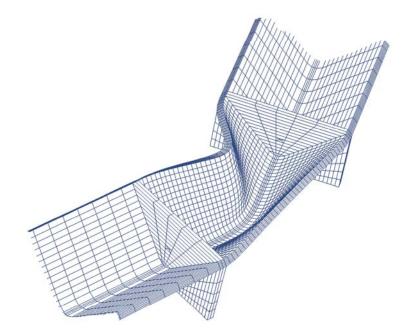


The Finite Element Method for the Analysis of Non-Linear and Dynamic Systems

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Prof. Dr. Michael Havbro Faber Dr. Nebojša Mojsilović Swiss Federal Institute of Technology ETH Zurich, Switzerland

Contents of Today's Lecture

- Analysis of direct integration methods
- Stability analysis (liner analysis)
- Accuracy analysis (liner analysis)
- Some practical considerations
- Examples

Analysis of direct integration methods

Dynamic equilibrium equation

$$\mathbf{M}\ddot{\mathbf{U}}(t) + \mathbf{C}\dot{\mathbf{U}}(t) + \mathbf{K}\mathbf{U}(t) = \mathbf{R}(t)$$

- Solution procedures: mode superposition and direct integration
- Integration schemes: central difference method, the Houbolt method, the Wilson θ method and the Newmark integration procedure

Analysis of direct integration methods

- Key to all integration procedures: time step Δt
- Small enough to obtain the accurate solution
- Large enough to save computational time
- How to select the appropriate time step for direct integration?
- Two fundamental concepts are to be considered: stability and accuracy of the integration scheme

- Relation between mode superposition and direct integration is of interest
- Mode superposition: change of basis from the nodal displacements to the basis of eigenvectors of generalised eigenproblem

$$\mathbf{K}\boldsymbol{\Phi} = \boldsymbol{\omega}^2 \mathbf{M}\boldsymbol{\Phi}$$

• Setting $U(t) = \Phi X(t)$ in dynamic equilibrium equation (columns in Φ are free-vibration modes ϕ_1, \dots, ϕ_n) one obtains

$$\ddot{\mathbf{X}}(t) + \Delta \dot{\mathbf{X}}(t) + \mathbf{\Omega}^2 \mathbf{X}(t) = \mathbf{\Phi}^T \mathbf{R}(t)$$

Analysis of direct integration methods

- where Ω^2 is diagonal matrix with free-vibration frequencies squared $\omega_1^2, ..., \omega_n^2$, Δ is also a diagonal matrix of damping Δ = diag($2\omega_i \xi_i$), ξ_i being the damping ratio in i_{th} mode
- The obtained equation system consists of n uncoupled equations and can be solved using one of the direct integration procedures we already mentioned
- We can use for each equation an appropriate time step since the periods of vibrations are known
- Choosing the same time step for all n equations we obtain the same solution as if we were using direct integration analysis
- In this way both procedures deliver the same solution

Analysis of direct integration methods

- So, to study the accuracy of direct integration we can consider the equations of mode superposition using the same time step by integration
- Thus, the variables that have to be considered in the stability and accuracy analysis of of the direct integration method are only Δt, ω_i and ξ_i for i=1,...n.
- Furthermore, we need to consider only one of the equations (since they are similar):

$$\ddot{x} + 2\xi\omega\dot{x} + \omega^2x = r$$

 This equation represents equilibrium equation governing motion of single degree of freedom system with free-vibration period T, dumping ξ and applied load r

Direct integration approximation and load operators

 Having solutions for the discrete times 0, ∆t, 2∆t,...t-∆t, t the solution for the time t+∆t can be obtained using recursive relationship in the specific integration method considered

$$^{\iota+\Delta\iota}\mathbf{\hat{X}} = \mathbf{A}^{\iota}\mathbf{\hat{X}} + \mathbf{L}(^{\iota+\nu}r)$$

Vectors X are storing solution quantities (displacements, velocities), r is a load at time t+ν, ν being 0, Δt ot θΔt depending on the chosen method. The matrix A is the integration approximation and vector L is load operator.

