

Introduction to Finite Element Method



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Chapter 12 Weighted Residual Methods

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12.1 Weighted Residual Method

- Consider the one-dimensional differential equation

$$Lu + g = 0, \quad a \leq x \leq b$$

- L : a differential operator
- $u(x)$: unknown function
- $g(x)$: known function
- If L is chosen, it specifies the actual form of the differential equation

e.g. $L = \frac{d}{dx} \Rightarrow \frac{du}{dx} + g = 0$

$$L = \frac{d^2}{dx^2} + 1 \Rightarrow \frac{d^2u}{dx^2} + u + g = 0$$

- BCs

$$u(a) = u_a, \quad u(b) = u_b$$

- u_a, u_b : known quantities

- Only certain problems can be solved analytically in terms of exact closed-form solutions

12.1 Weighted Residual Method

- To seek methods for solving the above differential equation for arbitrary expressions of L and g , consider the following approximation procedure:
 - **weighted integral**
 - $\int_a^b v(Lu + g)dx = 0$
 - where v is arbitrary
 - This is **not** a weak formulation unless an integration by parts is performed to reduce the order of differentiation of the unknown function

12.1 Weighted Residual Method

- Assume the approximation fulfilling the BCs
 - $u^{app} = \psi_1 a_1 + \psi_2 a_2 + \dots + \psi_n a_n$
 - a_1, a_2, \dots, a_n : unknown parameters
 - $\psi_1, \psi_2, \dots, \psi_n$: trial functions, specified in advance
 - Trial functions are functions of x , may be taken as any approximation and this is why they are termed trial functions
 - Once a_1, a_2, \dots, a_n are known, the approximate solution is given by u^{app}
 - $u^{app} = \boldsymbol{\psi} \mathbf{a}$
 - $\boldsymbol{\psi} = [\psi_1 \quad \psi_2 \quad \dots \quad \psi_n]$, $\mathbf{a} = [a_1 \quad a_2 \quad \dots \quad a_n]^T$

12.1 Weighted Residual Method

- Recall

$$- \int_a^b v(Lu + g)dx = 0$$

- Substitute u by u^{app}

$$- \int_a^b v(Lu^{app} + g)dx = 0$$

- u^{app} will not satisfy the equation exactly in general

$$- Lu^{app} + g = e$$

- e : residual, a measure for the error

$$- \int_a^b v e dx = 0 \quad (1)$$

- The residual $e(x)$ is given a certain weight $v(x)$
- The integral of the weighted residual $v(x)e(x)$ over the region of interest is required to be zero

12.1 Weighted Residual Method

- Consider the general form of the weight function
 - $v = V_1 c_1 + V_2 c_2 + \dots + V_n c_n$
 - c_1, c_2, \dots, c_n : certain parameters
 - V_1, V_2, \dots, V_n : known functions of x , specified in advance
 - The numbers of terms in v and u^{app} are the same
 - $v = \mathbf{V}\mathbf{c} = (\mathbf{V}\mathbf{c})^T = \mathbf{c}^T \mathbf{V}^T \quad (2)$
 - $\mathbf{V} = [V_1 \quad V_2 \quad \dots \quad V_n]$, $\mathbf{c} = [c_1 \quad c_2 \quad \dots \quad c_n]^T$
 - As the weight function v is arbitrary and \mathbf{V} is known, \mathbf{c} is arbitrary
 - $(2) \Rightarrow (1): \mathbf{c}^T \int_a^b \mathbf{V}^T e dx = 0$
 - which holds for arbitrary \mathbf{c}^T
 - $\int_a^b \mathbf{V}^T e dx = 0 \quad (3)$

12.1 Weighted Residual Method

- $\int_a^b \mathbf{V}^T e dx = 0 \quad (3)$

- The column matrix \mathbf{V}^T has the dimension $n \times 1$

$$\int_a^b V_1 e dx = 0$$

$$\int_a^b V_2 e dx = 0 \quad (4)$$

⋮

$$\int_a^b V_n e dx = 0$$

- The residual e depends on the unknowns a_1, a_2, \dots, a_n . Thus, eq.(4) serves as a system of equations to determine the n unknowns

