

# A simple FEM solver and its data parallelism

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# Partial differential equation

## Considered Problem Classes

Find  $u$  such that

$$\begin{aligned} Lu(x) &= f(x) & \forall x \in \Omega \\ lu(x) &= g(x) & \forall x \in \partial\Omega \end{aligned}$$

variational  $\Downarrow$  formulation

Find  $u \in \mathbb{V}$ :

$$a(u, v) = \langle F, v \rangle \quad \forall v \in \mathbb{V}$$

FEM, FDM  $\Downarrow$  FVM, FIT

Solve

$$K_h \cdot \underline{u}_h = \underline{f}_h \quad \underline{u}_h \in \mathbb{R}^{N_h}$$

► (linear) 2<sup>nd</sup> order problem.

- Poisson equation  
(temperature)
- Lamé equation  
(deformation)
- Maxwell's equations  
(magnetic field)

► Matrix  $K_h$  is sparse, positive definite

(symmetric, large dimension)

► non-linear and time-dependent problems.

## Second order PDE

Find  $u \in X := C^2(\Omega) \cap C^1(\Omega \cup \Gamma_2 \cup \Gamma_3) \cap C(\Omega \cup \Gamma_1)$  such that the partial differential equation

$$-\sum_{i,j=1}^m \frac{\partial}{\partial x_i} \left( a_{ij}(x) \frac{\partial u}{\partial x_j} \right) + \sum_{i=1}^m a_i(x) \frac{\partial u}{\partial x_i} + a(x)u(x) = f(x) \quad (1)$$

holds for all  $x \in \Omega$  and that the Boundary Conditions (BC)

- ▶  $u(x) = g_1(x), \forall x \in \Gamma_1$  (Dirichlet (1<sup>st</sup>-kind) BC),
- ▶  $\frac{\partial u}{\partial N} := \sum_{i,j=1}^m a_{ij}(x) \frac{\partial u(x)}{\partial x_j} n_i(x) = g_2(x), \forall x \in \Gamma_2$   
(Neumann (2<sup>nd</sup>-kind) BC),
- ▶  $\frac{\partial u}{\partial N} + \alpha(x)u(x) = g_3(x), \forall x \in \Gamma_3$  (Robin (3<sup>rd</sup>-kind) BC).

are satisfied.

with  $u(x)$  as classical continuous solution of the PDE.

## Variational formulation

Choose the space of test functions  $V_0 = \{v \in V = H^1(\Omega) : v = 0 \text{ on } \Gamma_1\}$ , where  $V = H^1(\Omega)$  is the basic space

Find  $u \in V_g$  such that  $a(u, v) = \langle F, v \rangle \quad \forall v \in V_0$ , where

$$a(u, v) := \int_{\Omega} \left( \sum_{i,j=1}^m a_{ij} \frac{\partial u}{\partial x_j} \frac{\partial v}{\partial x_i} + \sum_{i=1}^m a_i \frac{\partial u}{\partial x_i} v + a u v \right) dx + \int_{\Gamma_3} \alpha u v ds,$$

$$\langle F, v \rangle := \int_{\Omega} f v dx + \int_{\Gamma_2} g_2 v ds + \int_{\Gamma_3} g_3 v ds,$$

$$V_g := \{v \in V = H^1(\Omega) : v = g_1 \text{ on } \Gamma_1\},$$

$$V_0 := \{v \in V : v = 0 \text{ on } \Gamma_1\}.$$

(2)

with  $u(x)$  as weak continuous solution of the PDE.

## Finite Elements

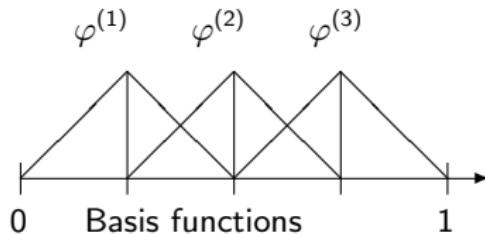
Continuous solution  $u(x) \rightarrow$  discrete solution  $\underline{u}_h$  from the finite dimensional space

$$V_h = \text{span} \left\{ \varphi^{(i)} : i \in \bar{\omega}_h \right\} = \left\{ v_h = \sum_{i \in \bar{\omega}_h} v^{(i)} \varphi^{(i)} \right\} = \text{span } \bar{\Phi} \subset V \quad (3)$$

spanned by the (linear independent) basis functions

$\bar{\Phi} = [\varphi^{(i)} : i \in \bar{\omega}_h] = [\varphi_1, \dots, \varphi_{N_h}]$  with  $\bar{\omega}_h$  as indices of basis functions.

