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Introduction

This short lecture series...

LECTURE 1

Context

Implementation and Exercises

1. Numerical Solution of Poisson's Equation via the Finite Element Method

Let us consider numerical solutions of (variable coefficient) Poisson's equation:

(2.1a)
$$-\nabla \cdot (\beta \nabla u) = f \in \Omega$$

$$(2.1b) u = p \in \partial \Omega_d$$

$$\beta \nabla u \cdot \hat{n} = q \quad \in \partial \Omega_n$$

Here, $\Omega \subset \mathbb{R}^d$ is open; $\partial \Omega_d$ and $\partial \Omega_n$ partition the boundary of Ω and define where Dirichlet (d) and Neumann (n) boundary conditions are applied, respectively; \hat{n} is the outward-pointing unit normal on $\partial \Omega_n$; $\beta, f: \Omega \to \mathbb{R}$, $p: \partial \Omega_d \to \mathbb{R}$, and $q: \partial \Omega_q \to \mathbb{R}$ are given, and β is bounded below by some positive number; and the goal is to solve for the unknown function $u: \Omega \to \mathbb{R}$. For now, we'll ignore the details of the function spaces that u, β, f, p , and q are assumed to belong in.

For the numerical solution of (2.1), we'll use a discretization based on the Finite Element Method (FEM). Briefly, an FEM discretization focuses on the weak formulation of the partial differential equation. The derivation involves multiplying the differential equation by a test function (chosen from some suitable function space), integrating by parts, and applying boundary conditions:

$$(2.2a) -\nabla \cdot (\beta \nabla u) = f$$

(2.2b)
$$\Rightarrow \int_{\Omega} -\nabla \cdot (\beta \nabla u) \, v = \int_{\Omega} f v$$

$$(2.2c) \qquad \qquad \Rightarrow \int_{\Omega} \beta \nabla u \cdot \nabla v = \int_{\Omega} f v + \int_{\partial \Omega} \left(\beta \nabla u \cdot \hat{n} \right) v$$

$$(2.2\mathrm{d}) \qquad \qquad \Rightarrow \int_{\Omega} \beta \nabla u \cdot \nabla v = \int_{\Omega} f v + \int_{\partial \Omega_n} q v$$

where the derivation from (2.2c) to (2.2d) is possible by stipulating that $v \equiv 0$ on $\partial\Omega_d$. Clearly, then, if u satisfies (2.1), then u also satisfies (2.2d) for all appropriate v. Under suitable conditions, it turns out to be the case that the converse holds as well. Since (2.2d) involves only first order derivatives in u, its analysis is often preferable to (2.1).

We then suppose that u can be approximated by a function in a finite-dimensional function space (the *(discretized) solution space*) (e.g., continuous piecewise linear functions over some tessellation of Ω), i.e., $u \approx \tilde{u} = \sum_{j} u_{j} e_{j}$ for basis elements $\{e_{j}\}$. Likewise, v is selected from some other (often related) finite-dimensional function

space (the *(discretized) test space*), thus giving a system of equations for the coefficients u_j . For example, if v is taken from the same space as \tilde{u} , and we let v be, in turn, each of the basis functions, then we obtain

(2.3)
$$\sum_{j} \left(\int_{\Omega} \beta \nabla e_{j} \cdot \nabla e_{i} \right) u_{j} = \int_{\Omega} f e_{i} + \int_{\partial \Omega_{n}} q e_{i}$$

for all i, which may be expressed as a linear system of equations:

$$(2.4) A\vec{u} = \vec{b},$$

where

(2.5a)
$$A_{ij} = \int_{\Omega} \beta \nabla e_i \cdot \nabla e_j$$

$$(2.5b) \vec{u}_i = u_i$$

(2.5c)
$$\vec{b}_i = \int_{\Omega} f e_i + \int_{\partial \Omega_n} q e_i$$

Strictly speaking, the above is only correct if v varies precisely throughout the same function space which \tilde{u} belongs to. This is generally not the case. For example, in the presence of Dirichlet boundary conditions, \tilde{u} will generally be nonzero along Ω_d , however v will be chosen from a function space which vanishes along Ω_d (hence these functions spaces will only align under homogeneous Dirichlet boundary conditions). How this manifests itself in the linear system will be investigated in the exercises.

Note that the sparsity of A is directly related to the overlap of the supports of the e_j 's. Typically we choose the e_j 's such that the support of one basis element overlaps the supports of only a small constant number of other basis elements, in which case A will have a number of nonzero entries proportional to the number of basis elements.

Let's now focus on dimension d=1 and use linear finite elements. Let Ω be an interval (a,b), and let $\{x_j\}$ be grid points in (a,b), with $a=x_0 < x_1 < \cdots < x_n = b$. Often times, the x_j 's will be equally spaced, which makes the analysis simpler, but it's not necessary. We'll take \tilde{u} (the discretized solution) and v (the discretized test function) to be continuous and piecewise linear on each segment (x_{j-1},x_j) , with v(a) and/or v(b) vanishing if a Dirichlet condition is specified at a and/or b, respectively. This space of continuous piecewise linear functions is spanned by the nodal basis functions (also called "hat" functions, because of the shape of their graphs) e_j , (uniquely) defined such that

(2.6)
$$e_{j}(x_{k}) = \delta_{jk} = \begin{cases} 1, & j = k \\ 0, & j \neq k \end{cases}$$

1.1. Exercises

Unless otherwise noted, assume dimension d=1 linear finite elements with the basis $\{e_j\}$ given by (2.6), although these results extend to higher order and higher dimensional finite elements.