- $e = L(\boldsymbol{\psi} \mathbf{a}) + g = L(\boldsymbol{\psi}) \mathbf{a} + g \quad (5)$

- where \mathbf{a} is independent of x

12.1 Weighted Residual Method

- (5) \Rightarrow (3): $\int_a^b \mathbf{V}^T L(\boldsymbol{\psi}) dx \mathbf{a} = -\int_a^b \mathbf{V}^T g dx$ $\int_a^b \mathbf{V}^T e dx = 0$ (3)
- Define $\mathbf{K} = \int_a^b \mathbf{V}^T L(\boldsymbol{\psi}) dx$ $e = L(\boldsymbol{\psi}) \mathbf{a} + g$ (5)
- $\mathbf{f} = -\int_a^b \mathbf{V}^T g dx \quad \Rightarrow \mathbf{K} \mathbf{a} = \mathbf{f}$

$$\mathbf{K} = \begin{bmatrix} \int_a^b V_1 L(\psi_1) dx & \int_a^b V_1 L(\psi_2) dx & \cdots & \int_a^b V_1 L(\psi_n) dx \\ \int_a^b V_2 L(\psi_1) dx & \int_a^b V_2 L(\psi_2) dx & \cdots & \int_a^b V_2 L(\psi_n) dx \\ \vdots & \vdots & & \vdots \\ \int_a^b V_n L(\psi_1) dx & \int_a^b V_n L(\psi_2) dx & \cdots & \int_a^b V_n L(\psi_n) dx \end{bmatrix}, \quad \mathbf{f} = - \begin{bmatrix} \int_a^b V_1 g dx \\ \int_a^b V_2 g dx \\ \vdots \\ \int_a^b V_n g dx \end{bmatrix}$$

- \mathbf{K} is a square matrix with dimension $n \times n$
- The system consists of n linear equations from which the n unknowns a_1, a_2, \dots, a_n , i.e. \mathbf{a} , can be determined
- When \mathbf{a} is obtained, $u^{app} = \boldsymbol{\psi} \mathbf{a}$ provides the approximate solution

12.1 Weighted Residual Method

- The procedure described above applies to all weighted residual methods
- A variety of **different weighted residual methods** is obtained depending on the choice of the weight function v , i.e. the **choice for V**

12.2 Point Collocation Method

- Dirac delta function

- $\delta(x - x_i) = \begin{cases} \infty & \text{if } x = x_i \\ 0 & \text{otherwise} \end{cases} \quad (6)$

- $\int_{-\infty}^{\infty} \delta(x - x_i) dx = 1$

- x_i : a given fixed value

- Alternatively, $\int_{x_i^-}^{x_i^+} \delta(x - x_i) dx = 1$

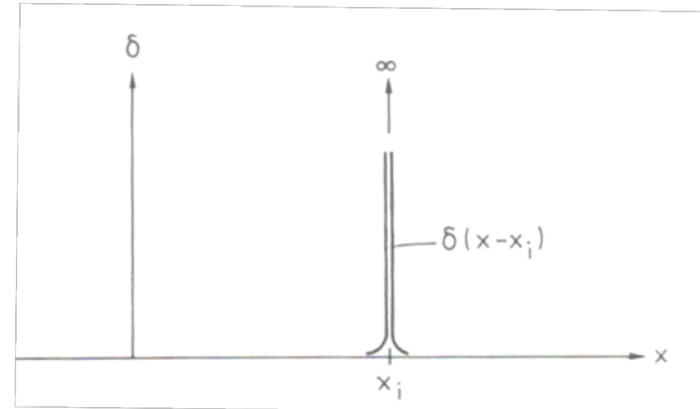
- x_i^+ and x_i^- denote x -values slightly larger than and smaller than x_i

- The weight function v is chosen such that

- $V = [\delta(x - x_1) \quad \delta(x - x_2) \quad \dots \quad \delta(x - x_n)]$

