Project in TMA4220

Candidatenumber: 10000 & 10028

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0.1 Poisson solver

We here present a solver for the Poisson equation,

$$\nabla^2 u = f,$$

in both 2 and 3 dimensions, with both Neumann and Dirichlet boundary conditions, while using the element method. We let $u(r) = \sin(2\pi r^2)$ and $f(r) = \nabla^2 u(r) = -8\pi \cos(2\pi r^2) + 16r^2\pi^2 \sin(2\pi r^2)$.

0.1.1 2D solver

The program poisson2d.m solves the Poisson equation in 2 dimensions, with Dirichlet boundary conditions, and created the plots in figure 1.

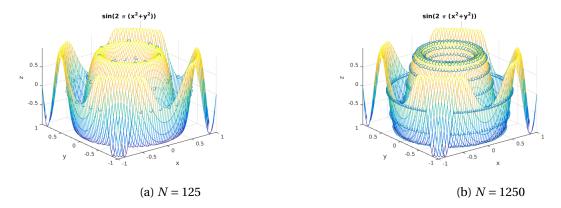


Figure 1: Scatterplotts of the numerical solution of the Poisson equation in 2 dimensions, with Dirichlet boundary condition and different number of points, *N*, together with a surfplot of the correct solution.

The program poissonBnd2d.m solves the Poisson equation in 2 dimensions, with Dirichlet boundary conditions on half the domain, and Neumann on the other half, and created the plots in figure 2.

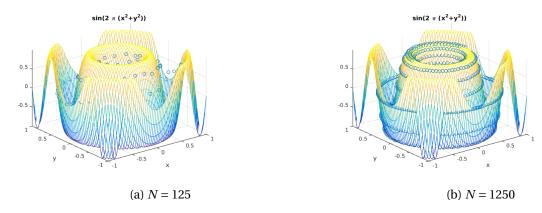


Figure 2: Scatterplotts of the numerical solution of the Poisson equation in 2 dimensions, with Dirichlet boundary condition on half the boundary, and Neumann on the other half, with and different number of points, *N*, together with a surfplot of the correct solution.



Figure 3: A plot of the numerical solution of the Poisson equation in 3 dimensions, with Dirichlet boundary condition, with different number of points, *N*. The correct solution is a perfect sphere.

0.1.2 3D solver

The program poisson3d.m solves the Poisson equation in 3 dimensions, with Dirichlet boundary conditions, and created the plots in figure 3.

The program poissonBnd3d.m solves the Poisson equation in 3 dimensions, with Dirichlet boundary condition on half the boundary, and Neumann on the other half, and created the plots in figure 4

0.1.3 Discussion

As we can see from figure 1 and figure 2, the 2 dimensional case works quite good, and seems to be convergent. Thus poisson2d.m and poissonBnd2d.m are working solvers for the Poisson equation.

Also the 3 dimensional case with Dirichelt boundary the solver poisson3d.m converges. But when including the Neumann boundary condition the *A* matrix became singular because we needed to include

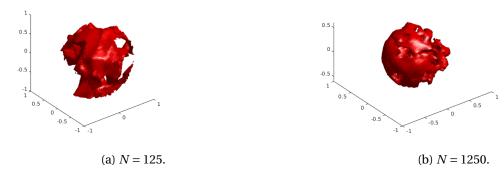


Figure 4: A plot of the numerical solution of the Poisson equation in 3 dimensions, with Dirichlet boundary condition on half the boundary, and Neumann on the other half, with different number of points, *N*. The correct solution is a perfect sphere.

some points on the boundary which we did not need while only having Dirichelt boundary. There also seems to be a problem with while adding the Neumann conditions to b. Thus we conclude that the program poisson3dBnd.m does not work correctly. We have no clue why the Neumann conditions works in 2 dimension and not in 3 dimensions, considering that the implementation is almost identical.

0.2 Meat equation

We are here going to present a solution to the heat equation in 3 dimensions using Neumann conditions on the edges, with forward Euler in time, and element-method in space.

0.2.1 Weak formulation

The heat equation

$$\frac{\partial u}{\partial t} = \alpha \nabla^2 u \tag{1}$$

Using Greens formula we obtain

$$\int_{\Omega} \frac{\partial u(\mathbf{x})}{\partial t} v(\mathbf{x}) d\Omega = \int_{\Omega} \Delta u(\mathbf{x}) v(\mathbf{x}) d\Omega = -\int_{\Omega} \nabla u(\mathbf{x}) \nabla v(\mathbf{x}) d\Omega + \int_{\Gamma} \frac{\partial u}{\partial n} (\mathbf{x}) v(\mathbf{x}) d\gamma.$$

This has an unique solution if continuous and weakly coercive. $f \equiv 0$ for our equation.

