# Series 2 Warmup



Numerical methods for PDEs Last edited: March 14, 2017 Due date: None at 23:59

Template codes are available on the course's webpage at https://moodle-app2.let.ethz.ch/course/view.php?id=3089.

This is a warmup problem. You do **NOT need to hand in** this problem.

# Exercise 1 Linear Finite Elements for the Poisson equation in 2D

We consider the problem

$$-\Delta u = f(\boldsymbol{x}) \quad \text{in } \Omega \subset \mathbb{R}^2 \tag{1}$$

$$u(\mathbf{x}) = 0 \quad \text{on } \partial\Omega$$
 (2)

where  $f \in L^2(\Omega)$ .

1a)

Write the variational formulation for (1)-(2).

**Solution:** We multiply both the left handside and right handside of (1) by a test function v. Applying Green's formula for integration by parts on the left handside we get:

$$-\int_{\Omega} \Delta u(\boldsymbol{x}) v(\boldsymbol{x}) = \int_{\Omega} \nabla u(\boldsymbol{x}) \cdot \nabla v(\boldsymbol{x}) \ d\boldsymbol{x} - \int_{\partial \Omega} \frac{\partial u}{\partial \boldsymbol{n}}(\boldsymbol{x}) v(\boldsymbol{x}) \ d\boldsymbol{x}.$$

Since u satisfies Dirichlet boundary conditions, the test functions belong to  $H_0^1(\Omega)$  and thus the

boundary integral in the above expression vanishes. The variational formulation results then:

Find  $u \in V = H_0^1(\Omega)$  such that

$$\int_{\Omega} \nabla u(\boldsymbol{x}) \cdot \nabla v(\boldsymbol{x}) = \int_{\Omega} f(\boldsymbol{x}) v(\boldsymbol{x}) \ d\boldsymbol{x} \text{ for all } v \in H_0^1(\Omega),$$

We solve (1)-(2) by means of linear finite elements on triangular meshes of  $\Omega$ . Let us denote by  $\varphi_N^i$ ,  $i=0,\ldots,N-1$  the finite element basis functions (hat functions) associated to the vertices of a given mesh, with  $N=N_V$  the total number of vertices. The finite element solution  $u_N$  to (1) can thus be expressed as

$$u_N(\boldsymbol{x}) = \sum_{i=0}^{N-1} \mu_i \varphi_N^i(\boldsymbol{x}), \tag{3}$$

where  $\boldsymbol{\mu} = \{\mu_i\}_{i=0}^{N-1}$  is the vector of coefficients. Notice that we don't know  $\mu_i$  if i is an interior vertex, but we know that  $\mu_i = 0$  if i is a vertex on the boundary  $\partial\Omega$ .

Hint: Here and in the following, we use zero-based indices in contrast to the lecture notes.

Inserting  $\varphi_N^i$ , i = 0, ..., N-1 as test functions in the variational formulation from subproblem **1a**) we obtain the linear system of equations

$$\mathbf{A}\boldsymbol{\mu} = \mathbf{F},\tag{4}$$

with  $\mathbf{A} \in \mathbb{R}^{N \times N}$  and  $\mathbf{F} \in \mathbb{R}^N$ .

1b)

Write an expression for the entries of  $\mathbf{A}$  and  $\mathbf{F}$  in (4).

Solution: We have

$$\mathbf{A}_{ij} = \int_{\Omega} 
abla arphi_N^j(m{x}) \cdot 
abla arphi_N^i(m{x}) \; dm{x} \quad ext{ and } \mathbf{F}_i = \int_{\Omega} f(m{x}) arphi_N^i(m{x}) \; dm{x},$$

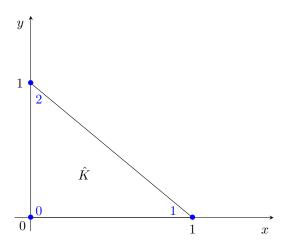
for  $i, j = 0, \dots, N - 1$ .