- The fixed points x_1, x_2, \dots, x_n are collocation points chosen arbitrarily within the region $a \leq x \leq b$

- $\int_a^b V_i e(x) dx = \int_a^b \delta(x - x_i) e(x) dx = 0, \quad i = 1, 2, \dots, n$



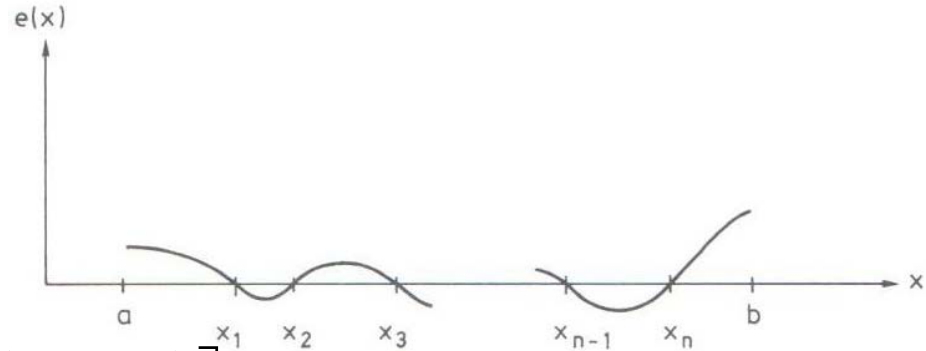
12.2 Point Collocation Method

- As Dirac delta function is zero unless $x = x_i$, we have

$$\int_a^b \delta(x - x_i) e(x) dx = \int_{x_i^-}^{x_i^+} \delta(x - x_i) e(x) dx = e(x_i) \int_{x_i^-}^{x_i^+} \delta(x - x_i) dx = e(x_i) \quad (7)$$

$$\int_a^b V_i e(x) dx = e(x_i) = 0, \quad i = 1, 2, \dots, n$$

- By analogy,



$$\mathbf{K} = \begin{bmatrix} L(\psi_1(x_1)) & L(\psi_2(x_1)) & \cdots & L(\psi_n(x_1)) \\ L(\psi_1(x_2)) & L(\psi_2(x_2)) & \cdots & L(\psi_n(x_2)) \\ \vdots & \vdots & & \vdots \\ L(\psi_1(x_n)) & L(\psi_2(x_n)) & \cdots & L(\psi_n(x_n)) \end{bmatrix}, \quad \mathbf{f} = - \begin{bmatrix} g(x_1) \\ g(x_2) \\ \vdots \\ g(x_n) \end{bmatrix}$$

- The differentiation indicated by the operator L , for instance $L(\psi_1(x_2))$, should be understood as follows: the differentiation of the function $\psi_1(x)$ is first carried out and then the value $x = x_2$ is inserted

Example

- Consider the following differential equation

$$\frac{d^2u}{dx^2} + u + x = 0, 0 \leq x \leq 1$$

BCs: $u(0) = u(1) = 0$

The exact solution is $u = \frac{\sin x}{\sin 1} - x$

- Find the approximation u^{app} to the problem
 - Recall $Lu + g = 0$
 - $L = \frac{d^2}{dx^2} + 1, g = x$

Example

- Express the approximation in the following form:

$$u^{app} = \psi_1 a_1 + \psi_2 a_2 + \cdots + \psi_n a_n$$

- Consider

$$u^{app} = b_0 + (a_1 \sin cx + b_1 \cos cx) + (a_2 \sin 2cx + b_2 \cos 2cx) + \dots + (a_n \sin ncx + b_n \cos ncx)$$

– BCs

- $u^{app}(0) = 0 \Rightarrow b_0 = b_1 = b_2 = \cdots = b_n = 0$
- $u^{app}(1) = 0 \Rightarrow c = \pi$