Direct integration approximation and load operators

 To study the accuracy and stability of the integration methods we will need the relation which gives the solution for the time t+n∆t and which is obtained in applying the above relation recursively:

$${}^{t+n\Delta t} \mathbf{\hat{X}} = \mathbf{A}^{n} {}^{t} \mathbf{\hat{X}} + \mathbf{A}^{n-1} \mathbf{L} ({}^{t+\nu} r) + \mathbf{A}^{n-2} \mathbf{L} ({}^{t+\Delta t+\nu} r) + \cdots$$

$$+ \mathbf{A} \mathbf{L} ({}^{t+(n-2)\Delta t+\nu} r) + \mathbf{L} ({}^{t+(n-1)\Delta t+\nu} r)$$

 Now, we proceed and calculate operators A and L for different direct integration methods

The central difference method

• We approximate acceleration and velocity at time t. The equilibrium equation is also considered in time t. Thus we obtain

$$\begin{bmatrix} t+\Delta t \\ t \\ t \end{bmatrix} = \mathbf{A} \begin{bmatrix} t \\ t \\ t-\Delta t \\ x \end{bmatrix} + \mathbf{L}^{t} r$$

• where

$$\mathbf{A} = \begin{bmatrix} \frac{2 - \omega^2 \,\Delta t^2}{1 + \xi \omega \,\Delta t} & -\frac{1 - \xi \omega \,\Delta t}{1 + \xi \omega \,\Delta t} \\ 1 & 0 \end{bmatrix} \qquad \mathbf{L} = \begin{bmatrix} \frac{\Delta t^2}{1 + \xi \omega \,\Delta t} \\ 0 \end{bmatrix}$$

The Houbolt method

 The equilibrium equation is considered in time t+∆t and two backward formulas are used for the acceleration and velocity at time t+∆t. Thus we obtain

$$\begin{bmatrix} t + \Delta t \\ t \\ t \\ t - \Delta t \\ t \end{bmatrix} = \mathbf{A} \begin{bmatrix} t \\ t \\ t - \Delta t \\ t \end{bmatrix} + \mathbf{L} t + \Delta t \mathbf{r}$$

• where $\mathbf{A} = \begin{bmatrix} \frac{5\beta}{\omega^2 \Delta t^2} + 6\kappa & -\left(\frac{4\beta}{\omega^2 \Delta t^2} + 3\kappa\right) & \frac{\beta}{\omega^2 \Delta t^2} + \frac{2\kappa}{3} \\ 1 & 0 & 0 \\ 0 & 1 & 0 \end{bmatrix} \quad \mathbf{L} = \begin{bmatrix} \frac{\beta}{\omega^2} \\ 0 \\ 0 \\ 0 \end{bmatrix}$ $\beta = \left(\frac{2}{\omega^2 \Delta t^2} + \frac{11\xi}{3\omega \Delta t} + 1\right)^{-1}; \quad \kappa = \frac{\xi\beta}{\omega \Delta t}$

Method of Finite Elements II

The Wilson $\boldsymbol{\theta}$ method

The equilibrium equation is considered in time t+θΔt. Thus we obtain

$$\begin{bmatrix} t+\Delta t \ddot{x} \\ t+\Delta t \dot{x} \\ t+\Delta t \chi \end{bmatrix} = \mathbf{A} \begin{bmatrix} t \ddot{x} \\ t \dot{x} \\ t \chi \end{bmatrix} + \mathbf{L} t+\theta \Delta t r$$

• where

$$\mathbf{A} = \begin{bmatrix} 1 - \frac{\beta\theta^2}{3} - \frac{1}{\theta} - \kappa\theta & \frac{1}{\Delta t}(-\beta\theta - 2\kappa) & \frac{1}{\Delta t^2}(-\beta) \\ \Delta t \left(1 - \frac{1}{2\theta} - \frac{\beta\theta^2}{6} - \frac{\kappa\theta}{2}\right) & 1 - \frac{\beta\theta}{2} - \kappa & \frac{1}{\Delta t}\left(-\frac{\beta}{2}\right) \\ \Delta t^2 \left(\frac{1}{2} - \frac{1}{6\theta} - \frac{\beta\theta^2}{18} - \frac{\kappa\theta}{6}\right) & \Delta t \left(1 - \frac{\beta\theta}{6} - \frac{\kappa}{3}\right) & 1 - \frac{\beta}{6} \end{bmatrix} \qquad \mathbf{L} = \begin{bmatrix} \frac{\beta}{\omega^2 \Delta t^2} \\ \frac{\beta}{2\omega^2 \Delta t} \\ \frac{\beta}{6\omega^2} \end{bmatrix}$$
$$\boldsymbol{\beta} = \left(\frac{\theta}{\omega^2 \Delta t^2} + \frac{\xi\theta^2}{\omega \Delta t} + \frac{\theta^3}{6}\right)^{-1}; \qquad \kappa = \frac{\xi\beta}{\omega \Delta t}$$

