

# Finite Element Method for 1D second order elliptic equation

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# Outline

Finite Element  
Method for 1D  
second order  
elliptic equation

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Weak/Galerkin  
formulation

FE space

FE discretization

Boundary  
treatment

FE method

- 1 Weak/Galerkin formulation
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- 3 FE discretization
- 4 Boundary treatment
- 5 FE method

# Target problem

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- Solve

$$-\frac{d}{dx} \left( c(x) \frac{du(x)}{dx} \right) = f(x), \quad a < x < b,$$
$$u(a) = g_a, \quad u(b) = g_b.$$

for  $u(x)$ .

- Why do we start from this problem?

An easy look at the basic idea of the finite element method.

## Numerical methods for partial differential equations:

- finite element method
- finite difference method
- finite volume method
- boundary element method
- ...

which use different techniques to discretize partial differential equations.

# Weak formulation

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- First, multiply a function  $v(x)$  on both sides of the original equation,

$$-\frac{d}{dx} \left( c(x) \frac{du(x)}{dx} \right) = f(x), \quad a < x < b$$

$$\Rightarrow -\frac{d}{dx} \left( c(x) \frac{du(x)}{dx} \right) v(x) = f(x)v(x), \quad a < x < b$$

$$\Rightarrow -\int_a^b \frac{d}{dx} \left( c(x) \frac{du(x)}{dx} \right) v(x) dx = \int_a^b f(x)v(x) dx.$$

- $u(x)$  is called a **trial** function and  $v(x)$  is called a **test** function.

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- Second, using integration by parts, we obtain

$$\begin{aligned}& \int_a^b \frac{d}{dx} \left( c(x) \frac{du(x)}{dx} \right) v(x) dx \\&= \int_a^b (cu')' v dx \\&= \int_a^b v d(cu') \\&= cu' v \Big|_a^b - \int_a^b cu' dv \\&= c(b)u'(b)v(b) - c(a)u'(a)v(a) - \int_a^b cu' v' dx.\end{aligned}$$

- Then

$$-c(b)u'(b)v(b) + c(a)u'(a)v(a) + \int_a^b cu'v' dx = \int_a^b fv dx.$$

- Since the solution at  $x = a$  and  $x = b$  are given by  $u(a) = g_a$ ,  $u(b) = g_b$ , then we can choose the test function  $v(x)$  such that  $v(a) = v(b) = 0$ .

- Hence

$$\int_a^b cu'v' dx = \int_a^b fv dx.$$

- What spaces should  $u$  and  $v$  belong to? **Sobolev spaces!**

# 1D Sobolev spaces

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## Definition (Support)

*If  $u$  is a function, then its support  $\text{supp}(u)$  is the closure of the set on which  $u$  is nonzero.*

## Definition (Compactly supported)

*If  $u$  is a function defined on an open interval  $I$  and  $\text{supp}(u)$  is a compact subset (that is, a closed and bounded subset), then  $u$  is said to be compactly supported in  $I$ .*

## Lemma

*A function compactly supported in an open interval  $I$  is zero on and near the boundary of  $I$ .*



## Definition

$C_0^\infty(I)$  is the set of all functions that are infinitely differentiable on  $I$  and compactly supported in  $I$ .

- Recall integration by parts:

$$\begin{aligned}\int_a^b u' v \, dx &= uv \Big|_a^b - \int_a^b uv' \, dx \\ &= u(b)v(b) - u(a)v(a) - \int_a^b uv' \, dx.\end{aligned}$$

- For  $v \in C_0^\infty(I)$ , we have  $v(a) = v(b) = 0$ . Then

$$\int_a^b u' v \, dx = - \int_a^b uv' \, dx.$$

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## Definition (weak derivative)

*Suppose  $u$  is a real-valued function defined on an open interval  $I = (a, b)$  and that  $u$  is integrable over every compact subset of  $I$ . If there exists another locally integrable function  $w$  defined on  $I$  such that*

$$\int_a^b wv \, dx = - \int_a^b uv' \, dx$$

*for all  $v \in C_0^\infty(I)$ , then  $u$  is said to be weakly differentiable and  $w$  is called the weak derivative of  $u$ .*

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## Lemma

*If  $u$  is differentiable, then  $u$  is weakly differentiable and its weak derivative is  $u'$ .*

## Remark

*In the Sobolev spaces, which will be defined below,  $u'$  is used to represent the weak derivative.*

### Definition ( $L^2$ space)

$$L^2(I) = \left\{ v : I \rightarrow \mathbb{R} : \int_a^b v^2 dx < \infty \right\}, \quad I = (a, b).$$

### Definition ( $H^1$ space)

$$H^1(I) = \{ v \in L^2(I) : v' \in L^2(I) \}, \quad I = (a, b).$$

### Definition ( $H_0^1$ space)

$$H_0^1(I) = \{ v \in H^1(I) : v(a) = v(b) = 0 \}, \quad I = (a, b).$$

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- Weak formulation: find  $u \in H^2(I)$  such that

$$\int_a^b cu'v' dx = \int_a^b fv dx$$

for any  $v \in H_0^1(I)$ .

- Let

$$a(u, v) = \int_a^b cu'v' dx, \quad (f, v) = \int_a^b fv dx.$$

- Weak formulation: find  $u \in H^2(I)$  such that

$$a(u, v) = (f, v)$$

for any  $v \in H_0^1(I)$ .

