

# Poisson's Equation

## Computer Lab 2

### Poisson's Equation

Consider Poisson's equation: Solve  $u(\mathbf{x})$  from

$$-\nabla \cdot (\kappa(\mathbf{x}) \nabla u(\mathbf{x})) = f(\mathbf{x}), \quad \mathbf{x} \in \Omega, \quad (1)$$

$$\mathbf{n} \cdot \kappa(\mathbf{x}) \nabla u(\mathbf{x}) = \gamma(g(\mathbf{x}) - u(\mathbf{x})), \quad \mathbf{x} \in \partial\Omega, \quad (2)$$

where  $f(\mathbf{x})$  and  $g(\mathbf{x})$  are given functions from  $L_2(\Omega)$ ,  $\kappa(\mathbf{x}) \in L_2(\Omega)$  is a positive function and  $\gamma$  is a positive parameter, and  $\Omega$  is a polygonal domain with boundary  $\partial\Omega$  and outward pointing unit normal  $\mathbf{n}$ . The boundary condition (2) is a so called Robin boundary condition, which is a combination of Dirichlet and Neumann boundary conditions.

### Finite Element Approximation

The weak form of (1) reads: Find  $u \in H^1$  such that

$$\int_{\Omega} \kappa \nabla u \cdot \nabla v \, d\mathbf{x} + \int_{\partial\Omega} \gamma u v \, ds = \int_{\Omega} f v \, d\mathbf{x} + \int_{\partial\Omega} \gamma g v \, ds, \quad (3)$$

for all  $v \in H^1$ , where  $H^1$  is the Hilbert space  $H^1 = \{v : \|v\| + \|\nabla v\| < \infty\}$ .

Let  $V_h \subset H^1$  be the space of all continuous piecewise linear functions on a partition  $\mathcal{K} = \{K\}$  of  $\Omega$  into triangles with diameter (longest edge)  $h_K$ .

The finite element approximation of (3) takes the form: Find  $u_h \in V_h$  such that

$$\int_{\Omega} \kappa \nabla u_h \cdot \nabla v \, d\mathbf{x} + \int_{\partial\Omega} \gamma u_h v \, ds = \int_{\Omega} f v \, d\mathbf{x} + \int_{\partial\Omega} \gamma g v \, ds, \quad (4)$$

for all  $v \in V_h$ .

**Problem 1.** Derive the weak form (3).

**How to obtain the finite element solution.** Recall that the set of hat functions  $\{\varphi_i\}_{i=1}^N$ , where  $N$  is the number of nodes within the triangulation  $\mathcal{K}$ , is a basis for  $V_h$ .

The finite element approximation (4) is equivalent to

$$\int_{\Omega} \kappa \nabla u_h \cdot \nabla \varphi_i \, d\mathbf{x} + \int_{\partial\Omega} \gamma u_h \varphi_i \, ds = \int_{\Omega} f \varphi_i \, d\mathbf{x} + \int_{\partial\Omega} \gamma g \varphi_i \, ds, \quad (5)$$

for  $i = 1, 2, \dots, N$ . This is due to the fact that if (5) holds for each basis function  $\varphi_i$  independently, then it also holds for any linear combination of them, and thus for any  $v \in V_h$ .

Since  $u_h \in V_h$  it can be written as the sum

$$u_h = \sum_{j=1}^N x_j \varphi_j, \quad (6)$$

for some coefficients  $x_j$  to be determined.

