

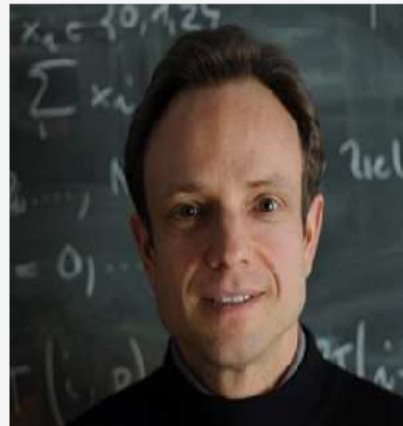
Homework Problem Video Tutorial

Problem 2-5: Triangular linear FEM for 2D reaction-diffusion BVP

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Essential prerequisite: Sect 2.4.



$$u \in H^1(\Omega): \underbrace{\int_{\Omega} \text{grad } u \cdot \text{grad } v + uv \, dx}_{:=a(u,v)} = \underbrace{\int_{\Omega} f v \, dx}_{:=\ell(v)} \quad \forall v \in H^1(\Omega). \quad (2.5.1)$$

"reaction term"

a) Strong form of (2.5.1):
l.b.p. (Green's formula)

$$\int_{\Omega} \underbrace{(-\text{div grad } u + u)}_{=0 \text{ by testing with } v \in C_0^\infty(\Omega)} v \, dx + \int_{\partial\Omega} \underbrace{\text{grad } u \cdot n}_{=0 \text{ by testing with } v \in C^\infty(\bar{\Omega})} v \, dS = \int_{\Omega} f v \, dx \quad \forall v \in H^1(\Omega)$$

$$\begin{aligned} \triangleright \quad & -\Delta u + u = f \quad \text{in } \Omega \\ & \text{grad } u \cdot n = 0 \quad \text{on } \partial\Omega \end{aligned}$$

b) Element matrix M_K for $(u, v) \rightarrow \int_{\Omega} uv \, dx$ for general triangle K .

[Setting: FE space $S_1^0(\mathcal{M})$]

$$(M_K)_{i,j} = \int_K \underbrace{\lambda_i \lambda_j}_{\text{barycentric coordinate fcts.}} \, dx$$

Lemma 2.7.5.5. Integration of powers of barycentric coordinate functions

For any non-degenerate d -simplex K with barycentric coordinate functions $\lambda_1, \dots, \lambda_{d+1}$ and exponents $\alpha_j \in \mathbb{N}, j = 1, \dots, d+1$,

$$\int_K \lambda_1^{\alpha_1} \dots \lambda_{d+1}^{\alpha_{d+1}} dx = d!|K| \frac{\alpha_1! \alpha_2! \dots \alpha_{d+1}!}{(\alpha_1 + \alpha_2 + \dots + \alpha_{d+1} + d)!} \quad \forall \alpha \in \mathbb{N}_0^{d+1}. \quad (2.7.5.6)$$

$d = 2$: $(M_K)_{ii} \stackrel{!}{=} \alpha_i = 2, \alpha_2 = \alpha_3 = 0$
 $\triangleright = \frac{1}{6} |K|$
 $(M_K)_{ij} \stackrel{!}{=} \alpha_i = \alpha_j = 1, \alpha_3 = 0 \quad (i \neq j)$
 $\triangleright = \frac{1}{12} |K|$

$$M_K = \frac{1}{12} |K| \begin{bmatrix} 2 & 1 & 1 \\ 1 & 2 & 1 \\ 1 & 1 & 2 \end{bmatrix}$$

$$\mathbb{1}^T M_K \mathbb{1} = \int_K 1 \cdot dx = |K|$$

c_1 Computation of $M_K \in \mathbb{R}^{3,3}$

d_1

Implement the function

```
double L2Error(const TriangularMesh2D& mesh,
               const Eigen::VectorXd& uFEM,
               const std::function<double(double, double)> exact);
```

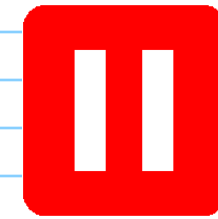
to compute the error $\|u_h - u\|_{L^2(\Omega)}$, where u is the exact solution to (2.5.1), passed through the function handle `exact`, and $u_h \in \mathcal{S}_1^0(\mathcal{M})$ is the finite element Galerkin solution passed through its coefficient vector `uFEM` with respect to the customary basis of tent functions. The argument `mesh` contains the mesh data structure.