# Galerkin formulation

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- Assume there is a finite dimensional subspace  $U_h \subset H^1[a, b]$ .
- Galerkin formulation: find  $u_h \in U_h$  such that

$$a(u_h, v_h) = (f, v_h) \quad \text{i.e.,} \quad \int_a^b c u_h' v_h' dx = \int_a^b f v_h dx$$

for any  $v_h \in U_h$ .

- Basic idea of Galerkin formulation: use **finite** dimensional space to **approximate infinite** dimensional space.
- Question: How to obtain  $U_h$ ?

# Mesh

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- Assume that we have a uniform partition of  $[a, b]$  into  $N$  elements with mesh size  $h = \frac{b-a}{N}$ .
- Let  $x_i = a + (i-1)h$  ( $i = 1, \dots, N+1$ ) denote the mesh nodes.
- Let  $E_n = [x_n, x_{n+1}]$  ( $n = 1, \dots, N$ ) denote the mesh element.

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- Define 1D linear finite element space:

$$U_h = \{\phi \in C[a, b] : \phi(x) \text{ is linear on each } E_n (n = 1, 2, \dots, N)\}.$$

- $U_h$  is actually a piecewise linear function space based on the mesh generated before.



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## Theorem

$U_h$  is an  $(N + 1)$ -dimensional subspace of  $C[a, b]$ .

Proof:

- First, it is easy to verify that  $U_h$  is a subspace of  $C[a, b]$ .
- If we can find a continuous piecewise linear basis of  $N + 1$  functions for  $U_h$ , then the proof is completed.
- Consider  $\phi_j(x) \in U_h$  such that

$$\phi_j(x_i) = \delta_{ij} = \begin{cases} 0, & j \neq i, \\ 1, & j = i \end{cases}$$

for  $i, j = 1, \dots, N + 1$ .

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Continued proof:

- In fact,

$$\phi_1(x) = \begin{cases} \frac{x_2 - x}{h}, & x_1 \leq x \leq x_2, \\ 0, & \text{otherwise,} \end{cases} \quad (1)$$

$$\phi_j(x) = \begin{cases} \frac{x - x_{j-1}}{h}, & x_{j-1} \leq x \leq x_j, \\ \frac{x_{j+1} - x}{h}, & x_j \leq x \leq x_{j+1}, \\ 0, & \text{otherwise,} \end{cases} \quad (2)$$

$$\phi_{N+1}(x) = \begin{cases} \frac{x - x_N}{h}, & x_N \leq x \leq x_{N+1}, \\ 0, & \text{otherwise,} \end{cases} \quad (3)$$

- In order to show that  $\phi_j(x) (j = 1, \dots, N+1)$  form a basis of  $U_h$ , we need to show the linear independence of  $\{\phi_j\}_{j=1}^{N+1}$  and  $U_h = \text{span}\{\phi_j\}_{j=1}^{N+1}$ .

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Continued proof:

**Linear independence:** consider

$$\sum_{j=1}^{N+1} c_j \phi_j(x) = 0$$

for any  $x \in [a, b]$ .

- Let  $x = x_i (i = 1, \dots, N+1)$ , then

$$\phi_j(x_i) = \delta_{ij} \Rightarrow c_i = 0 \quad (i = 1, \dots, N+1).$$

- So  $\phi_j(x) (j = 1, \dots, N+1)$  are linearly independent.

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Continued proof:

**Span:** Given any  $f \in U_h$ , let  $c_j = f(x_j)$  and consider

$$g(x) = \sum_{j=1}^{N+1} c_j \phi_j(x) = \sum_{j=1}^{N+1} f(x_j) \phi_j(x).$$

- First,  $g(x_i) = c_i = f(x_i)$  ( $i = 1, \dots, N+1$ ).
- Second, both  $f(x)$  and  $g(x)$  are linear in each piece  $[x_i, x_{i+1}]$  ( $i = 1, \dots, N$ ).
- Hence  $f(x) = g(x)$  in each piecewise  $[x_i, x_{i+1}]$  ( $i = 1, \dots, N$ ).

- Then  $f(x) = g(x) = \sum_{j=1}^{N+1} c_j \phi_j(x).$

- This implies  $U_h = \text{span}\{\phi_j\}_{j=1}^{N+1}.$

- Therefore  $\phi_j(x)$  ( $j = 1, \dots, N+1$ ) form a basis of  $U_h$ .

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- Recall the Galerkin formulation: find  $u_h \in U_h$  such that

$$a(u_h, v_h) = (f, v_h) \Leftrightarrow \int_a^b c u_h' v_h' dx = \int_a^b f v_h dx$$

for any  $v_h \in U_h$ .

- Since  $u_h \in U_h = \text{span}\{\phi_j\}_{j=1}^{N+1}$ , then

$$u_h = \sum_{j=1}^{N+1} u_j \phi_j$$

for some coefficients  $u_j (j = 1, \dots, N+1)$ .