1D linear basis functions with **finite** support on the neighboring **elements** are presented in the following picture:



## Our example: Laplace equation

$$\begin{aligned} \text{Find } u \text{ such that} \quad -\Delta u(x) &= f(x) & \forall x \in \Omega = [0, 1]^2 \\ u(x) &= 0 & \forall x \in \partial\Omega \end{aligned}$$

variational  $\Downarrow$  formulation

$$\begin{aligned} \text{Find } u \in \mathbb{V}: \quad a(u, v) &:= \int_{\Omega} \nabla^T v(x) \cdot \nabla u(x) dx \\ \langle F, v \rangle &:= \int_{\Omega} f(x) v(x) dx \end{aligned}$$

FEM, FDM  $\Downarrow$  FVM, FIT

$$\text{Solve} \quad K_h \cdot \underline{u}_h = \underline{f}_h \quad \underline{u}_h \in \mathbb{R}^{N_h}$$

$$\text{with } K^{ij} := \int_{\Omega} \nabla^T \varphi_j(x) \cdot \nabla \varphi_i(x) dx = \sum_{\tau_e \in \text{supp } \varphi_i \cap \text{supp } \varphi_j} \int_{\tau_e} \nabla^T \varphi_j(x) \cdot \nabla \varphi_i(x) dx$$

## How to solve Laplace equation?

1. Generate a finite element mesh.
2. Determine matrix pattern (sparse matrix!) and allocate storage.
3. Calculate Matrix  $K_h$  and r.h.s.  $\underline{f}_h$  for each element.

$$\int_{\tau_e} \nabla^T \varphi_j(x) \cdot \nabla \varphi_i(x) dx$$

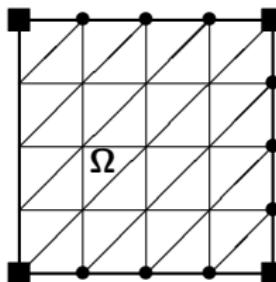
4. Accumulate the element entries.

$$\sum_{\tau_e \in \text{supp } \varphi_i \cap \text{supp } \varphi_j}$$

5. Solve the system of equations  $K_h \cdot \underline{u}_h = \underline{f}_h$ .

## Discretizing the domain $[xl, xr] \times [yb, yt]$

$nx=ny=4$  intervals



- ▶ triangular elements
- ▶ linear shape functions
- ▶ `GetMesh(nx, ny, xl, xr, yb, yt, nnode, xc, nelem, ia);`
- ▶ OUTPUT:
  - ▶ `nnode` : number of nodes
  - ▶ `xc[2*nnode]` : node coordinates
  - ▶ `nelem` : number of finite elements
  - ▶ `ia[3*nelem]` : element connectivity (3 node numbers per element)

## Storing the sparse matrix

**CRS**: compressed row storage The matrix

$$K_{n \times m} = \begin{pmatrix} 10 & 0 & 0 & -2 \\ 3 & 9 & 0 & 0 \\ 0 & 7 & 8 & 7 \\ 3 & 0 & 8 & 7 \end{pmatrix}$$

can be stored using just two integer vectors and one real/double vector.

Values :  $sk =$ 

10	-2	3	9	7	8	7	3	8	7
----	----	---	---	---	---	---	---	---	---

Column index :  $ik =$ 

1	4	1	2	2	3	4	1	3	4
---	---	---	---	---	---	---	---	---	---

Starting index of row :  $id =$ 

1	3	5	8	11
---	---	---	---	----

Dimensions for  $n$  rows and  $nnz$  non-zero elements in matrix:

$sk[nn]$  ,  $ik[nn]$  ,  $id[n+1]$

Note that (in C/C++)  $id[n] = nnz$ .

also: Compressed Column Storage (CCS), Compressed Diagonal Storage (CDS), Jagged Diagonal Storage (JDS),

ELLPACK, ...

## Matrix generation in code

- ▶ Determine matrix pattern and allocate memory for CRS

```
Get_Matrix_Pattern(nelem, 3, ia, nnz, id, ik, sk);
```

- ▶ nnz : number of non-zero elements in matrix
- ▶ id [nnode+1], ik [nnz] allocated and initialized
- ▶ sk [nnz] allocated

- ▶ Calculate Matrix entries and accumulate them

```
GetMatrix (nelem, 3, ia, nnode, xc, nnz, id, ik, sk, f);
```

- ▶ sk [nnz] matrix values initialized
- ▶ f [nnode] r.h.s. initialized

- ▶ Apply Dirichlet boundary conditions

```
ApplyDirichletBC(nx, ny, neigh, u, id, ik, sk, f);
```

