



CMHL SJTU COMPUTATIONAL MARINE HYDRODYNAMICS LAB 上海交大船舶与海洋工程计算水动力学研究中心

#### **Class-4**

#### NA26018

# Finite Element Analysis of Solids and Fluids



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2020年

# Introduction

In this section,

- we develop the finite element models of one-dimensional timedependent problems and
- describe time approximation schemes to convert ordinary differential equations in time to algebraic equations.

We consider finite element models of the time-dependent version of the differential equations studied previously. These include
The second-order (in space) parabolic equations (i.e, first time derivative)

• The second-order hyperbolice quations (i.e, second time derivative)

• The fourth-order hyperbolic equations arising in connection with the bending of beams



Finite element models of time-dependent problems can be developed in two alternative ways:

(a) coupled formulation in which the time t is treated as an additional coordinate along with the spatial coordinate x
 (b) decoupled formulation where time and spatial variations are assumed to be separable

Thus, the approximation in the two formulations takes the form  $u(x,t) \approx u_h^e(x,t) = \sum_{\substack{j=1 \ n}}^n \hat{u}_j^e \hat{\psi}_j^e(x,t)$  (coupled formulation)  $u(x,t) \approx u_h^e(x,t) = \sum_{\substack{j=1 \ n}}^n u_j^e(t) \psi_j^e(x)$  (coupled formulation)

- $\hat{\psi}_{j}^{e}(x,t)$  are time-space (two-dimensional) interpolation functions
- $\hat{u}_{j}^{e}$  are the nodal values that are independent of x and t,
- $\psi_j^e(x)$ ) are the usual one-dimensional interpolation functions in spatial coordinate x only
- $u_j^e(t)$  are functions of time t only



Space-time coupled finite element formulations are not common, and they have not been adequately studied. In this section, we consider the space-time decoupled formulation only

The space-time decoupled finite element formulation of timedependent problems involves 2 steps :

#### **1.** Spatial approximation,

- The solution *u* of the equation under consideration is approximate by decoupled form, and the spatial finite element model of the equation is developed using the procedures of static or steady-state problems while carrying all time-dependent terms in the formulation.
- This step results in a set of ordinary differential equations (i.e, a semidiscrete system of equations) in time for the nodal variables  $u_j^e(t)$  of the element. Decoupled Eq. represents the spatial approximation of u for any time t
- When the solution is separable into functions of time only and space only, u(x,t) = T(t)X(x), the approximation is clearly justified



- Even when the solution is not separable, decoupled Eq. can represent a good approximation of the actual solution provided a sufficiently small time step is used
- 2. Temporal approximation
  - The system of ordinary differential equations are further approximated in time, often using finite difference formulae for the time derivatives
  - This step allows conversion of the system of ordinary differential equations into a set of algebraic equations among u at time  $t_{s+1} = (s+1)\Delta t$ , where  $\Delta t$  is the time increment and s is a nonnegative integer

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All time approximation schemes seek to find  $u_j$  at time  $t_s + 1$  using the values of  $u_j$  from previous times :

compute  $\{u\}_{s+1}$  using $\{u\}_s, \{u\}_{s-1}, \dots$ 

Thus, at the end of the two-stage approximation, we have a continuous spatial solution at discrete intervals of time

$$u(x,t_s) \approx u_h^e(x,t_s) = \sum_{i=1}^n u_j^e(t_s)\psi_j^e(t_s) \ (s=0,1,...)$$

Note: the approximate solution has the same form as that in the separation-of variables technique used to solve boundary value and initial value problems

$$u(x, t_s) \approx u_h^e(x, t_s) = \sum_{j=1}^n u_j^e(t_s) \psi_j^e(t_s) \ (s = 0, 1, ...)$$

Note: By taking nodal values to be functions of time, we see that the spatial points in an element take on different values for different times



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- We study the details of the two steps by considering a model differential equation that contains both
- second-and fourth-order spatial derivatives
- first-and second-order time derivatives

$$-\frac{\partial}{\partial x}\left(a\frac{\partial u}{\partial x}\right) + \frac{\partial^2}{\partial x^2}\left(b\frac{\partial^2 u}{\partial x^2}\right) + c_0u + c_1\frac{\partial u}{\partial t} + c_2\frac{\partial^2 u}{\partial t^2} = f(x,t)$$