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- If we can set up a linear algebraic system for  $u_j (j = 1, \dots, N+1)$  and solve it, then we can obtain the finite element solution  $u_h$ .
- Therefore, we choose the test function  $v_h = \phi_i (i = 1, \dots, N+1)$ . Then the finite element formulation gives

$$\int_a^b c \left( \sum_{j=1}^{N+1} u_j \phi_j \right)' \phi_i' dx = \int_a^b f \phi_i dx, \quad i = 1, \dots, N+1$$

$$\Rightarrow \sum_{j=1}^{N+1} u_j \left( \int_a^b c \phi_j' \phi_i' dx \right) = \int_a^b f \phi_i dx, \quad i = 1, \dots, N+1$$

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- Define the stiffness matrix

$$\mathbf{A} = [a_{ij}]_{i,j=1}^{N+1} = \left[ \int_a^b c \phi_j' \phi_i' dx \right]_{i,j=1}^{N+1}$$

- Define the load vector

$$\mathbf{b} = [b_i]_{i=1}^{N+1} = \left[ \int_a^b f \phi_i dx \right]_{i=1}^{N+1}.$$

- Define the unknown vector

$$\mathbf{x} = [u_j]_{j=1}^{N+1}.$$

- Then we obtain the linear algebraic system

$$\mathbf{Ax} = \mathbf{b}.$$

- Here  $\mathbf{A}$  is symmetric positive-definite if the original elliptic equation is symmetric positive-definite.

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## Remark

- *In fact, since*

$$\phi_j(x_k) = \delta_{jk},$$

*then*

$$u_h(x_k) = \sum_{j=1}^{N+1} u_j \phi_j(x_k) = u_k.$$

- *Hence the coefficient  $u_j$  is actually the numerical solution at the node  $x_j (j = 1, \dots, N+1)$ .*
- *Once  $\mathbf{x} = [u_j]_{j=1}^{N+1}$  is obtained, the finite element solution*  
$$u_h = \sum_{j=1}^{N+1} u_j \phi_j$$
*and the numerical solution at all the mesh nodes are obtained.*



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- In this section we will first introduce the matrix and vector assembly by using a special method. In the later section “FE method”, we will discuss a different universal framework.
- From the definition of  $\phi_j (j = 1, \dots, N+1)$ , we can see that  $\phi_j$  are non-zero only on the elements adjacent to the node  $x_j$ , but 0 on all the other elements.
- This observation motivates us to think about

$$a_{ij} = \int_a^b c \phi_j' \phi_i' dx = \sum_{n=1}^N \int_{x_n}^{x_{n+1}} c \phi_j' \phi_i' dx, \quad i, j = 1, \dots, N+1.$$

- It is easy to see that most of  $\int_{x_n}^{x_{n+1}} c \phi_j' \phi_i' dx$  will be 0.
- So we only need to use numerical integration to compute those nonzero integrals.

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Case 1: when  $|i-j| > 1$ ,  $x_i$  and  $x_j$  are not neighbouring mesh nodes.

- On any element  $[x_n, x_{n+1}]$  ( $n = 1, \dots, N$ ), at least one of  $\phi_j$  and  $\phi_i$  is 0.
- Hence,

$$\int_{x_n}^{x_{n+1}} c \phi_j' \phi_i' dx = 0 \quad (n = 1, \dots, N)$$

$$\Rightarrow a_{ij} = \sum_{n=1}^N \int_{x_n}^{x_{n+1}} c \phi_j' \phi_i' dx = 0.$$

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Case 2: when  $i = j + 1 (j = 1, \dots, N)$

- the only element, on which both  $\phi_j$  and  $\phi_i$  are not zero, is  $[x_j, x_{j+1}]$ .
- Hence,

$$\int_{x_n}^{x_{n+1}} c \phi'_j \phi'_i dx = 0 (n = 1, \dots, j-1, j+1, \dots, N)$$

$$\Rightarrow a_{ij} = \sum_{n=1}^N \int_{x_n}^{x_{n+1}} c \phi'_j \phi'_i dx = \int_{x_j}^{x_{j+1}} c \phi'_j \phi'_i dx$$

$$\Rightarrow a_{j+1,j} = \int_{x_j}^{x_{j+1}} c(x) \left( \frac{x_{j+1} - x}{h} \right)' \left( \frac{x - x_j}{h} \right)' dx = -\frac{1}{h^2} \int_{x_j}^{x_{j+1}} c(x) dx.$$

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Case 3: when  $i = j-1 (j = 2, \dots, N+1)$

- the only element, on which both  $\phi_j$  and  $\phi_i$  are not zero, is  $[x_{j-1}, x_j]$ .
- Hence,

$$\int_{x_n}^{x_{n+1}} c \phi_j' \phi_i' dx = 0 \quad (n = 1, \dots, i-1, i+1, \dots, N)$$

$$\Rightarrow a_{ij} = \sum_{n=1}^N \int_{x_n}^{x_{n+1}} c \phi_j' \phi_i' dx = \int_{x_{j-1}}^{x_j} c \phi_j' \phi_i' dx$$

$$\Rightarrow a_{j-1,j} = \int_{x_{j-1}}^{x_j} c(x) \left( \frac{x - x_{j-1}}{h} \right)' \left( \frac{x_j - x}{h} \right)' dx = -\frac{1}{h^2} \int_{x_{j-1}}^{x_j} c(x) dx.$$

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Case 4: when  $i = j (j = 2, \dots, N)$

