

FEM Homework Report

Yue Wu*

September 26, 2022

1 Introduction

Make a program to solve the two-point boundary value problem with purely Neumann boundary condition:

$$\begin{cases} -u'' = f, & 0 < x < 1 \\ u'(0) = u'(1) = 0 \end{cases} \quad (1)$$

Use an equidistant mesh and a piecewise linear polynomial space V_h as the finite element space. Use $f(x) = -12x^2 + 12x - 2$, $u(x) = x^2(x - 1)^2$ to test your program, and compute the following errors:

$$\|u - u_h\|_{L^2[0,1]}, \quad \|u - u_h\|_{H^1[0,1]} \quad (2)$$

2 Method

2.1 The Galerkin approximation

We use the following equidistant mesh:

$$x_j = jh, \quad j = 0, 1, \dots, N, \quad I_j = [x_{j-1}, x_j], \quad h_j = x_j - x_{j-1}, \quad h = \max_j h_j \quad (3)$$

and we use the following finite element space:

$$V_h = \{v \in C[0, 1] \mid v|_{I_j} \in \mathcal{P}^1(I_j), j = 1, \dots, N\} \quad (4)$$

The Galerkin approximation to the problem using approximation space V_h is given by the following:

$$\begin{aligned} &\text{find } u_h \in V_h, \text{ such that:} \\ &\forall v_h \in V_h, \quad (u'_h, v'_h) = (f, v_h) + u'(1)v(1) - u'(0)v(0) = (f, v_h) \end{aligned} \quad (5)$$

*School of the Gifted Young, University of Science and Technology of China, Hefei, Anhui, China.
E-mail: pilotjohnwu@mail.ustc.edu.cn

Since the boundary condition is purely Neumann, it is obvious that the solution space is an affine space. Using the Fredholm theory for second-order elliptic equations, we need to modify the problem into the following one with an Lagrange multiplier:

$$\begin{aligned} & \text{find } u_h \in V_h \text{ and } \lambda \in \mathbb{R} \text{ such that:} \\ & \forall v_h \in V_h, \begin{cases} (u'_h, v'_h) + (\lambda, v_h) = (f, v_h) \\ (1, u_h) = c \end{cases} \end{aligned} \quad (6)$$

To approximate integrals of non-polynomial functions **in computing the right-hand-side term f** , we apply the 2-point Gauss-Lobatto quadrature rule **in both cases**.

To approximate integrals of non-polynomial functions **in computing the errors**, we apply the 2-point Gauss-Legendre quadrature rule and the 2-point Gauss-Lobatto quadrature rule **respectively**.

Here, we give the 2-point Gauss-Legendre quadrature rule and the 3-point Gauss-Lobatto quadrature rule:

$$\int_{-1}^1 f(x)dx \approx f\left(-\frac{1}{\sqrt{3}}\right) + f\left(\frac{1}{\sqrt{3}}\right) \quad (7)$$

$$\int_{-1}^1 f(x)dx \approx f(-1) + f(1) \quad (8)$$

2.2 The global basis and matrix

The global (nodal) basis for the finite element space V_h is given by:

$$\phi_0(x) = \begin{cases} \frac{x_0-x}{h_1}, & x \in I_1 \\ 0, & \text{otherwise} \end{cases} \quad (9a)$$

$$\phi_j(x) = \begin{cases} \frac{x-x_{j-1}}{h_j}, & x \in I_j \\ \frac{x_j-x}{h_{j+1}}, & x \in I_{j+1} \\ 0, & \text{otherwise} \end{cases}, \quad j = 1, \dots, N-1 \quad (9b)$$

$$\phi_N(x) = \begin{cases} \frac{x-x_{N-1}}{h_N}, & x \in I_N \\ 0, & \text{otherwise} \end{cases} \quad (9c)$$

Using global basis, each function $v_h \in V_h$ has the following nodal representation:

$$v_h(x) = \sum_{k=1}^{2N-1} v_h(x_{\frac{k}{2}}) \phi_k(x) \quad (10)$$

In this manner, the global stiffness matrix is $\mathbf{A} = \left(\left(\phi'_i, \phi'_j \right) \right)_{(N+1) \times (N+1)}$ with entries $A_{i,j} = \left(\phi'_{i-1}, \phi'_{j-1} \right)$, $i, j = 1, \dots, N+1$.

Lemma 1. *The matrix \mathbf{A} is symmetric and positive semi-definite. The null space of \mathbf{A} is precisely spanned by $\mathbf{1}$ (the vector with unit entries).*

Proof. Since we are using nodal representations, it is straight-forward to check that $\mathbf{A}\mathbf{1} = \mathbf{0}$. We next show that the null space cannot be larger.