- ▶ sk [nnz] matrix values adapted to B.C.
- ▶ f [nnode] r.h.s. adapted to B.C.
- ▶ nx, ny represent the geometry a input
- ▶ neigh represents neighboring domains in parallel context

## Solve the system of equations via Jacobi iteration

We solve  $K\underline{u} = \underline{f}$  by the Jacobi iteration ( $\omega = 1$ )

$$\underline{u}^{k+1} := \underline{u}^{k+1} + \omega D^{-1} (\underline{f} - K \cdot \underline{u}^k)$$

`JacobiSolve(nnnode, id, ik, sk, f, u );`

until the relative error in the  $KD^{-1}K$ -norm is smaller than  $\varepsilon = 10^{-5}$ .

```
D := diag(K)
\u := 0
\u_r := \u - K \cdot \u^0
\u_w := D^{-1} \cdot \u_r
\sigma := \sigma_0 := (\u_w, \u_r)
k := 0
while \sigma > \varepsilon^2 \cdot \sigma_0 do
    k := k + 1
    \u^k := \u^{k-1} + \omega \cdot \u_w           // vector arithmetics
    \u_r := \u - K \cdot \u^k                  // sparse matrix-times-vector + vector arithmetics
    \u_w := D^{-1} \cdot \u_r                  // vector arithmetics
    \sigma := (\u_w, \u_r)                    // inner product
end
```

# Data Parallelism for distributed memory

## Decomposing the mesh

The f.e. mesh is partitioned into  $P$  non-overlapping subdomains.  
(METIS, PARMETIS; SCOTCH, PT-SCOTCH)

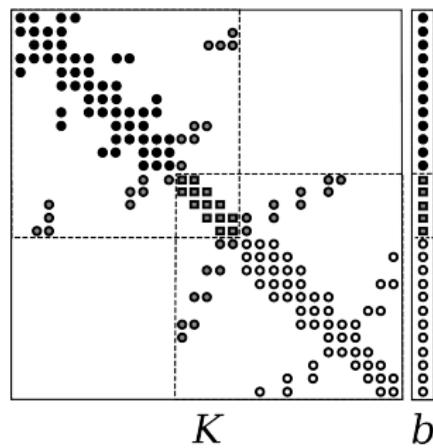
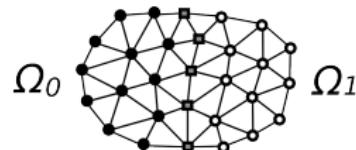
Unique mapping of an element to exactly one subdomain.

Decompose linear system

$$K^{ij} = \sum_{\tau_h} \int \nabla \varphi_i \cdot \nabla \varphi_j$$

into two subsystems  $K_0$  and  $K_1$ :