- the only elements, on which both  $\phi_j$  and  $\phi_i$  are not zero, are  $[x_{j-1}, x_j]$  and  $[x_j, x_{j+1}]$ .
- Hence,

$$\begin{aligned} & \int_{x_n}^{x_{n+1}} c \phi_j' \phi_i' dx = 0 \quad (n = 1, \dots, j-2, j+1, \dots, N) \\ \Rightarrow a_{ij} &= \sum_{n=1}^N \int_{x_n}^{x_{n+1}} c \phi_j' \phi_i' dx = \int_{x_{j-1}}^{x_j} c \phi_j' \phi_i' dx + \int_{x_j}^{x_{j+1}} c \phi_j' \phi_i' dx \\ \Rightarrow a_{jj} &= \int_{x_{j-1}}^{x_j} c(x) \left( \frac{x - x_{j-1}}{h} \right)' \left( \frac{x - x_{j-1}}{h} \right)' dx \\ & \quad + \int_{x_j}^{x_{j+1}} c(x) \left( \frac{x_{j+1} - x}{h} \right)' \left( \frac{x_{j+1} - x}{h} \right)' dx \\ &= \frac{1}{h^2} \int_{x_{j-1}}^{x_j} c(x) dx + \frac{1}{h^2} \int_{x_j}^{x_{j+1}} c(x) dx. \end{aligned}$$

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Case 5: when  $i = j = 1$

- the only element, on which both  $\phi_j$  and  $\phi_i$  are not zero, are  $[x_1, x_2]$ .
- Hence,

$$\begin{aligned} \int_{x_n}^{x_{n+1}} c \phi_1' \phi_1' dx &= 0 (n = 2, \dots, N) \\ \Rightarrow a_{11} &= \sum_{n=1}^N \int_{x_n}^{x_{n+1}} c \phi_1' \phi_1' dx = \int_{x_1}^{x_2} c \phi_1' \phi_1' dx \\ \Rightarrow a_{jj} &= \int_{x_1}^{x_2} c(x) \left( \frac{x_2 - x}{h} \right)' \left( \frac{x_2 - x}{h} \right)' dx \\ &= \frac{1}{h^2} \int_{x_1}^{x_2} c(x) dx. \end{aligned}$$

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Case 6: when  $i = j = N + 1$

- the only elements, on which both  $\phi_j$  and  $\phi_i$  are not zero, is  $[x_N, x_{N+1}]$ .
- Hence,

$$\int_{x_n}^{x_{n+1}} c \phi'_{N+1} \phi'_{N+1} dx = 0 \quad (n = 1, \dots, j-2, j+1, \dots, N-1)$$

$$\Rightarrow a_{N+1, N+1} = \sum_{n=1}^N \int_{x_n}^{x_{n+1}} c \phi'_{N+1} \phi'_{N+1} dx = \int_{x_N}^{x_{N+1}} c \phi'_{N+1} \phi'_{N+1} dx$$

$$\begin{aligned} \Rightarrow a_{N+1, N+1} &= \int_{x_N}^{x_{N+1}} c(x) \left( \frac{x - x_N}{h} \right)' \left( \frac{x - x_N}{h} \right)' dx \\ &= \frac{1}{h^2} \int_{x_N}^{x_{N+1}} c(x) dx. \end{aligned}$$

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- From the above discussion, we can see that most of the elements  $a_{ij}(i, j = 1, \dots, N+1)$  are 0.
- Hence the stiffness matrix  $\mathbf{A}$  is a sparse matrix.
- We can also see that we only need to compute the integrals on local elements instead of the whole domain, which later will lead to the “local assembly” idea of finite elements.



# Assembly of the stiffness matrix

## Algorithm 1

- Initialize the matrix:  $\mathbf{A} = \text{sparse}(N + 1, N + 1)$ ;
- Compute the integrals and assemble them into  $\mathbf{A}$ :

FOR  $j = 1, \dots, N + 1$ :

IF  $j \leq N$ , THEN

$$\text{Compute } A(j + 1, j) = -\frac{1}{h^2} \int_{x_j}^{x_{j+1}} c(x) dx$$

END

IF  $j \geq 2$ , THEN

$$\text{Compute } A(j - 1, j) = -\frac{1}{h^2} \int_{x_{j-1}}^{x_j} c(x) dx$$

END

IF  $2 \leq j \leq N$ , THEN

$$\text{Compute } A(j, j) = \frac{1}{h^2} \int_{x_{j-1}}^{x_{j+1}} c(x) dx$$

END

END

$$\text{Compute } A(1, 1) = \frac{1}{h^2} \int_{x_1}^{x_2} c(x) dx$$

$$\text{Compute } A(N + 1, N + 1) = \frac{1}{h^2} \int_{x_N}^{x_{N+1}} c(x) dx$$

# Assembly of the load vector

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The idea for the assembly of the load vector is similar. We have

$$b_i = \int_a^b f \phi_i dx = \sum_{n=1}^N \int_{x_n}^{x_{n+1}} f \phi_i dx, i = 1, \dots, N+1.$$

- Case 1: when  $2 \leq i \leq N$ , the only two elements, on which  $\phi_i$  is not zero, are  $[x_{i-1}, x_i]$  and  $[x_i, x_{i+1}]$ . Then

$$\int_{x_n}^{x_{n+1}} f \phi_i dx = 0, \quad n = 1, \dots, i-2, i+1, \dots, N,$$