– Thus,

$$u^{app} = a_1 \sin \pi x + a_2 \sin 2\pi x + \cdots + a_n \sin n\pi x$$

Example

- Consider one term in the series as a simple approximation
 - $u^{app} = a_1 \sin \pi x$
 - or $u^{app} = \psi_1 a_1 = \boldsymbol{\psi} \boldsymbol{a}$
 - $\boldsymbol{\psi} = [\psi_1] = [\sin \pi x]$
 - $\boldsymbol{a} = [a_1]$
- Use point collocation method to determine a_1

Example

- Recall

- $L = \frac{d^2}{dx^2} + 1, g = x$

- $\psi_1 = \sin \pi x$

- $\mathbf{K} \mathbf{a} = \mathbf{f}$

$$\mathbf{K} = \left[L(\psi_1(x_1)) \right] = \left[-\pi^2 \sin \pi x_1 + \sin \pi x_1 \right] = \left[(-\pi^2 + 1) \sin \pi x_1 \right]$$

$$\mathbf{f} = -\left[g(x_1) \right] = -\left[x_1 \right]$$

- Choose the collocation point, for instance, the midpoint of the interval $x_1 = \frac{1}{2}$

$$\left(-\pi^2 + 1 \right) \sin \frac{\pi}{2} a_1 = -\frac{1}{2} \Rightarrow a_1 = \frac{1}{2(\pi^2 - 1)} \approx 0.0564$$

12.3 Subdomain Collocation Method

- In the point collocation method, n points are chosen
- In the subdomain collocation method, the region is divided into n subregions
 - Each subregion is given by $x_i \leq x \leq x_{i+1}$, where both x_i and x_{i+1} are located in the region $a \leq x \leq b$
 - V is chosen such that
 - $V_i = \begin{cases} 1 & \text{if } x_i \leq x \leq x_{i+1} \\ 0 & \text{otherwise} \end{cases} \quad i = 1, 2, \dots, n$

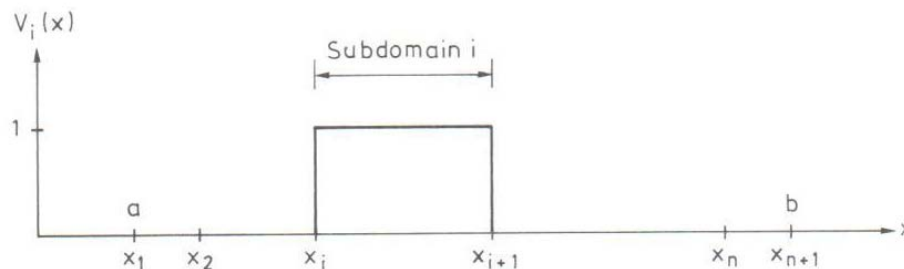
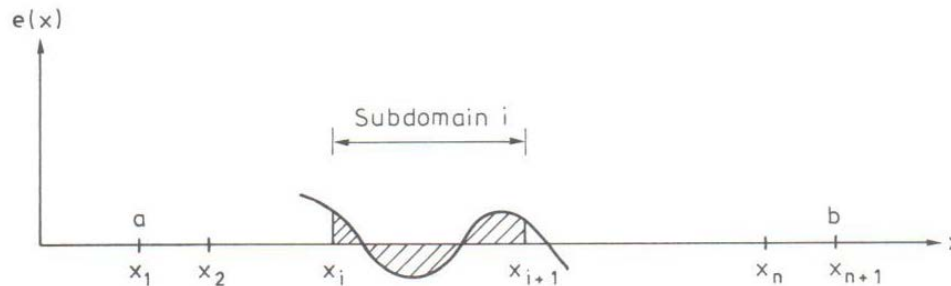


Figure 8.3 Weight function in subdomain collocation method

12.3 Subdomain Collocation Method

$$\int_a^b V_i e(x) dx = \int_{x_i}^{x_{i+1}} e(x) dx = 0, \quad i = 1, 2, \dots, n$$