1. Non-overlapping decomposition of finite elements.
2. Overlapping nodes on boundary between subdomains.



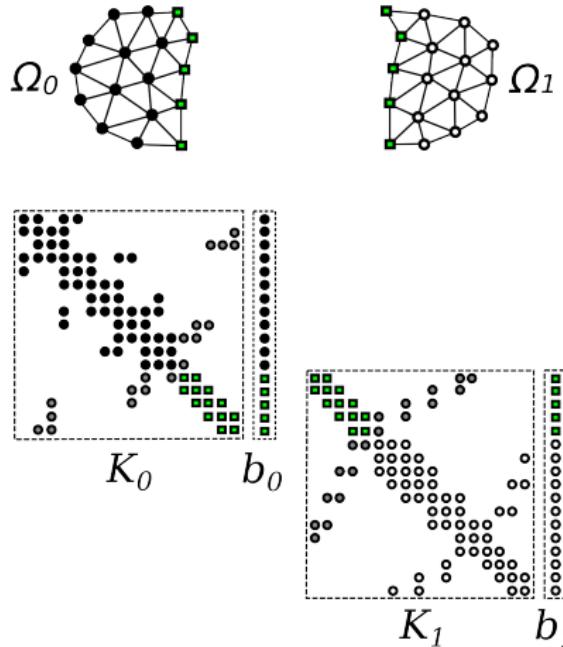
## Decomposition of matrix I

Local system

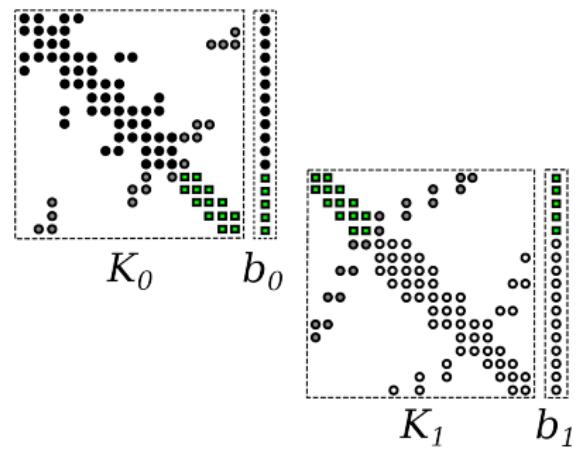
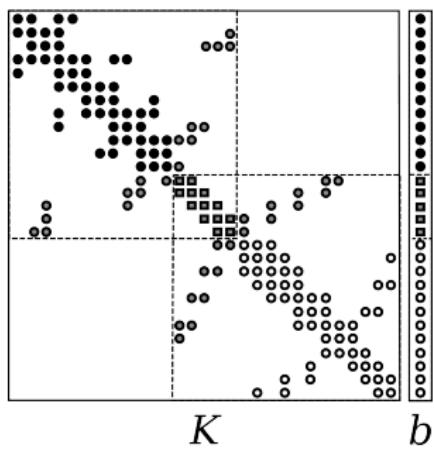
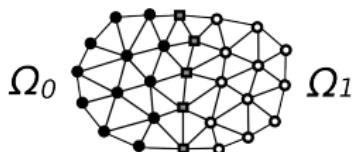
$$K_s^{ij} = \sum_{\tau_h \cap \Omega_s} \int_{\tau_h} \nabla \varphi_i \cdot \nabla \varphi_j$$

assembled locally:

- ▶ Distribute geometry
- ▶ Compute local stiffness matrix
- ▶ Assemble local *distributed* equation system.

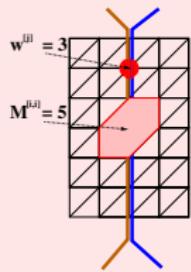


## Decomposition of matrix II



## Data representations

accumulated

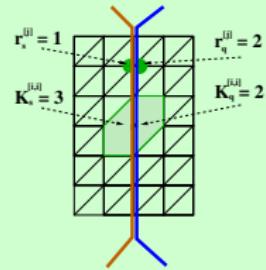


$$\underline{u}_s = A_s \underline{u}$$

$$\mathcal{K}_s = A_s \mathcal{K} A_s^T$$

$$\mathcal{K}^{ij} = \sum_{\tau_h} \int_{\tau_h} \nabla \varphi_i \cdot \nabla \varphi_j$$

distributed



$$\underline{r} = \sum_{s=1}^P A_s^T \underline{r}_s$$

$$\mathbf{K} = \sum_{s=1}^P A_s^T \mathbf{K}_s A_s$$

$$\mathbf{K}_s^{ij} = \sum_{\tau_h \cap \Omega_s} \int_{\tau_h} \nabla \varphi_i \cdot \nabla \varphi_j$$

# Parallel Linear Algebra

Global-to-local map

$$A_i = \begin{bmatrix} 1 & & & \\ & \ddots & & \\ & & 1 & \\ & & & \end{bmatrix}$$

Scalar product

$$\langle \textcolor{red}{w}, \textcolor{green}{r} \rangle = \textcolor{red}{w}^T \cdot \textcolor{green}{r} = \textcolor{red}{w}^T \cdot \sum_{i=1}^P A_i^T \textcolor{green}{r}_i = \sum_{i=1}^P (A_i \textcolor{red}{w})^T \textcolor{green}{r}_i = \sum_{i=1}^P \langle \textcolor{red}{w}_i, \textcolor{green}{r}_i \rangle$$

Matrix-vector product

$$\textcolor{green}{f} := \sum_{i=1}^P A_i^T \textcolor{green}{f}_i = \sum_{i=1}^P A_i^T \textcolor{green}{K}_i \textcolor{red}{u}_i = \sum_{i=1}^P A_i^T \textcolor{green}{K}_i A_i \textcolor{red}{u} = \textcolor{green}{K} \cdot \textcolor{red}{u}$$

Jacobi iteration

$$\textcolor{red}{u} := \textcolor{red}{u} + \omega \mathfrak{D}^{-1} \sum_{k=1}^P A_k^T (\textcolor{green}{f}_k - \textcolor{green}{K}_k \textcolor{red}{u}_k)$$

# Parallel Linear Algebra

no communication

$$\begin{aligned}\underline{v} &\leftarrow \mathbf{K} \cdot \underline{s} \\ \underline{r} &\leftarrow \underline{f} + \alpha \cdot \underline{v} \\ \underline{w} &\leftarrow \underline{u} + \alpha \cdot \underline{s} \\ \underline{r} &\leftarrow R^{-1} \cdot \underline{w}\end{aligned}$$

global communication

$$\langle \underline{w}, \underline{r} \rangle = \sum_{s=1}^P \langle \underline{w}_s, \underline{r}_s \rangle$$

next neighbor comm.