The above equation is subject to appropriate boundary and initial conditions. The boundary conditions are of the form

specify 
$$u(x,t)$$
 or  $-a\frac{\partial u}{\partial x} + \frac{\partial}{\partial x}\left(b\frac{\partial^2 u}{\partial x^2}\right)$   
specify  $\frac{\partial u}{\partial x}$  or  $b\frac{\partial^2 u}{\partial x^2}$ 

at x = 0, L, and the initial conditions involve specifying

$$c_2 u(x,0)$$
 and  $c_2 \dot{u}(x,0) + c_1 u(x,0)$   $\dot{u} \equiv \partial u / \partial t$ 



$$-\frac{\partial}{\partial x}\left(a\frac{\partial u}{\partial x}\right) + \frac{\partial^2}{\partial x^2}\left(b\frac{\partial^2 u}{\partial x^2}\right) + c_0u + c_1\frac{\partial u}{\partial t} + c_2\frac{\partial^2 u}{\partial t^2} = f(x,t)$$

- Above equation describes following physical roblems
- a. Heat transfer and fluid flow:

 $c_2 = 0 and b = 0$ 

**b.** Transverse motion of a cable:

 $a = T, c_0 = 0, b = 0, c_1 = \rho, c_2 = 0$ 

- c. The longitudinal motion of a rod: a = EA, b = 0; if damping is not considered,  $c_1 = 0, c_2 = \rho A$
- d. The transverse motion of an Euler-Bernoulli beam:

 $a = 0, b = EI, c_0 = k, c_1 = 0, c_2 = \rho A$ 



# **Semidiscrete Finite Element Models**

The semidiscrete formulation involves approximation of the spatial variation of the dependent variable. The formulation follows essentially the same steps as described in previous

- The first step involves the construction of the weak form of the equation over a typical element
- In the second step, we develop the finite element model by seeking approximation of the decoupled form



$$-\frac{\partial}{\partial x}\left(a\frac{\partial u}{\partial x}\right) + \frac{\partial^2}{\partial x^2}\left(b\frac{\partial^2 u}{\partial x^2}\right) + c_0u + c_1\frac{\partial u}{\partial t} + c_2\frac{\partial^2 u}{\partial t^2} = f(x,t)$$

$$\begin{split} 0 &= \int_{x_a}^{x_b} w \left[ -\frac{\partial}{\partial x} \left( a \frac{\partial u}{\partial x} \right) + \frac{\partial^2}{\partial x^2} \left( b \frac{\partial^2 u}{\partial x^2} \right) + c_0 u + c_1 \frac{\partial u}{\partial t} + c_2 \frac{\partial^2 u}{\partial t^2} - f \right] dx \\ &= \int_{x_a}^{x_b} \left[ \frac{\partial w}{\partial x} a \frac{\partial u}{\partial x} + \frac{\partial^2 w}{\partial x^2} b \frac{\partial^2 u}{\partial x^2} + c_0 w u + c_1 w \frac{\partial u}{\partial t} + c_2 w \frac{\partial^2 u}{\partial t^2} - w f \right] dx \\ &\quad + \left[ w \left[ \left( -a \frac{\partial u}{\partial x} \right) + \frac{\partial}{\partial x} \left( b \frac{\partial^2 u}{\partial x^2} \right) \right] + \frac{\partial w}{\partial x} \left( -b \frac{\partial^2 u}{\partial x^2} \right) \right]_{x_a}^{x_b} \\ &= \int_{x_a}^{x_b} \left[ a \frac{\partial w}{\partial x} \frac{\partial u}{\partial x} + b \frac{\partial^2 w}{\partial x^2} \frac{\partial^2 u}{\partial x^2} + c_0 w u + c_1 w \frac{\partial u}{\partial t} + c_2 w \frac{\partial^2 u}{\partial t^2} - w f \right] dx \\ &\quad - \hat{Q}_1 w(x_a) - \hat{Q}_3 w(x_b) - \hat{Q}_2 \left( -\frac{\partial w}{\partial x} \right) \right|_{x_a} - \hat{Q}_4 \left( -\frac{\partial w}{\partial x} \right) \right|_{x_b} \end{split}$$