1c)

Complete the template file shape.hpp implementing the function

which computes the the value a local shape function  $\lambda_i(\mathbf{x})$ , with i that can assume the values 0, 1 or 2, on the reference element depicted in Fig. 1 at the point  $\mathbf{x} = (x, y)$ .

The convention for the local numbering of the shape functions is that  $\lambda_i(\mathbf{x}_j) = \delta_{i,j}$ , i, j = 0, 1, 2, with  $\delta_{i,j}$  denoting the Kronecker delta.



**Figure 1:** Reference element  $\hat{K}$  for 2D linear finite elements.

**Solution:** See listing 1 for the code.

Listing 1: Implementation for lambda

#### #pragma once

```
//! The shape function (on the reference element)
//!
//! We have three shape functions.
//!
//! lambda(0, x, y) should be 1 in the point (0,0) and zero in (1,0) and (0,1)
//! lambda(1, x, y) should be 1 in the point (1,0) and zero in (0,0) and (0,1)
//! lambda(2, x, y) should be 1 in the point (0,1) and zero in (0,0) and (1,0)
//!
//! @param i integer between 0 and 2 (inclusive). Decides which shape function to
   \hookrightarrow return.
//! @param x x coordinate in the reference element.
//! Oparam y y coordinate in the reference element.
inline double lambda(int i, double x, double y) {
 //// NPDE_START_TEMPLATE
   if (i == 0) {
       return 1 - x - y;
   } else if (i == 1) {
       return x;
   } else {
```

```
return y;
}
//// NPDE_RETURN_TEMPLATE
//// NPDE_END_TEMPLATE
}
```

#### 1d)

Complete the template file grad\_shape.hpp implementing the function

```
inline Eigen::Vector2d gradientLambda(const int i, double x, double y)
```

which returns the value of the derivatives (i.e. the gradient) of a local shape functions  $\lambda_i(\boldsymbol{x})$ , with i that can assume the values 0,1 or 2, on the reference element depicted in Fig. 1 at the point  $\boldsymbol{x}=(x,y)$ .

**Solution:** See listing 2 for the code.

Listing 2: Implementation for gradientLambda

```
#pragma once
#include <Eigen/Core>
//! The gradient of the shape function (on the reference element)
//! We have three shape functions
//!
//! @param i integer between 0 and 2 (inclusive). Decides which shape function to
   \hookrightarrow return.
//! Oparam x x coordinate in the reference element.
//! @param y y coordinate in the reference element.
inline Eigen::Vector2d gradientLambda(const int i, double x, double y) {
 //// NPDE_START_TEMPLATE
   return Eigen::Vector2d(-1 + (i > 0) + (i==1),
                         -1 + (i > 0) + (i==2));
   //// NPDE_END_TEMPLATE
   return Eigen::Vector2d(0,0); //remove when implemented
}
```

The routine makeCoordinateTransform contained in the file coordinate\_transform.hpp computes the Jacobian matrix of the linear map  $\Phi_l : \mathbb{R}^2 \to \mathbb{R}^2$  such that

$$\Phi_l \begin{pmatrix} 1 \\ 0 \end{pmatrix} = \begin{pmatrix} a_{11} \\ a_{12} \end{pmatrix} = \boldsymbol{a}_1, \quad \Phi_l \begin{pmatrix} 0 \\ 1 \end{pmatrix} = \begin{pmatrix} a_{21} \\ a_{22} \end{pmatrix} = \boldsymbol{a}_2,$$

where  $a_1, a_2 \in \mathbb{R}^2$  are the two input arguments.

### 1e)

Complete the template file stiffness\_matrix.hpp implementing the routine

that returns the element stiffness matrix for the bilinear form associated to (1) and for the triangle with vertices a, b and c.

Hint: Use the routine gradientLambda from subproblem 1d) to compute the gradients and the routine makeCoordinateTransform to transform the gradients and to obtain the area of a triangle.