The scheme is weakly coersive if $a(u,u) + \lambda \|u\|_{L^2(\Omega)} \ge \alpha \|u\|_V$ for $\lambda \ge 0$, $\alpha > 0$, with $a(u,v) = \int \nabla u(\mathbf{x}) \nabla v(\mathbf{x}) d\Omega$. $|a(u,v)| = |\int \nabla u \nabla v d\Omega| \le \sqrt{\int (\nabla u)^2 d\Omega} \sqrt{\int (\nabla v)^2 d\Omega} = \|u\|_V \|v\|_V$ so the the scheme is continuous. $|a(v,v)| = \int \nabla v \nabla v d\Omega = \|v\|_V^2$, so this scheme is weakly coercive for $\alpha \le 1$. Therefore there exists a unique solution.

Trying to solve our problem using v and u in the space $H^1(\Omega)$, approximated by $X_h^1 = \{v_h \in C^0(\bar{\Omega}) : v_h|_K \in \mathbb{P}_1 \forall K \in \tau_h\}$ in both 2 and 3 dimentions. $V_h = \{v_h \in X_h^1 : v \text{ satisfying boundary conditions}\}$. Using Lagrangian functions $\phi_j \in V_h$ with the property

$$\phi_j(x_i) = \delta_{ij} = \begin{cases} 1 & i = j \\ 0 & i \neq j \end{cases}$$

0.2.2 Galerkin

We can make a Galerkin problem using

$$u_h(\mathbf{x},t) = \sum_{j=1}^{N_h} u_j(t)\phi(\mathbf{x}) \text{ and } v \in span\{\phi_1,\phi_2,\cdots,\phi_{N_h}\}.$$

This gives

$$\int \sum_{j=1}^{N_h} \frac{\partial u_j}{\partial t} \phi_j \phi_i d\Omega + a(\sum_{j=1}^{N_h} u_j(t) \phi_j, \phi_i) = \int_{\Gamma} \sum_{j=1}^{N_h} \frac{\partial u_j(t) \phi_j}{\partial n} \phi_i d\gamma$$

$$\sum_{i=1}^{N_h} \frac{\partial u_j}{\partial t} \int \phi_j \phi_i d\Omega + \sum_{i=1}^{N_h} u_j(t) a(\phi_j, \phi_i) = \sum_{i=1}^{N_h} \frac{\partial u_j}{\partial t} m_{ij} + \sum_{i=1}^{N_h} u_j(t) a_{ij} = b_i$$

This is the linear system, $M\dot{\mathbf{u}}(t) + A\mathbf{u}(t) = \mathbf{0}$, the boundary conditions get inserted differently.

The Lagrangian functions ϕ is on the form $\phi_i = \alpha_i x \beta_i y + \gamma_i$. For K_j ϕ_i is 1 in $(x, y) = (x_i, y_i)$ and 0 in the other corners in K. The K elements are triangles in 2 dimentions and tetrahedrons in 3 dimentions, and together they cover our domain Ω .

The Galerkin method let's us finally construct forward Euler method iterations on the form $M \frac{\mathbf{u}^{k+1} - \mathbf{u}^k}{\Delta t} + A\mathbf{u}^k = 0$.

0.2.3 Stability of forward euler galerkin

The stability analysis for forward euler in time and FEM in space we find by using w_h^j which is eigenvectors of $a(\cdot, \cdot)$ instead of ϕ , the step-size h may vary in space.

$$\frac{1}{\Delta t} \sum_{j=1}^{N_h} [u_j^{k+1} - u_j^k] \int w_h^j w_h^i + \sum_{j=1}^{N_h} u_j^k a(w_h^j, w_h^i) = 0$$

The eigenvalues of the matrix can be calculated

$$a(w_h^j, w_h^i) = \int \nabla w_h^j \nabla w_h^i d\Omega = \lambda_h^j \int w_h^j w_h^i d\Omega = \lambda_h^j (w_h^j, w_h^i) = \lambda_h^j \delta_{ij} = \lambda_h^i$$

giving

$$\frac{u_j^{k+1} - u_j^k}{\Delta t} + u_j^k \lambda_h^i = 0 \Rightarrow u_j^{k+1} = u_j^k (1 - \lambda_h^j \Delta t).$$

For the method to be absolutely stable we need $|1 - \lambda_h^j \Delta t| < 1$.

 $-1 < 1 - \lambda_h^j \Delta t < 1 \Rightarrow 0 < \Delta t < \frac{2}{\lambda_h^i}$. The eigenvalues for the stiffness matrix are

$$\max_{i} \lambda_{h}^{i} = \max_{i} \frac{a(w_{h}^{i}, w_{h}^{i})}{\|w_{h}^{i}\|_{L^{2}(\Omega)}^{2}} \leq \max_{i} \frac{\|w_{h}^{i}\|_{V}^{2}}{\|w_{h}^{i}\|_{L^{2}(\Omega)}^{2}} \simeq (1 + \frac{c}{h^{2}}), \text{ since } \|\nabla u\|_{L^{2}(\Omega)} \leq \frac{C\|u\|_{L^{2}(\Omega)}}{h}.$$

We get that this method only is stable if $\Delta t \leq Ch^2$.

0.2.4 Initial and boundary conditions

We assume that the body initially has a uniform temperature, T_i .