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$$\hat{Q}_{1} = \left[ -a \frac{\partial u}{\partial x} + \frac{\partial}{\partial x} \left( b \frac{\partial^{2} u}{\partial x^{2}} \right) \right]_{x_{a}}, \hat{Q}_{2} = \left[ b \frac{\partial^{2} u}{\partial x^{2}} \right]_{x_{a}}$$
$$\hat{Q}_{3} = -\left[ -a \frac{\partial u}{\partial x} + \frac{\partial}{\partial x} \left( b \frac{\partial^{2} u}{\partial x^{2}} \right) \right]_{x_{b}}, \hat{Q}_{4} = -\left[ b \frac{\partial^{2} u}{\partial x^{2}} \right]_{x_{b}}$$

Next, we assume that u is interpolated by an expression of the decoupled form:

$$u(x,t) \approx u_h^e(x,t) = \sum_{j=1}^n u_j^e(t)\psi_j^e(x)$$
 (decoupled formulation)

This equation implies that, at any arbitrarily fixed time t > 0, the function u can be approximated by a linear combination of the  $\psi_j^e$  and  $u_j^e(t)$ , with  $u_j^e(t)$  being the value of u at time t at the jth node of the element Omegae. In other words, the time and spatial variations of u are separable.

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- This assumption is not valid, in general, because it may not be possible to write the solution u(x,t) as the product of a function of time only and a function of space only
- However, with sufficiently small time steps, it is possible to obtain accurate solutions to even those problems for which the solution is not separable in time and space
- The finite element solution that we obtain at the end of the analysis is continuous in space but not in time
- We only obtain the finite element solution in the form

$$u(x,t_s) = \sum_{j=1}^n u_j^e(t_s)\psi_j^e(t_s) = \sum_{j=1}^n (u_j^s)^e \psi_j^e(x) \quad (s = 0,1,...)$$

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Where  $(u_j^s)^e$  is the value of u(x, t) at time  $t = t_s$  and node j of the element Omegae

Substituting  $w = \psi_j^e(x)$  (to obtain the ith equation of the system) and substitute decoupled approximation into weak form, we obtain

$$u(x,t) \approx u_h^e(x,t) = \sum_{j=1}^n u_j^e(t)\psi_j^e(x)$$
 (decoupled formulation)

$$0 = \int_{x_a}^{x_b} \left[ a \frac{\partial w}{\partial x} \frac{\partial u}{\partial x} + b \frac{\partial^2 w}{\partial x^2} \frac{\partial^2 u}{\partial x^2} + c_0 w u + c_1 w \frac{\partial u}{\partial t} + c_2 w \frac{\partial^2 u}{\partial t^2} - w f \right] dx$$
$$-\hat{Q}_1 w(x_a) - \hat{Q}_3 w(x_b) - \hat{Q}_2 \left( -\frac{\partial w}{\partial x} \right) \bigg|_{x_a} - \hat{Q}_4 \left( -\frac{\partial w}{\partial x} \right) \bigg|_{x_b}$$

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$$0 = \int_{x_a}^{x_b} \left[ a \frac{d\psi_i}{dx} \left( \sum_{j=1}^n u_j \frac{d\psi_j}{dx} \right) + b \frac{d^2\psi_i}{dx^2} \left( \sum_{j=1}^n u_j \frac{d^2\psi_j}{dx^2} \right) + c_0\psi_i \left( \sum_{j=1}^n u_j\psi_j \right) \right. \\ \left. + c_1\psi_i \left( \sum_{j=1}^n \frac{du_j}{dt}\psi_j \right) + c_2\psi_i \left( \sum_{j=1}^n \frac{d^2u_j}{dt^2}\psi_j \right) - \psi_i f \right] dx \\ \left. - \hat{Q}_1\psi_i(x_a) - \hat{Q}_3\psi_i(x_b) - \hat{Q}_2 \left( -\frac{d\psi_j}{dx} \right) \right|_{x_a} - \hat{Q}_4 \left( -\frac{d\psi_j}{dx} \right) \right|_{x_b} \\ \left. = \sum_{i=1}^n \left[ \left( K_{ij}^1 + K_{ij}^2 \right) u_j + M_{ij}^1 \frac{du_j}{dt} + M_{ij}^2 \frac{d^2u_j}{dt^2} \right] - F_i \right]$$