**Hint:** You do not have to analytically compute the integrals for the product of basis functions; instead, you can use the provided function integrate. It takes a function f(x, y) as a parameter, and it returns the value of  $\int_K f(x, y) dV$ , where K is the triangle with vertices in (0, 0), (1, 0) and (0, 1). Do not forget to take into account the proper coordinate transforms!

**Solution:** See listing 3 for the code.

**Listing 3:** Implementation for computeStiffnessMatrix

```
//! Evaluate the stiffness matrix on the triangle spanned by
//! the points (a, b, c).
//!
//! Here, the stiffness matrix A is a 3x3 matrix
//! A_{ij} = \int_{K} ( \lambda_i^K(x, y) \cdot \lambda_j^K(x, y) 
   \hookrightarrow \; dV$$
//!
//! where $K$ is the triangle spanned by (a, b, c).
//! @param[out] stiffnessMatrix should be a 3x3 matrix
//! At
                           the end, will contain the integrals above.
//!
//! @param[in] a the first corner of the triangle
//! @param[in] b the second corner of the triangle
//! @param[in] c the third corner of the triangle
template<class MatrixType, class Point>
void computeStiffnessMatrix(MatrixType& stiffnessMatrix,
                         const Point& a,
                         const Point& b,
                         const Point& c)
{
```

```
Eigen::Matrix2d coordinateTransform = makeCoordinateTransform(b - a, c - a);
   double volumeFactor = std::abs(coordinateTransform.determinant());
   Eigen::Matrix2d elementMap = coordinateTransform.inverse().transpose();
   //// NPDE_START_TEMPLATE
   for (int i = 0; i < 3; ++i) {
       for (int j = i; j < 3; ++j) {
           stiffnessMatrix(i, j) = integrate([&](double x, double y) {
              Eigen::Vector2d gradLambdaI = elementMap * gradientLambda(i, x, y);
              Eigen::Vector2d gradLambdaJ = elementMap * gradientLambda(j, x, y);
              auto lambdaI = lambda(i, x, y);
              auto lambdaJ = lambda(j, x, y);
              return volumeFactor * gradLambdaI.dot(gradLambdaJ);
          });
       }
   }
   // Make symmetric (we did not need to compute these value above)
   for (int i = 0; i < 3; ++i) {
       for (int j = 0; j < i; ++j) {
          stiffnessMatrix(i, j) = stiffnessMatrix(j, i);
   }
   //// NPDE_END_TEMPLATE
}
```

The routine integrate in the file integrate.hpp uses a quadrature rule to compute the approximate value of  $\int_{\hat{K}} f(\hat{x}) d\hat{x}$ , where f is a function, passed as input argument.

#### 1f)

Complete the template file load\_vector.hpp implementing the routine

that returns the element load vector for the linear form associated to (1), for the triangle with vertices a, b and c, and where f is a function handler to the right-hand side of (1).

Hint: Use the routine lambda from subproblem 1c) to compute values of the shape functions on the reference element, and the routines makeCoordinateTransform and integrate from the handout to map the points to the physical triangle and to compute the integrals.

**Solution:** See listing 4 for the code.

**Listing 4:** Implementation for computeLoadVector

```
//! Evaluate the load vector on the triangle spanned by
//! the points (a, b, c).
//! Here, the load vector is a vector $(v_i)$ of
//! three components, where
//! $v_i = \int_{K} \lambda_i^K(x, y) f(x, y) \
//!
//! where $K$ is the triangle spanned by (a, b, c).
//! @param[out] loadVector should be a vector of length 3.
//! At
                           the end, will contain the integrals above.
//!
//! @param[in] a the first corner of the triangle
//! @param[in] b the second corner of the triangle
//! @param[in] c the third corner of the triangle
//! @param[in] f the function f (LHS).
template < class Vector, class Point >
void computeLoadVector(Vector& loadVector,
                  const Point& a, const Point& b, const Point& c,
                  const std::function<double(double, double)>& f)
{
   Eigen::Matrix2d coordinateTransform = makeCoordinateTransform(b - a, c - a);
   double volumeFactor = std::abs(coordinateTransform.determinant());
   //// NPDE_START_TEMPLATE
   for (int i = 0; i < 3; ++i) {</pre>
       loadVector(i) = integrate([&](double x, double y) {
           Eigen::Vector2d z = coordinateTransform * Eigen::Vector2d(x, y) + Eigen
               \hookrightarrow :: Vector2d(a(0), a(1));
           return f(z(0), z(1)) * lambda(i, x, y) * volumeFactor;
       });
   }
   //// NPDE_END_TEMPLATE
}
```