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- Case 1: when  $2 \leq i \leq N$ , the only two elements, on which  $\phi_i$  is not zero, are  $[x_{i-1}, x_i]$  and  $[x_i, x_{i+1}]$ . Then

$$\int_{x_n}^{x_{n+1}} f \phi_i dx = 0, \quad n = 1, \dots, i-2, i+1, \dots, N,$$

$$\begin{aligned} \Rightarrow b_i &= \int_{x_{i-1}}^{x_i} f \phi_i dx + \int_{x_i}^{x_{i+1}} f \phi_i dx \\ &= \int_{x_{i-1}}^{x_i} f(x) \frac{x - x_{i-1}}{h} dx + \int_{x_i}^{x_{i+1}} f(x) \frac{x_{i+1} - x}{h} dx \end{aligned}$$

# Assembly of the load vector

- Case 2: when  $i = 1$ , the only element, on which  $\phi_1$  is not zero, is  $[x_1, x_2]$ . Then

$$\int_{x_n}^{x_{n+1}} f\phi_i dx = 0, \quad n = 2, \dots, N,$$

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# Assembly of the load vector

- Case 2: when  $i = 1$ , the only element, on which  $\phi_1$  is not zero, is  $[x_1, x_2]$ . Then

$$\int_{x_n}^{x_{n+1}} f \phi_i dx = 0, \quad n = 2, \dots, N,$$

$$\Rightarrow b_1 = \int_{x_1}^{x_2} f \phi_1 dx = \int_{x_1}^{x_2} f(x) \frac{x_2 - x}{h} dx$$

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- Case 2: when  $i = 1$ , the only element, on which  $\phi_1$  is not zero, is  $[x_1, x_2]$ . Then

$$\int_{x_n}^{x_{n+1}} f\phi_i dx = 0, \quad n = 2, \dots, N,$$

$$\Rightarrow b_1 = \int_{x_1}^{x_2} f\phi_1 dx = \int_{x_1}^{x_2} f(x) \frac{x_2 - x}{h} dx$$

- Case 3: when  $i = N + 1$ , the only element, on which  $\phi_{N+1}$  is not zero, is  $[x_N, x_{N+1}]$ . Then

$$\int_{x_n}^{x_{n+1}} f\phi_{N+1} dx = 0, \quad n = 1, \dots, N-1,$$

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- Case 2: when  $i = 1$ , the only element, on which  $\phi_1$  is not zero, is  $[x_1, x_2]$ . Then

$$\int_{x_n}^{x_{n+1}} f\phi_i dx = 0, \quad n = 2, \dots, N,$$

$$\Rightarrow b_1 = \int_{x_1}^{x_2} f\phi_1 dx = \int_{x_1}^{x_2} f(x) \frac{x_2 - x}{h} dx$$

- Case 3: when  $i = N + 1$ , the only element, on which  $\phi_{N+1}$  is not zero, is  $[x_N, x_{N+1}]$ . Then

$$\int_{x_n}^{x_{n+1}} f\phi_{N+1} dx = 0, \quad n = 1, \dots, N-1,$$

$$\Rightarrow b_{N+1} = \int_{x_N}^{x_{N+1}} f\phi_{N+1} dx = \int_{x_N}^{x_{N+1}} f(x) \frac{x - x_N}{h} dx$$

## Algorithm II

- Initialize the matrix:  $\mathbf{b} = \text{zeros}(N + 1, 1)$ ;
- Compute the integrals and assemble them into  $\mathbf{b}$ :

FOR  $i = 2, \dots, N$ :

$$\text{Compute } b(i) = \int_{x_{i-1}}^{x_i} f(x) \frac{x - x_{i-1}}{h} dx + \int_{x_i}^{x_{i+1}} f(x) \frac{x_{i+1} - x}{h} dx$$

END

$$\text{Compute } b(1) = \int_{x_1}^{x_2} f(x) \frac{x_2 - x}{h} dx$$

$$\text{Compute } b(N + 1) = \int_{x_N}^{x_{N+1}} f(x) \frac{x - x_N}{h} dx$$



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Basically, the Dirichlet boundary condition  $u(a) = g_a$ ,  $u(b) = g_b$  give the solutions at  $x_1 = a$  and  $x_{N+1} = b$ .

Since the coefficient  $u_j$  in the finite element solution  $u_h = \sum_{j=1}^{N+1} u_j \phi_j$  is actually the numerical solution at the node  $x_j (j = 1, \dots, N+1)$ , we actually know that  $u_1 = u(a) = g_a$  and  $u_{N+1} = u(b) = g_b$ .

Therefore, we don't really need the first and last equations in the linear system since they are set up for  $u_1$  and  $u_{N+1}$  by using  $\phi_1$  and  $\phi_{N+1}$ .

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One way to impose the Dirichlet boundary condition is to replace the first and last equations in the linear system by the following two equations

$$u_1 = g_a \Rightarrow 1 \cdot u_1 + 0 \cdot u_2 + \cdots + 0 \cdot u_{N+1} = g_a,$$

$$u_{N+1} = g_b \Rightarrow 0 \cdot u_1 + \cdots + 0 \cdot u_N + 1 \cdot u_{N+1} = g_b.$$

That is, the first and last rows of the matrix **A** should become

$$(1, 0, \cdots, 0) \text{ and } (0, \cdots, 0, 1)$$

respectively, and the first and last elements of the vector **b** should become  $g_a$  and  $g_b$  respectively.