(1) Here is one way to incorporate the Dirichlet boundary conditions into (2.5). For concreteness, suppose $\Omega_d = \{a\}$ (with $\Omega = (a, b)$). Then we still have $\tilde{u} = \sum_{j=0}^{n} u_j e_j$, but our space of test functions $\{v\}$ is slightly smaller, span $\{e_1, \ldots, e_n\}$, to ensure that v(a) = 0. Thus we obtain n equations of the form (2.3), plus one more equation from the Dirichlet boundary condition $(u_0 = u(a))$. This (ostensibly) gives enough equations to solve for the n+1 coefficients $\{u_j\}$.

Another conceptual way to incorporate the Dirichlet boundary conditions is to consider the function $w = u - u(a)e_0$, which satisfies a similar Poisson problem as u but now with homogeneous Dirichlet boundary conditions. Thus, the space of test functions $\{v\}$ and the solution space of w coincide, and one may use eqs (2.5) directly.

In practice, one can actually just ignore the Dirichlet boundary conditions when initially computing the entries of A and \vec{b} (i.e., just use (2.5a) and (2.5c)), then "correct" A and \vec{b} to account for the Dirichlet boundary conditions afterwards. Describe a *simple* algorithm to effect this "correction" procedure. Try to keep A symmetric.

- (2) Show that the linear system (2.4) has a unique solution if one specifies Dirichlet boundary conditions at either x=a or x=b or both (i.e., $u(a)=u_a$ and/or $u(b)=u_b$). For simplicity, you may assumed $\beta\equiv 1$. What happens if both boundary conditions are Neumann? Specifically address the conditions which \vec{b} must satisfy, how this translates into a condition on f, and "how uniquely" \tilde{u} is determined.
- (3) Compute the integral in (2.5a) when $\beta \equiv 1$. If you wish, for simplicitly, you may assume the x_j 's are equally spaced, i.e., $x_j = a + jh$, where h = (b-a)/n. What is the structure of the matrix A (e.g., describe the sparsity)?
- (4) Give an expression to reasonably approximate the integrals in (2.5a) and (2.5c) for aribtrary β , f, and q (assume β and f are sufficiently "well-behaved"). You may have to consider boundary and interior grid vertices separately.
- (5) Write some code (e.g., a Matlab script) to solve the following boundary value problem using linear finite elements:

(2.7a)
$$-((1+x)u'(x))' = (1+x)^{-2} \quad x \in (0,1)$$

(2.7b)
$$u(0) = 1$$

$$(2.7c) (1+(1)) u'(1) = -1/2$$

(6) The FEM terminology for the matrix A in (2.4) is the *(global) stiffness matrix* (the meaning of the term "global" will become obvious shortly). With dimension d=1 linear finite elements, it is relatively straightforward to compute the entries of the stiffness matrix, even given that one has to account for differences between boundary and interior grid vertices. However, the situation becomes more complicated with higher order or higher dimensional finite elements, and it becomes tedious to consider all the special boundary cases. As such, it is common practice to assemble the global stiffness matrix element by element (i.e., cell by cell), with each element contributing partially to several entries of the global stiffness

matrix. The collection of these contributions from a single element may be put into a small matrix, the *local stiffness matrix* for that element. The computation of the local stiffness matrix is identical to that of the global stiffness matrix upon replacing the integrations over Ω with integrations localized over the element. For example, back to dimension d=1 linear finite elements, the local stiffness matrix over $I_j=(x_{j-1},x_j)$ would consist of 2^2 (nonzero) entries (since only 2 basis functions are supported on I_j). Each of these entries would be added to accumulated global stiffness matrix entries:

(2.8a)
$$A_{j-1,j-1} += \int_{I_j} \beta \nabla e_{j-1} \cdot \nabla e_{j-1}$$

(2.8b)
$$A_{j-1,j}, A_{j,j-1} += \int_{I_i} \beta \nabla e_{j-1} \cdot \nabla e_j$$

$$(2.8c) A_{jj} += \int_{I_j} \beta \nabla e_j \cdot \nabla e_j$$

This suggests the following procedure to build the (global) stiffness matrix A. First, initialize the sparsity pattern of A (with zeros). Then iterate through each element and add the contributions of the element's local stiffness matrix to the corresponding entries in A.

Verify this procedure computes the same stiffness matrix as before, and modify your program from Exercise 5 to build the stiffness matrix in this fashion.

2. Neo-Hookean Elasticity with Quasistatics Evolution in Dimension 1

In this section, we will detail the numerical solution of the elasticity equations with the Neo-Hookean constitutive model. For simplicity, we will presently consider quasistatics evolution in dimension d = 1.

2.1. blah

$\begin{array}{c} \text{LECTURE 3} \\ \end{array}$ Further Reading

Bibliography