#### 1g)

Complete the template file stiffness\_matrix\_assembly.hpp implementing the routine

to compute the finite element matrix  $\mathbf{A}$  as in (4). The input argument vertices is a  $N_V \times 3$  matrix of which the *i*-th row contains the coordinates of the *i*-th mesh vertex,  $i=0,\ldots,N_V-1$ , with  $N_V$  the number of vertices. The input argument triangles is a  $N_T \times 3$  matrix where the *i*-th row contains the *indices* of the vertices of the *i*-th triangle,  $i=0,\ldots,N_T-1$ , with  $N_T$  the number of triangles in the mesh.

**Hint:** Use the routine computeStiffnessMatrix from subproblem 1e) to compute the local stiffness matrix associated to each element.

**Hint:** Use the sparse format to store the matrix A.

**Solution:** See listing 5 for the code.

**Listing 5:** Implementation for assembleStiffnessMatrix

```
//! Assemble the stiffness matrix
//! for the linear system
//! @param[out] A will at the end contain the Galerkin matrix
//! @param[in] vertices a list of triangle vertices
//! @param[in] triangles a list of triangles
template < class Matrix>
void assembleStiffnessMatrix(Matrix& A, const Eigen::MatrixXd& vertices,
                         const Eigen::MatrixXi& triangles)
{
   const int numberOfElements = triangles.rows();
   A.resize(vertices.rows(), vertices.rows());
   std::vector<Triplet> triplets;
   triplets.reserve(numberOfElements * 3 * 3);
   //// NPDE_START_TEMPLATE
   for (int i = 0; i < numberOfElements; ++i) {</pre>
       auto& indexSet = triangles.row(i);
       const auto& a = vertices.row(indexSet(0));
       const auto& b = vertices.row(indexSet(1));
       const auto& c = vertices.row(indexSet(2));
```

#### 1h)

Complete the template file load\_vector\_assembly.hpp implementing the routine

to compute the right-hand side vector  $\mathbf{F}$  as in (4). The input arguments vertices and triangles are as in subproblem  $\mathbf{1g}$ ), and f is an in subproblem  $\mathbf{1f}$ ).

**Hint:** Proceed in a similar way as for assembleStiffnessMatrix and use the routine computeLoadVector from subproblem 1f).

**Solution:** See listing 6 for the code.

Listing 6: Implementation for assembleLoadVector

```
F.resize(vertices.rows());
    F.setZero();
    //// NPDE_START_TEMPLATE
    for (int i = 0; i < numberOfElements; ++i) {</pre>
        const auto& indexSet = triangles.row(i);
        const auto& a = vertices.row(indexSet(0));
        const auto& b = vertices.row(indexSet(1));
        const auto& c = vertices.row(indexSet(2));
        Eigen::Vector3d elementVector;
        computeLoadVector(elementVector, a, b, c, f);
        for (int i = 0; i < 3; ++i) {</pre>
           F(indexSet(i)) += elementVector(i);
    }
     //// NPDE_END_TEMPLATE
}
The routine
void setDirichletBoundary(Eigen::VectorXd& u, Eigen::VectorXi& interiorVertexIndices,
                         const Eigen::MatrixXd& vertices,
                         const Eigen::MatrixXi& triangles,
                         const std::function<double(double, double)>& g)
```

implemented in the file dirichlet\_boundary.hpp provided in the handout does the following:

- it gets in input the matrices vertices and triangles as defined in subproblem 1g) and the function handle g to the boundary data, i.e. to q such that u = q on  $\partial\Omega$  (in our case  $q \equiv 0$ );
- it returns in the vector interior VertexIndices the indices of the interior vertices, that is of the vertices that are not on the boundary  $\partial\Omega$ ;
- if  $x_i$  is a vertex on the boundary, then it sets  $u(i)=g(x_i)$ , that is, in our case, it sets to 0 the entries of the vector u corresponding to vertices on the boundary.