- The average of the residual over each subdomain is forced to be equal to zero

$$\mathbf{K} = \begin{bmatrix} \int_{x_1}^{x_2} L(\psi_1) dx & \int_{x_1}^{x_2} L(\psi_2) dx & \cdots & \int_{x_1}^{x_2} L(\psi_n) dx \\ \int_{x_2}^{x_3} L(\psi_1) dx & \int_{x_2}^{x_3} L(\psi_2) dx & \cdots & \int_{x_2}^{x_3} L(\psi_n) dx \\ \vdots & \vdots & & \vdots \\ \int_{x_n}^{x_{n+1}} L(\psi_1) dx & \int_{x_n}^{x_{n+1}} L(\psi_2) dx & \cdots & \int_{x_n}^{x_{n+1}} L(\psi_n) dx \end{bmatrix}, \quad \mathbf{f} = - \begin{bmatrix} \int_{x_1}^{x_2} g dx \\ \int_{x_2}^{x_3} g dx \\ \vdots \\ \int_{x_n}^{x_{n+1}} g dx \end{bmatrix}$$

Example

- Consider the differential equation $\frac{d^2u}{dx^2} + u + x = 0$,
 $0 \leq x \leq 1$
- BCs: $u(0) = u(1) = 0$
- Use one term in the series as a simple approximation
 - $u^{app} = a_1 \sin \pi x = \boldsymbol{\psi} \boldsymbol{a}$
 - $\boldsymbol{\psi} = [\psi_1] = [\sin \pi x]$, $\boldsymbol{a} = [a_1]$
- Use subdomain collocation method to determine a_1

Example

- In this case, only one subdomain is involved and this is chosen as the entire region of interest

$$- L = \frac{d^2}{dx^2} + 1, g = x$$

$$- \psi_1 = \sin \pi x$$

$$- \mathbf{K} \mathbf{a} = \mathbf{f}$$

$$\mathbf{K} = \left[\int_0^1 L(\psi_1) dx \right] = \left[\int_0^1 (-\pi^2 + 1) \sin \pi x dx \right] = \left[\frac{2}{\pi} (-\pi^2 + 1) \right]$$

$$\mathbf{f} = - \left[\int_0^1 g dx \right] = - \left[\int_0^1 x dx \right] = - \left[\frac{1}{2} \right]$$

$$\frac{2}{\pi} (-\pi^2 + 1) a_1 = -\frac{1}{2} \Rightarrow a_1 = \frac{\pi}{4(\pi^2 - 1)} \approx 0.0885$$

12.4 Least-Squares Method

- Recall the residual $e = L(\boldsymbol{\psi})\mathbf{a} + g$ (5) $\Rightarrow e = e(x, a_1, a_2, \dots, a_n)$
- In the least-squares method, V is chosen such that

$$V_i = \frac{\partial e}{\partial a_i}, \quad i = 1, 2, \dots, n \quad (8)$$

$$\int_a^b V_i e(x) dx = \int_a^b \frac{\partial e}{\partial a_i} e(x) dx = 0, \quad i = 1, 2, \dots, n \quad (9)$$

- To evaluate this choice, consider $I = \int_a^b e^2(x, a_1, a_2, \dots, a_n) dx$
- As the integration is carried out over x , $I = I(a_1, a_2, \dots, a_n)$

$$\frac{\partial I}{\partial a_i} = 2 \int_a^b e \frac{\partial e}{\partial a_i} dx = 0, \quad i = 1, 2, \dots, n$$

- The weight function in eq.(8) implies that I is stationary. The stationary of I is a minimum.
- Consequently, the square of the error is a minimum, which gives the terminology of the least-squares method

