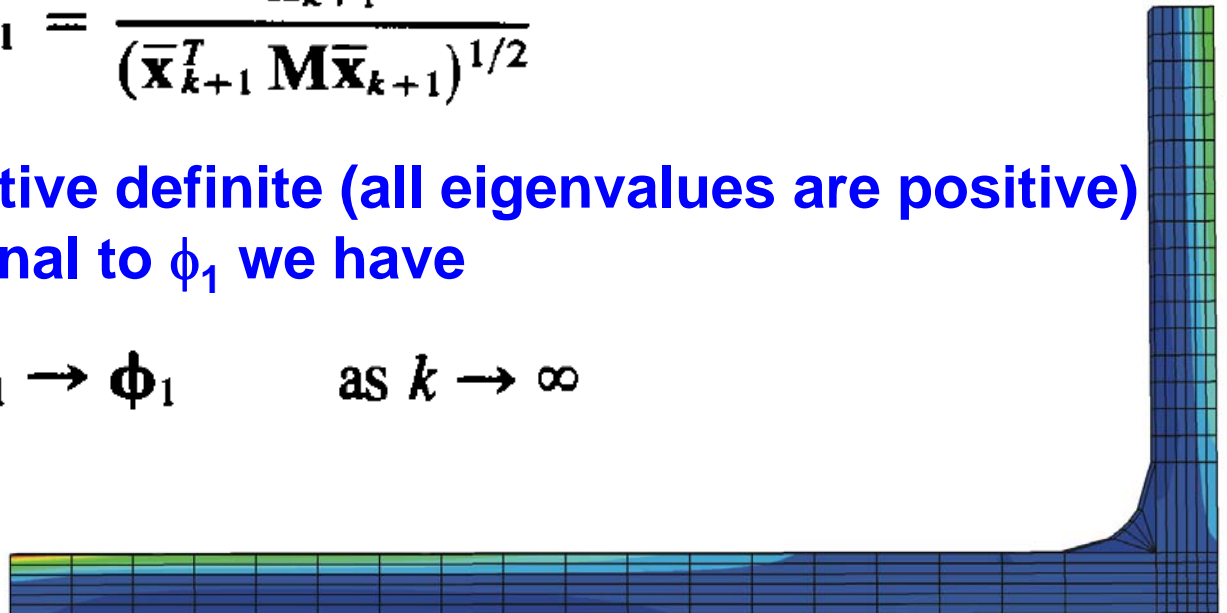
- Used to calculate an eigenvector (and later corresponding eigenvalue)
- Firstly, we assume starting vector \mathbf{x}_1 and thus in iteration step k we have

$$\mathbf{K}\bar{\mathbf{x}}_{k+1} = \mathbf{M}\mathbf{x}_k$$

$$\mathbf{x}_{k+1} = \frac{\bar{\mathbf{x}}_{k+1}}{(\bar{\mathbf{x}}_{k+1}^T \mathbf{M}\bar{\mathbf{x}}_{k+1})^{1/2}}$$

- Providing that \mathbf{K} is positive definite (all eigenvalues are positive) and \mathbf{x}_1 is not \mathbf{M} -orthogonal to ϕ_1 we have

$$\mathbf{x}_{k+1} \rightarrow \phi_1 \quad \text{as } k \rightarrow \infty$$



Forward iteration

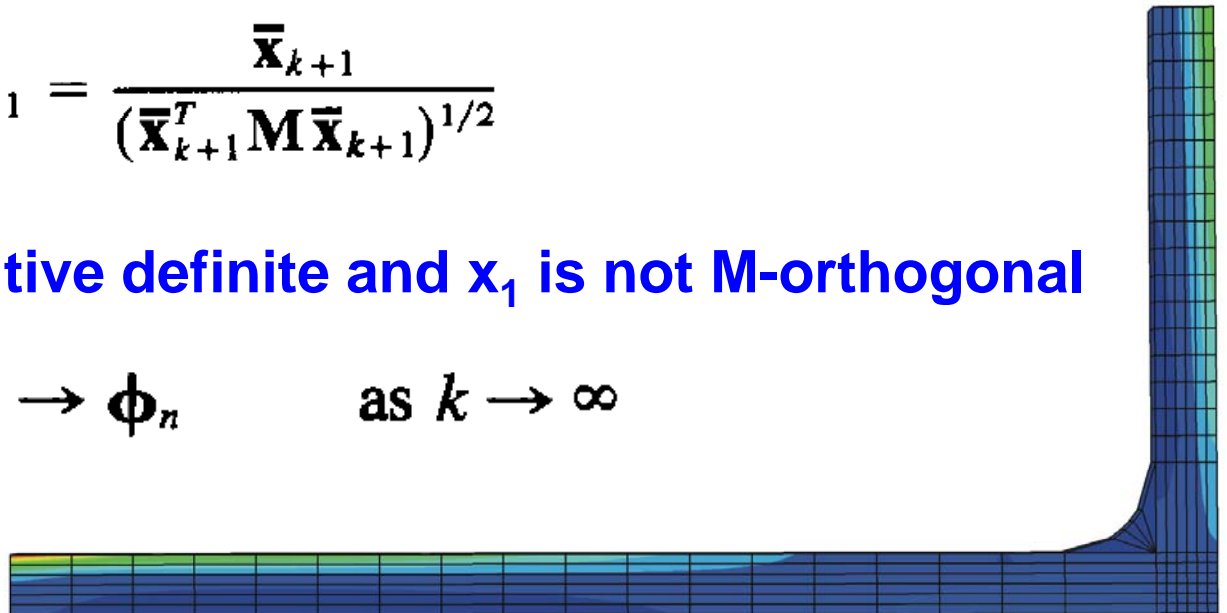
- Complementary to inverse iteration, as yielding the eigenvector corresponding to the largest eigenvalue
- Firstly, we assume starting vector \mathbf{x}_1 and thus in iteration step k we have

$$\mathbf{M}\bar{\mathbf{x}}_{k+1} = \mathbf{K}\mathbf{x}_k$$

$$\mathbf{x}_{k+1} = \frac{\bar{\mathbf{x}}_{k+1}}{(\bar{\mathbf{x}}_{k+1}^T \mathbf{M} \bar{\mathbf{x}}_{k+1})^{1/2}}$$

- Providing that \mathbf{M} is positive definite and \mathbf{x}_1 is not \mathbf{M} -orthogonal to ϕ_1 we have

$$\mathbf{x}_{k+1} \rightarrow \phi_n \quad \text{as } k \rightarrow \infty$$



Convergence of the inverse and forward iteration

- The convergence of both procedures can be proved
- Convergence rate can be improved by shifting, see also Ch. 10.2.3 (we perform a shift ρ on K matrix in order to accelerate the calculations of the required eigensystem)
- Additionally, a shift can be used to obtain convergence in inverse iteration procedure when K is positive semi definite (all eigenvalues are greater or equal zero) and in forward iteration procedure when M is diagonal with some zero diagonal elements