 $\omega^2 \Delta t$

 $\omega^2 \Delta t$

The Newmark method

 The equilibrium equation is considered in time t+Δt. For the velocity and displacements at time t+Δt two parameters, δ and α are to be chosen. Thus we obtain

$$\begin{bmatrix} t+\Delta t \ddot{x} \\ t+\Delta t \dot{x} \\ t+\Delta t x \end{bmatrix} = \mathbf{A} \begin{bmatrix} t \ddot{x} \\ t \dot{x} \\ t \\ t \end{bmatrix} + \mathbf{L} t+\Delta t r$$

• where

$$\mathbf{A} = \begin{bmatrix} -(\frac{1}{2} - \alpha)\beta - 2(1 - \delta)\kappa & \frac{1}{\Delta t}(-\beta - 2\kappa) & \frac{1}{\Delta t^{2}}(-\beta) \\ \Delta t [1 - \delta - (\frac{1}{2} - \alpha)\delta\beta - 2(1 - \delta)\delta\kappa] & 1 - \beta\delta - 2\delta\kappa & \frac{1}{\Delta t}(-\beta\delta) \\ \Delta t^{2}[\frac{1}{2} - \alpha - (\frac{1}{2} - \alpha)\alpha\beta - 2(1 - \delta)\alpha\kappa] & \Delta t(1 - \alpha\beta - 2\alpha\kappa) & (1 - \alpha\beta) \end{bmatrix} \mathbf{L}$$
$$\mathbf{\beta} = \left(\frac{1}{\omega^{2} \Delta t^{2}} + \frac{2\xi\delta}{\omega \Delta t} + \alpha\right)^{-1}; \qquad \kappa = \frac{\xi\beta}{\omega \Delta t}$$

Method of Finite Elements II

- In general, ∆t should be chosen according to the smallest period in system T_n; usually T_n/10 or smaller
- Practically, we need to consider only first p of total n equations and accordingly we need to perform finite element idealization such that the lowest p frequencies and mode shapes are predicted accurately
- Question of stability of one integration scheme: what happens when ∆t/T (T being natural period) is large?
- This question is answered by examining the behaviour of the numerical solution for arbitrary initial conditions
- To perform this we set r=0 thus getting

$$^{t+n\Delta t}\mathbf{\hat{X}} = \mathbf{A}^{n} \mathbf{\hat{X}}$$

- Considering the stability of integration method we have procedures that are:
- Unconditionally stable if the solution for any initial conditions does not grow without bound for any time step ∆t, in particular when ∆t/T is large
- Conditionally stable if the above holds only when ∆t/T is smaller or equal to a certain value, which is usually called stability limit
- For stability analysis one uses the spectral decomposition of A given by A=PJP⁻¹, where P is the matrix of eigenvectors of A and J is the Jordan canonical form of A with eigenvalues λ_i of A on its diagonal

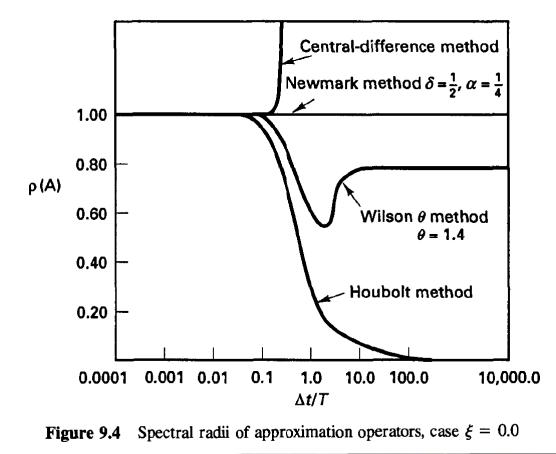
- Now we can write Aⁿ=PJⁿP⁻¹ and using this we can determine the stability of the time integration scheme
- If we now consider the spectral radius A defined as

$$\rho(\mathbf{A}) = \max_{i=1,2,\ldots} |\lambda_i|$$

- we can write the stability criterion:
- 1. If all eigenvalues are distinct we must have $\rho(A) \le 1$
- 2. If A contain multiple eigenvalues, we require that all such must be smaller than 1