12.4 Least-Squares Method

- $e = L(\boldsymbol{\psi})\mathbf{a} + g$ (5)

$$V_i = \frac{\partial e}{\partial a_i}, \quad i = 1, 2, \dots, n \quad (8)$$

– Insert eq.(5) into eq.(8) yields $V_i = L(\psi_i)$, $i = 1, 2, \dots, n$

$$\mathbf{K} = \begin{bmatrix} \int_a^b L(\psi_1)L(\psi_1)dx & \int_a^b L(\psi_1)L(\psi_2)dx & \cdots & \int_a^b L(\psi_1)L(\psi_n)dx \\ \int_a^b L(\psi_2)L(\psi_1)dx & \int_a^b L(\psi_2)L(\psi_2)dx & \cdots & \int_a^b L(\psi_2)L(\psi_n)dx \\ \vdots & \vdots & & \vdots \\ \int_a^b L(\psi_n)L(\psi_1)dx & \int_a^b L(\psi_n)L(\psi_2)dx & \cdots & \int_a^b L(\psi_n)L(\psi_n)dx \end{bmatrix}, \mathbf{f} = - \begin{bmatrix} \int_a^b L(\psi_1)gdx \\ \int_a^b L(\psi_2)gdx \\ \vdots \\ \int_a^b L(\psi_n)gdx \end{bmatrix}$$

– The coefficient matrix **K** is symmetric

Example

- Consider the differential equation $\frac{d^2u}{dx^2} + u + x = 0$,
 $0 \leq x \leq 1$
- BCs: $u(0) = u(1) = 0$
- Use one term in the series as a simple approximation
 - $u^{app} = a_1 \sin \pi x = \boldsymbol{\psi} \boldsymbol{a}$
 - $\boldsymbol{\psi} = [\psi_1] = [\sin \pi x]$, $\boldsymbol{a} = [a_1]$
- Use least-squares method to determine a_1

Example

- Recall

- $L = \frac{d^2}{dx^2} + 1, g = x, \psi_1 = \sin \pi x$

- $\mathbf{K} \mathbf{a} = \mathbf{f}$

$$\mathbf{K} = \left[\int_0^1 L(\psi_1) L(\psi_1) dx \right] = \left[\int_0^1 (-\pi^2 + 1)^2 \sin^2(\pi x) dx \right] = \left[\frac{1}{2} (-\pi^2 + 1)^2 \right]$$

$$\mathbf{f} = - \left[\int_0^1 L(\psi_1) g dx \right] = - \left[\int_0^1 (-\pi^2 + 1) x \sin \pi x dx \right] = - \left[\frac{1}{\pi} (-\pi^2 + 1) \right]$$

- where $\int \sin^2 x dx = \frac{x}{2} - \frac{1}{4} \sin 2x$

- Thus, $\int x \sin x dx = \sin x - x \cos x$

$$\frac{1}{2} (-\pi^2 + 1)^2 a_1 = - \frac{1}{\pi} (-\pi^2 + 1) \Rightarrow a_1 = \frac{2}{\pi (\pi^2 - 1)} \approx 0.0718$$

12.5 The Galerkin Method

- V is chosen as $V_i = \psi_i$, $i = 1, 2, \dots, n$
 - i.e., **weight (test) functions = trial functions**

$$\mathbf{K} = \begin{bmatrix} \int_a^b \psi_1 L(\psi_1) dx & \int_a^b \psi_1 L(\psi_2) dx & \cdots & \int_a^b \psi_1 L(\psi_n) dx \\ \int_a^b \psi_2 L(\psi_1) dx & \int_a^b \psi_2 L(\psi_2) dx & \cdots & \int_a^b \psi_2 L(\psi_n) dx \\ \vdots & \vdots & & \vdots \\ \int_a^b \psi_n L(\psi_1) dx & \int_a^b \psi_n L(\psi_2) dx & \cdots & \int_a^b \psi_n L(\psi_n) dx \end{bmatrix}, \quad \mathbf{f} = - \begin{bmatrix} \int_a^b \psi_1 g dx \\ \int_a^b \psi_2 g dx \\ \vdots \\ \int_a^b \psi_n g dx \end{bmatrix}$$