#### 1i)

Complete the template file fem\_solve.hpp with the implementation of the function

This function takes in input the matrices vertices, triangles as defined in the previous subproblems, and the function handle f to the right-hand side f in (1). The output argument u has to contain, at the end of the function, the finite element solution  $u_N$  to (1).

**Hint:** Use the routines assembleStiffnessMatrix and assembleLoadVector from subproblems 1g) and 1h), respectively, to obtain the matrix A and the vector F as in (4), and then use the provided routine setDirichletBoundary to set the boundary values of u to zero and to select the free degrees of freedom.

**Solution:** See listing 7 for the code.

Listing 7: Implementation for solveFiniteElement

```
//! Solve the FEM system.
//!
//! Cparam[out] u will at the end contain the FEM solution.
//! @param[in] vertices list of triangle vertices for the mesh
//! @param[in] triangles list of triangles (described by indices)
//! @param[in] f the RHS f (as in the exercise)
//! return number of degrees of freedom (without the boundary dofs)
int solveFiniteElement(Vector& u,
   const Eigen::MatrixXd& vertices,
   const Eigen::MatrixXi& triangles,
   const std::function<double(double, double)>& f)
{
   SparseMatrix A;
   //// NPDE_START_TEMPLATE
   assembleStiffnessMatrix(A, vertices, triangles);
   //// NPDE_END_TEMPLATE
   Vector F;
   //// NPDE_START_TEMPLATE
   assembleLoadVector(F, vertices, triangles, f);
   //// NPDE_END_TEMPLATE
   u.resize(vertices.rows());
   u.setZero();
   Eigen::VectorXi interiorVertexIndices;
   auto zerobc = [](double x, double y){ return 0;};
   // set homogeneous Dirichlet Boundary conditions
   //// NPDE_START_TEMPLATE
   setDirichletBoundary(u, interiorVertexIndices, vertices, triangles, zerobc);
   F -= A * u;
   //// NPDE_END_TEMPLATE
   SparseMatrix AInterior;
```

```
igl::slice(A, interiorVertexIndices, interiorVertexIndices, AInterior);
   Eigen::SimplicialLDLT<SparseMatrix> solver;
   Vector FInterior;
   igl::slice(F, interiorVertexIndices, FInterior);
   //initialize solver for AInterior
   //// NPDE_START_TEMPLATE
   solver.compute(AInterior);
   if (solver.info() != Eigen::Success) {
       throw std::runtime_error("Could not decompose the matrix");
   //// NPDE_END_TEMPLATE
   //solve interior system
   //// NPDE_START_TEMPLATE
   Vector uInterior = solver.solve(FInterior);
   igl::slice_into(uInterior, interiorVertexIndices, u);
   //// NPDE_END_TEMPLATE
   return interiorVertexIndices.size();
}
```

## 1j)

Run the code in the file fem2d.cpp to compute the finite element solution to (1) when  $\Omega = [0,1]^2$  is the unit square, the forcing term is given by  $f(x) = 2\pi^2 \sin(\pi x) \sin(\pi y)$  and the mesh is square\_5.  $\rightarrow$  mesh. Use then the routine plot\_on\_mesh.py to produce a plot of the solution.

**Solution:** See Fig. 2 for the plot.

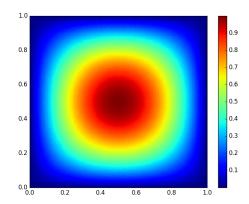


Figure 2: Solution plot for subproblem 1j).