A function $\in H^1(\Omega) \subset L^2(\Omega)$

Take into account : u in procedural form

$\triangleright$ Numerical quadrature mandatory

$\triangleright$ Cell-oriented computation



$$\|u_h - u\|_{L^2(\Omega)}^2 \approx \sum_{K \in \mathcal{M}} \frac{1}{3} |K| \left((u_h - u)^2(a_K^1) + (u_h - u)^2(a_K^2) + (u_h - u)^2(a_K^3) \right),$$

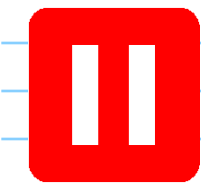
$u_h(a_K^i) \stackrel{!}{=} \text{entries of } uFEM$

e_1 Implement the function

```
double H1Error(const TriangularMesh2D& mesh, const Eigen::VectorXd&
               uFEM,
               const std::function<Eigen::Vector2d(double, double)>
               exact);
```

to compute the error norm $\|u_h - u\|_{H^1(\Omega)}$, where u is the exact solution of (2.5.1), whose gradient is passed in the function handle `exact` (that returns a column vector), and $u_h \in \mathcal{S}_1^0(\mathcal{M})$ is the finite element Galerkin solution, passed through the coefficient vector `u`. The input argument `Mesh` contains the mesh data structure.

③ Again: Numerical quadrature by 2D local trapezoidal rule



$$|u_h - u|_{H^1(\Omega)}^2 \approx \sum_{K \in \mathcal{M}} \frac{1}{3} |K| \left(\left\| \text{grad } u_h(a_K^1) - \text{grad } u(a_K^1) \right\|^2 + \left\| \text{grad } u_h(a_K^2) - \text{grad } u(a_K^2) \right\|^2 + \left\| \text{grad } u_h(a_K^3) - \text{grad } u(a_K^3) \right\|^2 \right).$$

↑
the same vector $\forall a_K^e$

Observe: $\text{grad } u_h$ is $\mathcal{M}$ -p.w. constant

$$\text{grad } u_h|_K = \sum_{e=1}^3 u_h(a_K^e) \text{grad } \lambda_e$$

↑
entry of VFEM

C++ code 2.4.5.11: Computation of gradients of barycentric coordinate functions on a triangle → GITLAB

```
1 Eigen::Matrix<double, 2, 3> gradbarycoordinates(const t_TriGeo& Vertices)
2 {
3   // Argument Vertices passes the vertex positions of the triangle
4   // as the rows of a 3x2-matrix, see Code 2.4.1.3...
5   // The function returns the components of the gradients as the
6   // columns of a 2x3-matrix
7
8   // Computation based on (2.4.5.10), solving for the
9   // coefficients of the barycentric coordinate functions.
10  Eigen::Matrix<double, 3, 3> X; // Temporary matrix
11  X.block<3, 1>(0, 0) = Eigen::Vector3d::Ones();
12  X.block<3, 2>(0, 1) = Vertices.transpose();
13  return X.inverse().block<2, 3>(1, 0);
14 }
```

f)

Implement a function

```
std::tuple<Eigen::VectorXd, double, double>
solve(const TriangularMesh2D& mesh);
```

that, given in input a mesh data structure `mesh`, computes the discrete solution $u_h \in \mathcal{S}_1^0(\mathcal{M})$ to (2.5.1) in the case that the exact solution is $u(x) = \cos(2\pi x_1) \cos(2\pi x_2)$, plots the mesh and u_h . The function returns the coefficient vector `U` of u_h , the L^2 -norm and the (approximate, due to quadrature) H^1 -seminorm of the discretization error $u - u_h$.

You can rely on the functions `assemLoad_LFE()` and `GalerkinAssembly()` given in the file `simple_linear_finite_elements.cc`. Those implement cell-oriented assembly according to the "distribute scheme", see [Lecture → Code 2.4.6.8] and [Lecture → Code 2.4.5.23].

How to make a given u a solution of the BVP

- Use strong form $\rightarrow f := -\Delta u + u$
- Verify b.c.

"method of manufactured solutions"







6