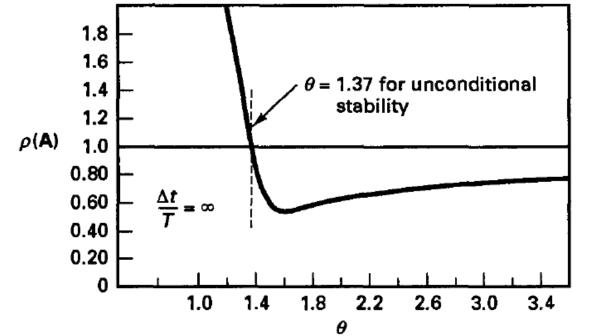
 Using simple Example (E 9.12) we can compare different integration methods using corresponding approximation operators



Method of Finite Elements II



 Furthermore, one can evaluate the optimum value of θ for the Wilson θ method



• For the Newmark method we obtain the unconditional stability for $\delta \ge 0.5$ and $\alpha \ge (0.25\delta+0.5)^2$

Accuracy analysis

- In general, using an unconditionally stable operator, the time step has to be chosen to yield an accurate and effective solution
- Integration accuracy can be assessed as a function of $\Delta t/T$, ξ and r, as we have seen before
- Considering the solution of the initial value problem defined by

$$\ddot{x} + \omega^2 x = 0$$

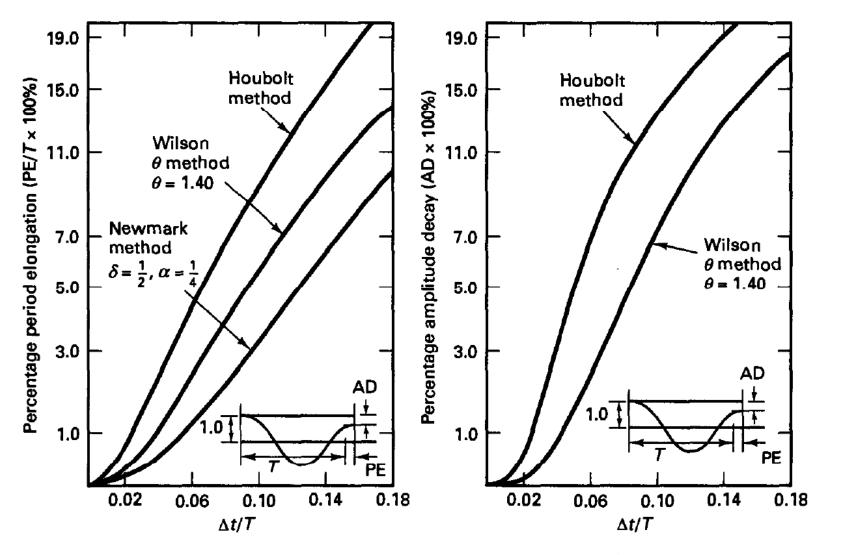
⁰x = 1.0; $^{0}\dot{x} = 0.0; \quad ^{0}\ddot{x} = -\omega^2$

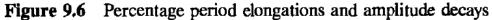
 and having an exact solution x=cos

 we can compare different integration methods using the errors in period elongation and amplitude decay as measures of accuracy



Accuracy analysis





Accuracy analysis

- All methods are accurate for very small ratios Δt/T (0.01). For larger ratios and for given Δt/T the Wilson θ method with θ=1.4 introduces less amplitude decay and period elongation than the Houbolt method, and the Newmark method the smallest percentage period elongation and no amplitude decay
- In sum, ∆t has to be chosen small enough so that the response in all modes that significantly contribute to the total structure response is calculated accurately; for central difference method ∆t should be smaller than or equal to ∆t_{cr}
- For practical analysis, both Newmark and Wilson θ method can be used



Some practical considerations

- To obtain an effective solution of a dynamic response an appropriate time integration scheme must be chosen. This choice depends on the finite element idealization which in turn depends on a physical problem to be analysed (structural dynamics or a wave propagation problem)
- Structural dynamics: only the lowest modes are considered; time step Δt should equal to $T_{co}/20$, where $T_{co}=2\pi/\omega_{co}$ and $\omega_{co}=4\omega_{u}$, ω_{u} being highest frequency significantly contained in the loading
- Wave propagation: large number of frequencies are excited in the system; "effective length" of finite element should be L_e=c∆t, where ∆t=t_w/n and t_w=L_w/c. L_w is a wavelength, c wave speed and n number of time steps needed to represent the travel of the wave