$$\begin{aligned}\underline{r}_s &\leftarrow A_s \sum_{k=1}^P A_k^T \underline{r}_k \\ \mathfrak{R}_s &\leftarrow A_s \left( \sum_{k=1}^P A_k^T \mathbf{K}_k A_k \right) A_s^T\end{aligned}$$

$$R = \text{diag}\{R_{ii}\}_{i=1}^N = \sum_{s=1}^P A_s \cdot A_s^T$$

$$\text{and } \mathbf{R}^{-1} \equiv \mathbf{I} = \sum_{s=1}^P A_s \mathbf{I}_s A_s^T \text{ (partition of unity)}$$

## Our example: Domain Decomposition

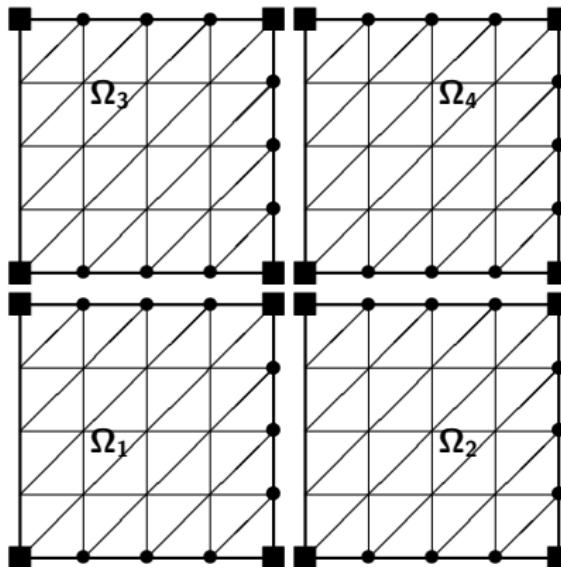


Figure : Non-overlapping elements.

## Parallel matrix generation

- ▶ Each process  $s$  possesses the elements of  $\Omega_s$ .

```
GetMesh(nx, ny, xl, xr, yb, yt, nnodes, xc, nelem, ia);
```

with individual  $xl$ ,  $xr$ ,  $yb$ ,  $yt$  in our example

- ▶ The local (distributed) matrix

$$K_s^{ij} := \sum_{\tau_h \cap \Omega_s} \int_{\tau_h} \nabla \varphi_i \cdot \nabla \varphi_j$$

is calculated by using directly the sequential routines

```
Get_Matrix_Pattern(nelem, 3, ia, nnz, id, ik, sk);
```

```
GetMatrix (nelem, 3, ia, nnodes, xc, nnz, id, ik, sk, f);
```

```
ApplyDirichletBC(nx, ny, neigh, u, id, ik, sk, f);
```

## Parallel Jacobi iteration for decomposed domain

We solve  $K\underline{u} = \underline{f}$  by the Jacobi iteration ( $\omega = 1$ )

$$\underline{u}^{k+1} := \underline{u}^{k+1} + \omega D^{-1} (\underline{f} - K \cdot \underline{u}^k)$$

on  $P$  processes with distributed data.

```
JacobiSolve(nnod, id, ik, sk, f, u );
❶  $\mathfrak{D} := \sum_{s=1}^P A_s^T \text{diag}(\mathbf{K}_s) A_s$  // next neighbor comm. of a vector
❷  $\underline{u} := 0$ 
❸  $\underline{r} := \underline{f} - \mathbf{K} \cdot \underline{u}^0$ 
❹  $\underline{w} := \mathfrak{D}^{-1} \cdot \sum_{s=1}^P A_s^T \underline{r}_s$  // next neighbor comm.
❺  $\sigma := \sigma_0 := (\underline{w}, \underline{r})$  // parallel reduction
❻  $k := 0$ 

while  $\sigma > \varepsilon^2 \cdot \sigma_0$  do
     $k := k + 1$ 
     $\underline{u}^k := \underline{u}^{k-1} + \omega \cdot \underline{w}$  // no comm.
     $\underline{r} := \underline{f} - \mathbf{K} \cdot \underline{u}^k$  // no comm.
     $\underline{w} := \mathfrak{D}^{-1} \cdot \sum_{s=1}^P A_s^T \underline{r}_s$  // next neighbor comm.
     $\sigma := (\underline{w}, \underline{r})$  // parallel reduction
end
```