On the boundary we have Neumann-conditions. We divide the boundary in two parts, $\partial \Omega_H$, $\partial \Omega_C$. $\partial \Omega_H$ is the part of the boundary that is in contact with temperature T_H , while $\partial \Omega_C$ is in contact with temperature T_C .

0.2.5 Numerical scheme

The resulting equations that needs to be solved is

$$\begin{split} M\frac{du}{dt} &= -\alpha Au & \text{on } \Omega \\ \frac{\partial u}{\partial v} &= \alpha_H (T_H - u) & \text{on } \partial \Omega_H \\ \frac{\partial u}{\partial v} &= \alpha_C (T_C - u) & \text{on } \partial \Omega_C. \end{split}$$

This is a simple set of ODEs, and is solved using forward Euler. The numerical scheme can then be written as

$$u^{i+1} = u^i - \alpha(M \setminus A)u^i$$
 on Ω
 $u^{i+1} = u^i + \alpha_H(T_H - u^i)$ on $\partial \Omega_H$.
 $u^{i+1} = u^i + \alpha_C(T_C - u^i)$ on $\partial \Omega_C$

Where α , α_H and α_C are the heat conductivity in the body, heat conductivity between $\partial\Omega_H$ and T_H , the heat conductivity between $\partial\Omega_C$ and T_C , respectivly.

0.2.6 Correctness of Method

To test if A, b and M where correct, we wanted to find

$$E_A = ||Au - b_1||_2,$$
 $b_1 = \int_{\Omega} f v d\Omega$
 $E_M = ||Mu - b_2||_2,$ $b_2 = \int_{\Omega} u v d\Omega$

with $u(r) = \sin(2\pi r^2)$, $f(r) = \nabla^2 u(r) = -8\pi\cos(2\pi r^2) + 16r^2\pi^2\sin(2\pi r^2)$ on the unit sphere. The program error3d.m was made for this purpose, and produced tabel 1, and figure 5. From table 1 it seems that E_A does not converge, but some experimentation with error3d.m shows that it decreases extremly slowly with increasing N. For the matrix M this is not an issue and it converges nicely. The 2 dimensional case will not be discussed in this report, but the program error2d.m can be used if this is of interest.

0.2.7 Example of use

As an example of use we wanted to cook a beef. For that we need a lot of different assumptions.

- The beef has a initial uniform temperature $T_i = 10^{\circ}C$, and is turned when the temperature on the middle poit reaches a temperature $T_f = 60^{\circ}C$. It is done when the lowest temperature in the beef is $T_f = 60^{\circ}C$. $T_H = 180^{\circ}C$, $T_C = 20^{\circ}C$.
- The beef is discretized with $n_1 = n_2 = n_3 = 5$ with the function getbeef.m.
- Thermal conductivities are independent of temperature, and as follows: $\alpha = 8.75 \cdot 10^{-8}$, $\alpha_H = 0.7$, $\alpha_C = 3 \cdot 10^{-4}$.

$$N = 125$$
 $N = 250$ $N = 1250$
 E_A 19.50 22.09 12.82
 E_M 0.1165 0.0556 0.0121

Table 1: Error of A and M, when compared to their corresponding b_1 and b_2 .

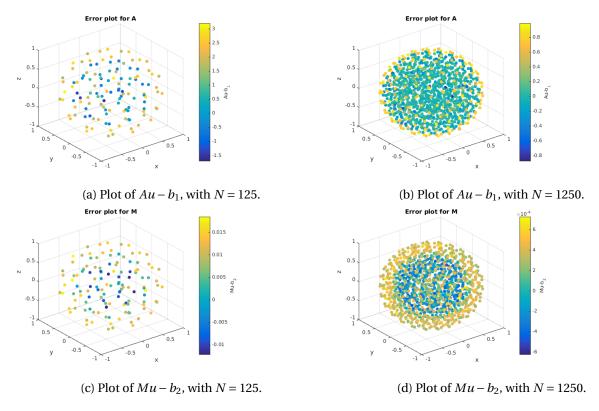


Figure 5: A scatterplott showing the pointwise error of *A* and *M*, with different number of points, *N*.

 α is given in the task, α_H is assumed to be quite large, while α_C is assumed to be somwhere in between α and α_H . The other constants are chosen freely.

Result

Running the program beef .m with these assumptions, gives out the following times:

- Time to turn the beef: 658 seconds
- Total time to the beef is done: 776 seconds

This is close to 13 minutes in total cookingtime, which seems about right for a normal steak. As we can see from figure 6, the edges are all cooked, while the upper middle is some what raw.

Discussion

Although the program beef.m gives a time-estimate similar to the reality, it is in no way an accurate description. The numerical scheme has it problems, the A matrix converges extremely slow with N. The program gives different cooking time with different values of n_i .

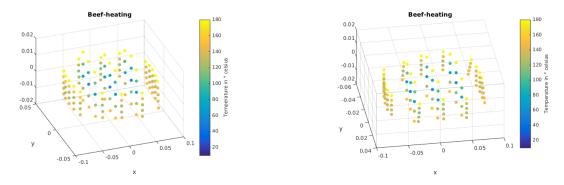


Figure 6: A scatterplott of a cooked beef from two different angles.