Inserting (6) into (5) we end up with a  $(N \times N)$  system of linear equations

$$\sum_{j=1}^N x_j \left( \int_{\Omega} \kappa \nabla \varphi_j \cdot \nabla \varphi_i \, d\mathbf{x} + \int_{\partial\Omega} \gamma \varphi_j \varphi_i \, ds \right) \quad (7)$$

$$= \int_{\Omega} f \varphi_i \, d\mathbf{x} + \int_{\partial\Omega} \gamma g \varphi_i \, ds, \quad i = 1, 2, \dots, N. \quad (8)$$

In matrix form we write this as

$$(A + R)x = b + r \quad (9)$$

where the entries of the left and right hand side matrices and vectors are given by

$$A_{ij} = \int_{\Omega} \kappa \nabla \varphi_i \cdot \nabla \varphi_j \, d\mathbf{x}, \quad (10)$$

$$R_{ij} = \int_{\partial\Omega} \gamma \varphi_i \varphi_j \, ds, \quad (11)$$

$$\mathbf{x} = (x_1, x_2, \dots, x_N)^T, \quad (12)$$

$$b_j = \int_{\Omega} f \varphi_j \, d\mathbf{x}, \quad (13)$$

$$r_j = \int_{\partial\Omega} \gamma g \varphi_j \, ds. \quad (14)$$

All index runs over  $i, j = 1, 2, \dots, N$ .

## Computer Implementation

Consider a single triangle  $K$  with nodes at its three corners  $N_1 = (x_1, y_1)$ ,  $N_2 = (x_2, y_2)$ , and  $N_3 = (x_3, y_3)$ . To each node  $N_i$ ,  $i = 1, 2, 3$ , there is a hat function  $\varphi_i$  associated, which takes on the value one at node  $N_i$  and zero at the other nodes. Each hat function is a linear function (i.e., a plane) on  $K$ . Hence

$$\varphi_i = a_i + b_i x + c_i y, \quad i = 1, 2, 3, \quad (15)$$

where the coefficients  $a_i$ ,  $b_i$ , and  $c_i$ , are determined from the requirement

$$\varphi_i(N_j) = \begin{cases} 1, & i = j, \\ 0, & i \neq j. \end{cases} \quad (16)$$

Thus, for example,  $a_1$ ,  $b_1$  and  $c_1$  are given by the equations  $\varphi_1(x_1, y_1) = 1$ ,  $\varphi_1(x_2, y_2) = 0$ , and  $\varphi_1(x_3, y_3) = 0$ . In matrix form need to solve the following system to compute the coefficients of  $\varphi_1$

$$\begin{pmatrix} 1 & x_1 & y_1 \\ 1 & x_2 & y_2 \\ 1 & x_3 & y_3 \end{pmatrix} \begin{pmatrix} a_1 \\ b_1 \\ c_1 \end{pmatrix} = \begin{pmatrix} 1 \\ 0 \\ 0 \end{pmatrix}. \quad (17)$$

Note that the gradient of  $\varphi_i$  is just the constant vector  $\nabla\varphi_i = (b_i, c_i)$ .

As usual the global stiffness matrix  $A$  is formed by summing local stiffness matrices  $A^K$ , which are found by performing the integration (10) successively over each triangle  $K$ .

We get the  $(3 \times 3)$  local stiffness matrix

$$A_{ij}^K = \int_K \kappa \nabla\varphi_i \cdot \nabla\varphi_j \, d\mathbf{x} = (b_i b_j + c_i c_j) \int_K \kappa \, d\mathbf{x}, \quad i, j = 1, 2, 3. \quad (18)$$

The  $3 \cdot 3 = 9$  entries of  $A^K$  are to be assembled (at the right locations) into  $A$ . In Matlab this can be done very elegantly using vectorized operations as shown below.

```
for K = 1:size(t,2); % loop over the triangles
    nodes = t(1:3,K); % find triangle K's nodes
    AK = ...          % compute the (3 x 3) stiffness matrix AK
    A(nodes,nodes) = A(nodes,nodes)+AK; % add AK(i,j), i,j=1,2,3,
                                     % to A(nodes(i),nodes(j))
end
```

**Problem 2.** Compute  $A^K$  on the reference triangle with corners at origo,  $(1, 0)$ , and  $(0, 1)$ . Verify that one of the eigenvalues of  $A^K$  is zero. Can you explain why?