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- Deal with the Dirichlet boundary conditions:

$$A(1, :) = 0;$$

$$A(1, 1) = 1;$$

$$A(N+1, :) = 0;$$

$$A(N+1, N+1) = 0;$$

$$b(1) = g_a;$$

$$b(N+1) = g_b;$$

# Basic algorithm

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- Input  $a, b$  and  $N$ , compute  $h = (b - a)/N$  and  $x_j = a + (j - 1)h$  ( $j = 1, \dots, N + 1$ ).
- Assemble the stiffness matrix  $\mathbf{A}$  by using Algorithm I.
- Assemble the load vector  $\mathbf{b}$  by using Algorithm II.
- Deal with the Dirichlet boundary condition by using Algorithm III.
- Solve  $\mathbf{Ax} = \mathbf{b}$  for  $\mathbf{x}$  by using a direct or iterative method.

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## Remark

*The above algorithm uses the Algorithms I, II and III, which are designed for some particular cases with a special method. It is not general enough to deal with different types of PDEs. Therefore, we will discuss a more universal framework in the following.*

# Universal framework of FEM

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- Generate the information matrix:  $P, T, E$ ;
- Assemble the matrices and vectors: local assembly based on  $P, T, E$  only;
- Deal with the boundary conditions: boundary information matrix and local assembly;
- Solve linear systems.

# Mesh information matrices

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- Define your global indices for all the mesh elements and mesh nodes. Let  $N$  denote the number of mesh elements and  $N_m$  denote the number of mesh nodes. Here  $N_m = N + 1$ .
- Define matrix  $\mathbf{P}$  to be an information matrix consisting the coordinates of all mesh nodes.
- Define matrix  $\mathbf{T}$  to be an information matrix consisting of the global node indices of the mesh nodes of all the mesh elements.

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For example, for the mesh used in this chapter, we can use the  $j^{th}$  column of the matrix  $\mathbf{P}$  to store the coordinates of the  $j^{th}$  mesh node and the  $n^{th}$  column of the matrix  $\mathbf{T}$  to store the global node indices of the mesh nodes of the  $n^{th}$  mesh elements:

$$\begin{aligned}\mathbf{P} &= \begin{pmatrix} x_1 & x_2 & \cdots & x_{N_m-1} & x_{N_m} \end{pmatrix} \\ &= \begin{pmatrix} x_1 & x_2 & \cdots & x_N & x_{N+1} \end{pmatrix}\end{aligned}$$

$$\begin{aligned}\mathbf{T} &= \begin{pmatrix} 1 & 2 & \cdots & N_m-2 & N_m-1 \\ 2 & 3 & \cdots & N_m-1 & N_m \end{pmatrix} \\ &= \begin{pmatrix} 1 & 2 & \cdots & N-1 & N \\ 2 & 3 & \cdots & N & N+1 \end{pmatrix}\end{aligned}$$

Note: The above mesh information matrices  $\mathbf{P}$  and  $\mathbf{T}$  are for the mesh nodes.



# Finite element information matrices

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We also need similar finite element information matrices  $\mathbf{P}_b$  and  $\mathbf{T}_b$  for the **finite elements nodes**, which are the nodes corresponding to the **finite basis function**.

For example, the finite element nodes of the linear finite element are the same as those mesh nodes since all the linear basis functions are corresponding to mesh nodes.

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We also need similar finite element information matrices  $P_b$  and  $T_b$  for the **finite elements nodes**, which are the nodes corresponding to the **finite basis function**.

For example, the finite element nodes of the linear finite element are the same as those mesh nodes since all the linear basis functions are corresponding to mesh nodes.

**Note:** For the nodal finite element basis functions, the correspondence between the finite elements nodes and the finite element basis nodes is one-to-one in a straightforward way. But it could be more complicated for other types of finite element basis functions in the future.

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- Define your global indices for all the mesh elements and finite element nodes (or the FE basis functions). Let  $N_b$  denote the total number of the FE basis functions (= the number of unknowns = the total number of the FE nodes). Here,  $N_b = N + 1$ .

- Then

$$u_h = \sum_{j=1}^{N_b} u_j \phi_j.$$

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- Define matrix  $\mathbf{P}_b$  to be an information matrix consisting of the **coordinates** of all **FE nodes**;
- Define matrix  $\mathbf{T}_b$  to be an information matrix consisting of the **global node indices** of the **FE nodes** of all the mesh elements.

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For example, for the linear FE we use here,  $\mathbf{P}_b = \mathbf{P}$  and  $\mathbf{T}_b = \mathbf{T}$  since the nodes of the linear FE basis functions are the same as those of the mesh. We can use the  $j^{th}$  column of the matrix  $\mathbf{P}_b$  to store the coordinates of the  $j^{th}$  FE node and the  $n^{th}$  column of the matrix  $\mathbf{T}_b$  to store the global node indices of the FE nodes of the  $n^{th}$  mesh elements:

$$\begin{aligned}\mathbf{P}_b &= \begin{pmatrix} x_1 & x_2 & \cdots & x_{N_b-1} & x_{N_b} \end{pmatrix} \\ &= \begin{pmatrix} x_1 & x_2 & \cdots & x_N & x_{N+1} \end{pmatrix}\end{aligned}$$

$$\begin{aligned}\mathbf{T}_b &= \begin{pmatrix} 1 & 2 & \cdots & N_b-2 & N_b-1 \\ 2 & 3 & \cdots & N_b-1 & N_b \end{pmatrix} \\ &= \begin{pmatrix} 1 & 2 & \cdots & N-1 & N \\ 2 & 3 & \cdots & N & N+1 \end{pmatrix}\end{aligned}$$