In matrix form, we have

$$[K]{u} + [M^1]{\dot{u}} + [M^2]{\ddot{u}} = \{F\}$$

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$$[K]{u} + [M^1]{\dot{u}} + [M^2]{\ddot{u}} = {F}$$
 (a)

where

$$\begin{split} & [K] = [K^{1}] + [K^{2}] + [M^{0}] \\ & M_{ij}^{0} = \int_{x_{a}}^{x_{b}} c_{0} \psi_{i} \psi_{j} dx, \\ & M_{ij}^{2} = \int_{x_{a}}^{x_{b}} c_{2} \psi_{i} \psi_{j} dx, \\ & K_{ij}^{1} = \int_{x_{a}}^{x_{b}} c_{2} \psi_{i} \psi_{j} dx, \\ & K_{ij}^{2} = \int_{x_{a}}^{x_{b}} b \frac{d^{2} \psi_{i}}{dx^{2}} \frac{d^{2} \psi_{j}}{dx^{2}} dx, \\ & F_{ij} = \int_{x_{a}}^{x_{b}} b \frac{d^{2} \psi_{i}}{dx^{2}} \frac{d^{2} \psi_{j}}{dx^{2}} dx, \\ & F_{i} = \int_{x_{a}}^{x_{b}} \psi_{i} f \, dx + \hat{Q}_{i} \end{split}$$

Equation (a) is a hyperbolic equation, and it contains the parabolic equation as a special case (set [M] = [0]). The time approximation of (a) for these two cases will be considered separately

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# **Parabolic Equations-Time Approximation**

- The time approximation is discussed with the help of a single first-order differential equation
- Suppose that we wish to determine u(t) for t > 0 such that u(t) satisfies

$$a \frac{du}{dt} + bu = f(t), \qquad 0 < t < T \text{ and } u(0) = u_0$$

where  $a \neq 0$ , *b*, and  $u_0$  are constants, and *f* is a function of time *t*.

 The exact solution of the problem consists of two parts: the homogeneous and particular solutions. The homogeneous solution is

$$u^h(t) = Ae^{-kt}, k = \frac{b}{a}$$

The particular solution is

$$u^{p}(t) = \frac{1}{a}e^{-kt}\left(\int_{0}^{t}e^{k\tau}f(\tau)\right)d\tau$$



The complete solution is given by

$$u(t) = e^{-kt} \left( A + \frac{1}{a} \int_0^t e^{k\tau} f(\tau) \right) d\tau$$

- In the finite difference solution of parabolic Eq., we replace the derivatives with their finite difference approximation
- The most commonly used scheme is the a family of approximation in which a weighted average of the time derivatives at two consecutive time steps is approximated by linear interpolation of the values of the variable at the two steps (as in Fig. next)

$$(1-\alpha)\dot{u}_s + \alpha\dot{u}_{s+1} = \frac{u_{s+1} - u_s}{\Delta t_{s+1}} \text{ for } 0 \le \alpha \le 1$$

us, denotes the value of u(t) at time  $\Delta t_s = t_s - t_{s-1}$  is the sth time step

 $t = t_s = \sum_{i=1}^{s} \Delta t_i$ 

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If the total time [0, T] is divided into equal time steps, then  $t_s = s\Delta t$ , and

$$(1 - \alpha)\dot{u}_s + \alpha\dot{u}_{s+1} = \frac{u_{s+1} - u_s}{\Delta t_{s+1}} \text{ for } 0 \le \alpha \le 1$$
 (a)

$$u_{s+1} = u_s + \Delta t \dot{u}_{s+\alpha}$$
$$\dot{u}_{s+\alpha} = (1-\alpha)\dot{u}_s + \alpha \dot{u}_{s+1} \text{ for } 0 \le \alpha \le 1$$
 (b)

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When  $\alpha = 0$ , Eq.(a) gives

$$\dot{u}_s = \frac{u_{s+1} - u_s}{\Delta t_{s+1}}$$

This is the slope of the function u(t) at time  $t = t_s$  based on the values of the function at time ts and  $t_s + 1$ 

 Since the value of the function from a step in front is used, it is termed a forward difference approximation

$$(1 - \alpha)\dot{u}_{s} + \alpha\dot{u}_{s+1} = \frac{u_{s+1} - u_{s}}{\Delta t_{s+1}} \text{ for } 0 \le \alpha \le 1$$
 (a)