Similar to the stiffness matrix  $A$ , the load vector  $b$  is formed by summing local load vectors  $b^K$ , which are found by restricting the integration (13) to one triangle  $K$  at a time. Using one point quadrature we get the approximation

$$b_i^K = \int_K f \varphi_i \approx f(x_c, y_c) \text{area}(K)/3, \quad i = 1, 2, 3, \quad (19)$$

where  $(x_c, y_c)$  is the centroid of  $K$ .

Finally, if two nodes of triangle  $K$  lies along the domain boundary  $\partial\Omega$ , then the edge between them will contribute to the line integrals (11) and (14) associated with the boundary conditions. The boundary mass matrix  $R$  is assembled by looping over all the triangle edges lying along the domain

boundary  $\partial\Omega$ . These are the columns of the mesh data matrix **e**, c.f. the matlab command **initmesh**.

If the edge  $E$  lies between the boundary nodes  $N_1$  and  $N_2$ , then we have the local boundary mass matrix

$$R_{ij}^E = \int_E \gamma \varphi_i \varphi_j \, ds = \frac{\gamma}{6} (1 + \delta_{ij}) \text{length}(E), \quad i, j = 1, 2, \quad (20)$$

and the local boundary vector

$$r_i^E = \int_E \gamma g \varphi_i \, ds \approx \frac{\gamma}{2} g(N_i) \text{length}(E), \quad i = 1, 2. \quad (21)$$

Note that these are the same as the one-dimensional local mass matrix and load vector.

**assemble.m** The following code is a template routine for performing the assembly of the system of equations (9). Besides completing the code the user must supply the routines, **f**, **gamma**, and **g**, defining the source term  $f$ , the boundary penalty parameter  $\gamma$ , and the boundary data  $g$ . For example if  $f = \sin x \sin y$ , then **f** would look like

```
function z=f(x,y)
z=sin(x).*sin(y)
```

```
function [A,R,b,r] = assemble(p,e,t)
N = size(p,2);
A = sparse(N,N);
R = sparse(N,N);
b = zeros(N,1);
r = zeros(N,1);
% assemble stiffness matrix A, and load vector b.
for K = 1:size(t,2);
    nodes = t(1:3,K);
    x = p(1,nodes);
    y = p(2,nodes);
```

```

    area_K = polyarea(x,y);
    AK = ???
    bK = ???
    A(nodes,nodes) = A(nodes,nodes) + AK;
    b(nodes) = b(nodes) + bK;
end
% assemble boundary mass matrix R, and the vector r.
for E = 1:size(e,2)
    nodes = e(1:2,E);
    x = p(1,nodes);
    y = p(2,nodes);
    length_E = sqrt((x(1)-x(2))^2+(y(1)-y(2))^2);
    R(nodes,nodes) = R(nodes,nodes) + gamma*length_E*[2 1;1 2]/6;
    r(nodes) = r(nodes) + gamma*length_E*g(x,y)'/2;
end

```

### Problem 3.

- a) Implement the finite element solver `MyPoissonSolver` described above. Test your code by solving a problem with known solution (e.g.,  $-\Delta u = 2\pi^2 \sin(\pi x) \sin(\pi y)$  in the unit square and with  $u = 0$  on the boundary. The exact solution is  $u = \sin(\pi x) \sin(\pi y)$ ).
- b) Compute the energy norm  $\|\nabla u_h\|_{L^2(\Omega)}^2 = x^T A x$  for the problem in a), and compare with theory. Make a convergence plot of the energy norm error (i.e.,  $\|\nabla(u - u_h)\|_{L^2(\Omega)}^2$ ) versus the mesh size  $h$ . Plot the result using a log-log diagram. Use `polyfit` to compute the slope of the graph. It should be roughly 2. Why?

**Problem 4.** Solve  $-\Delta u = 1$  on the unit square  $\Omega = [0, 1] \times [0, 1]$  with the boundary conditions  $u = \cos(2\pi y)$  on the line  $x = 0$ , and  $n \cdot \nabla u = 0$  on the rest of the boundary.