- In general, the coefficient matrix \mathbf{K} is not symmetric

Example

- Consider the differential equation $\frac{d^2u}{dx^2} + u + x = 0$,
 $0 \leq x \leq 1$
- BCs: $u(0) = u(1) = 0$
- Use one term in the series as a simple approximation
 - $u^{app} = a_1 \sin \pi x = \boldsymbol{\psi} \boldsymbol{a}$
 - $\boldsymbol{\psi} = [\psi_1] = [\sin \pi x]$, $\boldsymbol{a} = [a_1]$
- Use Galerkin method to determine a_1

Example

- Recall

$$- L = \frac{d^2}{dx^2} + 1, g = x, \psi_1 = \sin \pi x$$

$$- \mathbf{K} \mathbf{a} = \mathbf{f}$$

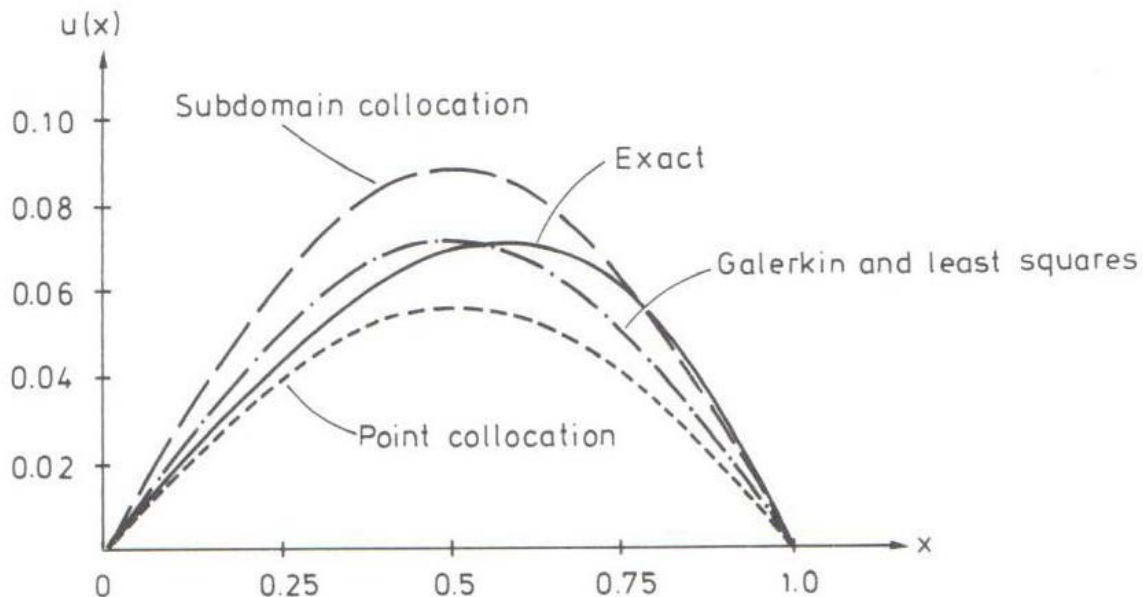
$$\mathbf{K} = \left[\int_0^1 \psi_1 L(\psi_1) dx \right] = \left[\int_0^1 (-\pi^2 + 1) \sin^2(\pi x) dx \right] = \frac{1}{2}(-\pi^2 + 1)$$

$$\mathbf{f} = - \left[\int_0^1 \psi_1 g dx \right] = - \left[\int_0^1 x \sin \pi x dx \right] = -\frac{1}{\pi}$$

$$\frac{1}{2}(-\pi^2 + 1) a_1 = -\frac{1}{\pi} \Rightarrow a_1 = \frac{2}{\pi(\pi^2 - 1)} \approx 0.0718$$

Example

- Comparison of different weighted residual methods
 - In this example, the Galerkin and least-squares methods provide the same results, and the superiority of the two methods is obvious for the example considered
 - The subdomain method is more accurate than the point collocation method



Example

- Comparison of different weighted residual methods
 - The least-squares method and the Galerkin method turn out to be very efficient
 - The least-squares method always results in a symmetric coefficient matrix, which is an advantage in numerical calculations
 - When the Galerkin method is used in combination with the weak formulation (in which an integration by parts is performed), a symmetric coefficient matrix arises