If $\mathbf{u} \notin \text{Span}\{\mathbf{1}\}$, the polynomial $u_h \in V_h$ corresponding to $\mathbf{u} \in \mathbb{R}^{N+1}$ is non-constant, thus $\mathbf{u}^T \mathbf{A} \mathbf{u} = a(u_h, u_h) > 0$. \square

2.3 The local basis and matrix

The local (nodal) shape functions within a reference element $I = [0, 1]$ are:

$$\psi^0(x) = 1 - x \quad (11a)$$

$$\psi^1(x) = x \quad (11b)$$

The local basis functions for an element I_j are:

$$\psi_j^{(k)} = \psi^{(k)}(\xi_j(x)), \quad k = 0, 1 \quad j = 1, \dots, N \quad (12)$$

with the affine transformation which maps I_j to $[0, 1]$:

$$\xi_j(x) = \frac{x - x_{j-1}}{h_j} \quad (13)$$

The local stiffness matrix $\mathbf{K}_j \in \mathbb{R}^{(N+1) \times (N+1)}$ is:

$$\mathbf{K}_j = \text{diag} \left(\mathbf{0}_{(j-1)}, h_j^{-1} \begin{pmatrix} 1 & -1 \\ -1 & 1 \end{pmatrix}, \mathbf{0}_{N-j} \right), \quad j = 1, \dots, N \quad (14)$$

and the local right-hand-side $\mathbf{F}_j \in \mathbb{R}^{2N-1}$ is:

$$\mathbf{F}_j = \begin{pmatrix} \mathbf{0}_{j-1} \\ \int_{I_j} f(x) \psi_j^0(x) dx \\ \int_{I_j} f(x) \psi_j^1(x) dx \\ \mathbf{0}_{N-j} \end{pmatrix}, \quad j = 1, \dots, N \quad (15)$$

With these local components, the global matrix and RHS is

$$\mathbf{A} = \sum_{j=1}^N \mathbf{K}_j \quad (16a)$$

$$\mathbf{F} = \sum_{j=1}^N \mathbf{F}_j \quad (16b)$$

and we use point values of u_h as the variable: $\mathbf{u} = (u_h(x_0), \dots, u_h(x_N))^T$.

2.4 Solving the saddle-point system

Since we used a Lagrange method for the purely Neumann problem, we need to solve the saddle-point system:

$$\begin{pmatrix} \mathbf{A} & \mathbf{B} \\ \mathbf{B}^T & 0 \end{pmatrix} \begin{pmatrix} \mathbf{u} \\ \lambda \end{pmatrix} = \begin{pmatrix} \mathbf{F} \\ c \end{pmatrix} \quad (17)$$

where $\mathbf{B} \in \mathbb{R}^{N+1}$ is given by:

$$\mathbf{B} = \sum_{j=1}^N \mathbf{B}_j = \sum_{j=1}^N \begin{pmatrix} \mathbf{0}_{j-1} \\ \int_{I_j} \psi_j^0(x) dx \\ \int_{I_j} \psi_j^1(x) dx \\ \mathbf{0}_{N-j} \end{pmatrix} \quad (18)$$

Theorem 1. *The saddle-point system above has a unique solution.*

Proof. We only need to show that when $\mathbf{F} = \mathbf{0}$ and $c = 0$, the above system only admits zero solution.

When $\mathbf{B}^T \mathbf{u} = 0$, we have $\mathbf{u} = \mathbf{B}_\perp \mathbf{v}$, where $\mathbf{B}_\perp \in \mathbb{R}^{(N+1) \times N}$ is orthogonal to \mathbf{B} , of rank N . Then, solving $\mathbf{A} \mathbf{u} + \lambda \mathbf{B} = \mathbf{F}$ is equivalent to minimizing $\frac{1}{2} \mathbf{u}^T \mathbf{A} \mathbf{u} - \mathbf{u}^T \mathbf{F}$, which is also to minimize $\frac{1}{2} \mathbf{v}^T \mathbf{B}^T \mathbf{A} \mathbf{B} \mathbf{v} - \mathbf{v}^T \mathbf{B}^T \mathbf{F}$. Using the null space of \mathbf{A} , we know that $\mathbf{B}^T \mathbf{A} \mathbf{B}$ is symmetric positive definite, thus the system is well-posed. \square

3 Results

We use $N = 10, 20, 40, 80$ equidistant elements for computation. By running the scripts in `main.m`, the outputs from the `matlab` program we coded are given as following:

N	L^2 error	order	H^1 error	order
10	7.7742e-04	-	2.7681e-02	-
20	1.8839e-04	2.0450	1.3151e-02	1.0738
40	4.6714e-05	2.0118	6.4853e-03	1.0199
80	1.1654e-05	2.0030	3.2313e-03	1.0051

Table 1: Table of error and order using Gauss-Legendre rule

N	L^2 error	order	H^1 error	order
10	1.5627e-03	-	4.6907e-02	-
20	3.7730e-04	2.0503	2.2641e-02	1.0509
40	9.3460e-05	2.0133	1.1216e-02	1.0134
80	2.3311e-05	2.0034	5.5946e-03	1.0034

Table 2: Table of error and order using Gauss-Lobatto rule

Remark 1. *In the test, since the exact solution $u(x)$ satisfies $\int_0^1 u(x)dx = \frac{1}{30}$, we choose $c = \frac{1}{30}$ in order to make the error checking procedure easier.*

We should emphasize that the value of c will not influence the solution, substituting c with $c + \Delta c$ will only output $u(x) + \Delta c/m([0, 1])$ instead of $u(x)$.

4 Discussion

From this experiment, we verified the theoretical 2-nd order convergence under the L^2 norm and 1-st order convergence under the H^1 norm for piecewise quadratic conforming FEM numerically. In the algorithm, we used the 2-point Gauss-Lobatto quadrature rule to approximate integration of non-polynomial functions in the right-hand-side. This does not cause the convergence order to degenerate since the quadrature is 2-th order accurate.