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## Remark

*For many types of FE, such as quadratic elements which will be discussed later and some elements which will be introduced in the next chapter,  $\mathbf{P}_b$  and  $\mathbf{T}_b$  are different from  $\mathbf{P}$  and  $\mathbf{T}$  since the nodes for the FE basis functions are different from those of the mesh.*

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Observation based on Algorithm I:

- All the non-zero entries in the stiffness matrix  $\mathbf{A}$  comes from the non-zero local integrals defined on the mesh elements;
- In each non-zero local integral, the trial and test basis functions are only corresponding to the nodes of the element which is the integral interval.
- On each element, all the local integrals, whose trial and test basis functions are corresponding to the nodes of this element, have non-trivial contribution to some non-zero entries of the stiffness matrix  $\mathbf{A}$ .

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## New assembly idea for the stiffness matrix $\mathbf{A}$ :

- Loop over all the mesh elements;
- Compute all non-zero local integrals on each element for  $\mathbf{A}$ ;
- Assemble these non-zero local integrals into the corresponding entries of  $\mathbf{A}$ .



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Compute all non-zero local integrals on each element for  $\mathbf{A}$ :

- On the  $n^{\text{th}}$  element  $E_n = [x_n, x_{n+1}]$ , we get non-zero local integrals only when the trial and test basis functions are corresponding to the FE nodes of the element.
- That is, we only consider the trial and test basis functions to be  $\phi_n$  or  $\phi_{n+1}$ .
- These are only four non-zero local integrals on  $E_n$  with the global basis function  $\phi_n$  and  $\phi_{n+1}$ :

$$\begin{aligned} \int_{x_n}^{x_{n+1}} c \phi'_n \phi'_n dx, & \quad \int_{x_n}^{x_{n+1}} c \phi'_{n+1} \phi'_n dx, \\ \int_{x_n}^{x_{n+1}} c \phi'_n \phi'_{n+1} dx, & \quad \int_{x_n}^{x_{n+1}} c \phi'_{n+1} \phi'_{n+1} dx. \end{aligned}$$

- They can be rewritten as

$$\int_{x_n}^{x_{n+1}} c \phi'_j \phi'_i dx (i, j = n, n+1).$$

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Recall

$$\phi_1(x) = \begin{cases} \frac{x_2 - x}{h}, & x_1 \leq x \leq x_2, \\ 0, & \text{otherwise} \end{cases}$$

$$\phi_j(x) = \begin{cases} \frac{x - x_{j-1}}{h}, & x_{j-1} \leq x \leq x_j, \\ \frac{x_{j+1} - x}{h}, & x_j \leq x \leq x_{j+1}, \\ 0, & \text{otherwise} \end{cases}$$

$$\phi_{N+1}(x) = \begin{cases} \frac{x - x_N}{h}, & x_N \leq x \leq x_{N+1}, \\ 0, & \text{otherwise} \end{cases}$$

Define two local linear basis functions:

$$\psi_{n1} = \phi_n|_{E_n} = \frac{x_{n+1} - x}{h}, \quad \psi_{n2} = \phi_{n+1}|_{E_n} = \frac{x - x_n}{h},$$

so in one element, the number of local basis functions  $N_{lb}=2$ .

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Then the only four non-zero local integrals become

$$\begin{aligned} \int_{x_n}^{x_{n+1}} c \psi'_{n1} \psi'_{n1} dx, & \quad \int_{x_n}^{x_{n+1}} c \psi'_{n2} \psi'_{n1} dx, \\ \int_{x_n}^{x_{n+1}} c \psi'_{n1} \psi'_{n2} dx, & \quad \int_{x_n}^{x_{n+1}} c \psi'_{n2} \psi'_{n2} dx. \end{aligned}$$

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Then the only four non-zero local integrals become

$$\begin{aligned} \int_{x_n}^{x_{n+1}} c \psi'_{n1} \psi'_{n1} dx, \quad \int_{x_n}^{x_{n+1}} c \psi'_{n2} \psi'_{n1} dx, \\ \int_{x_n}^{x_{n+1}} c \psi'_{n1} \psi'_{n2} dx, \quad \int_{x_n}^{x_{n+1}} c \psi'_{n2} \psi'_{n2} dx. \end{aligned}$$

That is, instead of the original four non-zero local integrals with the global basis functions  $\phi_n$  and  $\phi_{n+1}$ , we will compute the following four non-zero local integrals with the local basis functions  $\psi_{n1}$  and  $\psi_{n2}$ :

$$\int_{x_n}^{x_{n+1}} c \psi'_{n\alpha} \psi'_{n\beta} dx (\alpha, \beta = 1, 2)$$

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Question: How to compute these integrals?

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Question: How to compute these integrals?

**Gauss quadrature.** The needed information is stored in the matrices  $\mathbf{P}$  and  $\mathbf{T}$ .

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Assemble the non-zero local integrals into  $\mathbf{A}$ :

Based on Algorithm I, when the trial function is  $\phi_j$  and the test function is  $\phi_j$ , the corresponding non-zero local integrals should be assembled to  $a_{ij}$ .