When  $\alpha = 1$ , we obtain

$$\dot{u}_{s+1} = \frac{u_{s+1} - u_s}{\Delta t_{s+1}} \longrightarrow \dot{u}_s = \frac{u_s - u_{s-1}}{\Delta t_s}$$

which is termed, for obvious reason, the **backward difference** approximation



**Recall the parabolic Eq.:** 

$$a \frac{du}{dt} + bu = f(t), 0 < t < T \text{ and } u(0) = u_0$$

Note that it is valid for all times 0 < t < T. In particular, it is valid at times  $t = t_s$ , and  $t = t_s + 1$ . Hence,

$$\dot{u}_s = \frac{1}{a}(f_s - bu_s), \dot{u}_{s+1} = \frac{1}{a}(f_{s+1} - bu_{s+1})$$

Substituting the above expressions into finite difference approximation (a):

$$(1 - \alpha)\dot{u}_s + \alpha\dot{u}_{s+1} = \frac{u_{s+1} - u_s}{\Delta t_{s+1}} \text{ for } 0 \le \alpha \le 1$$
 (a)

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$$(1 - \alpha)(f_s - bu_s) + \alpha(f_{s+1} - bu_{s+1}) = \alpha \frac{u_{s+1} - u_s}{\Delta t_{s+1}}$$

Solving for  $u_s + 1$ , we obtain

 $[a + \alpha \Delta t_{s+1}b]u_{s+1} = [a - (1 - \alpha)\Delta t_{s+1}b]u_s + \Delta t_{s+1}[\alpha f_{s+1} + (1 - \alpha)f_s]$ 

$$u_{s+1} = \frac{a - (1 - \alpha)\Delta t_{s+1}b}{a + \alpha\Delta t_{s+1}b}u_s + \Delta t_{s+1}\frac{[\alpha f_{s+1} + (1 - \alpha)f_s]}{a + \alpha\Delta t_{s+1}b}$$

Thus, above Eq. can be used repeatedly to march in time and obtain the solution at times  $t = t_s + 1$ ,  $t_s + 2$ , ...,  $t_N$ , Ntime is the number of time steps required to reach the final T

At the very beginning, i. e s = 0, the solution u is calculated using the initial value  $u_0$ :

$$u_1 = \frac{a - (1 - \alpha)\Delta t_1 b}{a + \alpha \Delta t_1 b} u_0 + \Delta t_1 \frac{[\alpha f_1 + (1 - \alpha)f_0]}{a + \alpha \Delta t_1 b}$$

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We may also develop a time approximation scheme using the finite element method

• To this end, we consider the same parabolic problem

$$a \frac{du}{dt} + bu = f(t), 0 < t < T \text{ and } u(0) = u_0$$

• We wish to determine  $u_{s+1}$  in terms of  $u_s$ The weighted-integral form of the parabolic over the time interval  $(t_s, t_{s+1})$  is  $0 = \int_{0}^{t_{s+1}} u(t) \left(a \frac{du}{dt} + bu - f\right) dt$ 

$$0 = \int_{t_s}^{s+1} v(t) \left( a \frac{du}{dt} + bu - f \right) dt$$

where u is the weight function. Assuming a solution of the form

$$u(t) \approx \sum_{j=1}^{n} u_j \psi_j(t)$$

where  $\psi_j(t)$  are interpolation functions of order (n-1). The Galerkin finite element model is obtained by substituting the above approximation for u and  $v = \psi i$ . We obtain



$$[A]\{u\} = \{F\}$$

$$A_{ij} = \int_{t_s}^{t_{s+1}} \psi_i(t) \left( a \frac{d\psi_j}{dt} + b\psi_j \right) dt , F_i = \int_{t_s}^{t_{s+1}} \psi_i(t) f(t) dt$$

- Equation is valid with the time interval  $(t_s, t_{s+1})$ , and it represents a relationship be tween the values  $u_1, u_2, ..., u_n$ , which are the values of u at time  $t_s, t_s + \Delta t/(n-1), t_s + 2\Delta t/(n-1), ..., t_{s+1}$ , respectively
- This would yield a multistep approximation scheme
- To obtain a single-step approximation scheme, i.e., write  $u_s + 1$  in terms of  $u_s$  only, we assume linear approximation (i.e. n = 2)

$$u(t) = u_s \psi_1(t) + u_{s+1} \psi_2(t)$$

$$\psi_1(t) = \frac{t_{s+1} - t_s}{\Delta t}$$
 and  $\psi_2(t) = \frac{t - t_s}{\Delta t}$ 