For example,

- $\int_{x_n}^{x_{n+1}} c \phi'_n \phi'_n dx$  should be assembled to  $a_{nn}$ ;
- $\int_{x_n}^{x_{n+1}} c \phi'_{n+1} \phi'_n dx$  should be assembled to  $a_{n,n+1}$ ;
- $\int_{x_n}^{x_{n+1}} c \phi'_n \phi'_{n+1} dx$  should be assembled to  $a_{n+1,n}$ ;
- $\int_{x_n}^{x_{n+1}} c \phi'_{n+1} \phi'_{n+1} dx$  should be assembled to  $a_{n+1,n+1}$ .

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Therefore, if we find the global node indices of the trial and test basis functions, we can easily locate where to assemble a non-zero local integral.



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Therefore, if we find the global node indices of the trial and test basis functions, we can easily locate where to assemble a non-zero local integral.

Question: Since we compute

$$\int_{x_n}^{x_{n+1}} c \psi'_{n\alpha} \psi'_{n\beta} dx (\alpha, \beta = 1, 2)$$

instead of

$$\int_{x_n}^{x_{n+1}} c \phi'_j \phi'_i dx (i, j = n, n+1),$$

how do we obtain the corresponding **global node indices** of the local trial and test basis functions  $\psi_{n\alpha}$  and  $\psi_{n\beta}(\alpha, \beta = 1, 2)$ ?

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Therefore, if we find the global node indices of the trial and test basis functions, we can easily locate where to assemble a non-zero local integral.

Question: Since we compute

$$\int_{x_n}^{x_{n+1}} c \psi'_{n\alpha} \psi'_{n\beta} dx (\alpha, \beta = 1, 2)$$

instead of

$$\int_{x_n}^{x_{n+1}} c \phi'_j \phi'_i dx (i, j = n, n+1),$$

how do we obtain the corresponding **global node indices** of the local trial and test basis functions  $\psi_{n\alpha}$  and  $\psi_{n\beta} (\alpha, \beta = 1, 2)$ ?

**Information matrix  $\mathbf{T}_b$ !**

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Recall that the  $n^{\text{th}}$  column of the matrix  $\mathbf{T}_b$  stores the global node indices of the FE nodes of the  $n^{\text{th}}$  mesh element:

$$\mathbf{T}_b = \begin{pmatrix} 1 & 2 & \cdots & N-1 & N \\ 2 & 3 & \cdots & N & N+1 \end{pmatrix}.$$

Hence  $\mathbf{T}_b(\alpha, n)$  and  $\mathbf{T}_b(\beta, n)$  give the global node indices of the local trial and test basis function  $\psi_{n\alpha}$  and  $\psi_{n\beta}(\alpha, \beta = 1, 2)$ .

That is, for  $n = 1, \dots, N$ ,

$$\int_{x_n}^{x_{n+1}} c \psi'_{n\alpha} \psi'_{n\beta} dx(\alpha, \beta = 1, 2)$$

should be assembled to  $a_{ij}$ , where

$$i = \mathbf{T}_b(\beta, n), \quad j = \mathbf{T}_b(\alpha, n).$$

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## Algorithm IV:

- Initialize the matrix:  $\mathbf{A} = \text{sparse}(N_b, N_b)$ ;
- Compute the integrals and assemble them into  $\mathbf{A}$ :

FOR  $n = 1, \dots, N$ :

FOR  $\alpha = 1, \dots, N_{lb}$ :

FOR  $\beta = 1, \dots, N_{lb}$ :

Compute  $r = \int_{x_n}^{x_{n+1}} c \psi'_{n\alpha} \psi'_{n\beta} dx$

Add  $r$  to  $\mathbf{A}(\mathbf{T}_b(\beta, n), \mathbf{T}_b(\alpha, n))$

END

END

END

# Local assembly

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Algorithm IV (alternative version):

- Initialize the matrix:  $\mathbf{A} = \text{sparse}(N_b, N_b)$ ,  $\mathbf{S} = \text{zeros}(N_{lb}, N_{lb})$ ;
- Compute the integrals and assemble them into  $\mathbf{A}$ :

FOR  $n = 1, \dots, N$ :

FOR  $\alpha = 1, \dots, N_{lb}$ :

FOR  $\beta = 1, \dots, N_{lb}$ :

Compute  $S(\beta, \alpha) = \int_{x_n}^{x_{n+1}} c \psi'_{n\alpha} \psi'_{n\beta} dx$

END

END

$\mathbf{A}(\mathbf{T}_b(:, n), \mathbf{T}_b(:, n)) = \mathbf{A}(\mathbf{T}_b(:, n), \mathbf{T}_b(:, n)) + \mathbf{S}$ ;

END

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To make a general subroutine for different cases, more information are needed for computing and assembling the integral should be treated as input parameters or input function of this subroutine:

- the coefficient function  $c$ ;
- the Gauss quadrature points and weights for numerical integrals;
- the mesh information matrices  $\mathbf{P}$  and  $\mathbf{T}$ , which can also provide the number of the mesh elements  $N = \text{size}(\mathbf{T}, 2)$  and the number of mesh nodes  $N_m = \text{size}(\mathbf{P}, 2)$ ;
- the FE information matrices  $\mathbf{P}_b$  and  $\mathbf{T}_b$  for the trial and test functions respectively, which can also provide the number of local basis functions  $N_{lb} = \text{size}(\mathbf{T}_b, 1)$  and the number of the global basis functions  $N_b = \text{size}(\mathbf{P}_b, 2) = \text{the number of unknowns}$ ;
- the type of the basis function for the trial and test functions, respectively.