For this choice of approximation, the Matrix form can be written

1

as



$$\begin{pmatrix} a \\ 2 \begin{bmatrix} -1 & 1 \\ -1 & 1 \end{bmatrix} + \frac{b\Delta t}{6} \begin{bmatrix} 2 & 1 \\ 1 & 2 \end{bmatrix} \end{pmatrix} \begin{pmatrix} u_s \\ u_{s+1} \end{pmatrix} = \frac{\Delta t}{6} \begin{pmatrix} f_s \\ 2f_{s+1} \end{pmatrix}$$

Assuming that  $u_s$  is known, we solve for  $u_{s+1}$  from the second equation

$$\left(a + \frac{2b\Delta t}{3}\right)u_{s+1} = \left(a - \frac{b\Delta t}{3}\right)u_s + \Delta t\left(\frac{f_s}{3} + \frac{2f_{s+1}}{3}\right)$$

Recall, using Finite difference method, we get

 $[a + \alpha \Delta t_{s+1}b]u_{s+1} = [a - (1 - \alpha)\Delta t_{s+1}b]u_s + \Delta t_{s+1}[\alpha f_{s+1} + (1 - \alpha)f_s]$ 

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By comparison,

we find that the Galerkin scheme is a specia case of the a finite difference family of approximation, with  $\alpha = 2/3$ 



# Parabolic Equations-Stable and Conditionally Stable Schemes

$$u_{s+1} = \frac{a - (1 - \alpha)\Delta t_{s+1}b}{a + \alpha\Delta t_{s+1}b}u_s + \Delta t_{s+1}\frac{[\alpha f_{s+1} + (1 - \alpha)f_s]}{a + \alpha\Delta t_{s+1}b}$$

can be written in the form:

$$u_{s+1} = A(u_s) + F_{s,s+1}, A = \frac{a - (1 - \alpha)\Delta t_{s+1}b}{a + \alpha\Delta t_{s+1}b}$$
$$F_{s,s+1} = \Delta t_{s+1} \frac{[\alpha f_{s+1} + (1 - \alpha)f_s]}{a + \alpha\Delta t_{s+1}b}$$

The operator A is known as the amplification operator. Since us is an approximate solution, the error  $E_s = u_a(t_s) - u_s$  at time  $t_s$  (where  $u_a$  is the exact solution) will influence the solution at  $t_{s+1}$ 

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$$u_{s+1} = A(u_s) + F_{s,s+1}, A = \frac{a - (1 - \alpha)\Delta t_{s+1}b}{a + \alpha \Delta t_{s+1}b}$$
$$F_{s,s+1} = \Delta t_{s+1} \frac{[\alpha f_{s+1} + (1 - \alpha)f_s]}{a + \alpha \Delta t_{s+1}b}$$

- The error will grow (i.e, Es will be amplified) as we march in time if the magnitude of the operator is greater than 1, |A| > 1
- When the error grows without bound, the computational scheme becomes unstable (i.e, solution u<sub>s+1</sub> becomes unbounded with time)
- Therefore, in order for the scheme to be stable, it is necessary that  $|A| \le 1$ :

$$\left|\frac{a - (1 - \alpha)\Delta t_{s+1}b}{a + \alpha\Delta t_{s+1}b}\right| \le 1$$

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The above equation places a restriction on the magnitude of the time step for certain values of  $\alpha$ 

- When the error remains bounded for any time step (i.e., condition is trivially satisfied for any value of  $\Delta t$ , the scheme is stable
- If the error remains bounded only when the time step  $\Delta t$  remains below certain value, the scheme is said to be conditionally stable

For different values of  $\alpha$ , the time approximation scheme yields a different scheme. The following well-known time-approximation schemes along with their order of accuracy and stability should be noted

 $\alpha = \begin{cases} 0, & \text{The forward difference (or Euler) scheme (conditionally stable); order of accuracy=0(\Delta t) \\ \hline 2, & \text{The Crank-Nicolson scheme (stable);} 0(\Delta t)^2 \\ \hline 2, & \text{The Galerkin method (stable);} 0(\Delta t)^2 \\ \hline 1, & \text{The backward difference scheme (stable);} 0(\Delta t) \end{cases}$ 

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