

Parallel Scientific Computing

Course AMS301 — Fall 2022 — Lecture 7

Algebraic systems resulting from **finite element** discretizations

Algebraic systems resulting from finite element discretizations . . .

Problems with *algorithmic structures more complicated*

Writing/Implementing these algorithms require *manipulations of graphs*

Parallel implementation more complicated

Problem considered for this session

Let $\Omega \subset \mathbb{R}^d$ be an open bounded *domain* with a sufficiently regular border $\partial\Omega$
data $f \in L^2(\Omega)$ and $g \in L^2(\partial\Omega)$

$$\text{Find } u \in H^1(\Omega) \text{ such that } \begin{cases} -\Delta u + u = f & \text{dans } \Omega \\ \partial_n u|_{\partial\Omega} = g & \text{sur } \partial\Omega \end{cases}$$

After discretization with finite elements . . .

$$\text{Find } \mathbf{x} \in \mathbb{R}^N \text{ such that } \boxed{\mathbf{Ax} = \mathbf{f}}$$

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

*How to take advantage of the (sparse) structure of \mathbf{A}
for solving this problem in parallel?*

Algebraic systems resulting from finite element discretizations

Recap on finite elements

Sequential implementation

Parallel implementation

Recap on finite elements — Formulation

Exact differential formulation (DF)

$$\text{Find } u \in H^1(\Omega) \text{ such that } \begin{cases} -\Delta u + u = f & \text{dans } \Omega \\ \partial_n u|_{\partial\Omega} = g & \text{sur } \partial\Omega \end{cases}$$



Equivalent formulations

Exact variational formulation (VF)

Find $u \in H^1(\Omega)$ such that

$$\int_{\Omega} \nabla u \cdot \nabla v \, d\Omega + \int_{\Omega} uv \, d\Omega = \int_{\Omega} fv \, d\Omega + \int_{\partial\Omega} gv \, d\Omega, \quad \forall v \in H^1(\Omega)$$



Galerkin approximation:

$$V \longrightarrow V_h \subset V$$

Approximate variational formulation (VF)

Find $u_h \in V_h$ such that

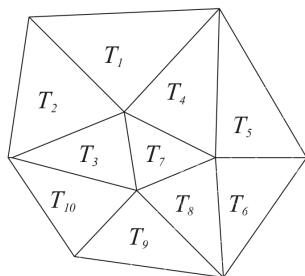
$$\int_{\Omega} \nabla u_h \cdot \nabla v_h \, d\Omega + \int_{\Omega} u_h v_h \, d\Omega = \int_{\Omega} f v_h \, d\Omega + \int_{\partial\Omega} g v_h \, d\Omega, \quad \forall v_h \in V_h$$

Recap on finite elements — Approximation space

For the sake of simplicity, we consider 2D cases with polygonal domains.

Mesh \mathcal{T}_h

- Set of cells/triangles $\mathcal{T}_h = (T_\ell)_{\ell=1 \dots L}$
- Set of vertices/nodes $\mathcal{M}_h = (M_i)_{i=1 \dots N}$
- Set of edges $\mathcal{E}_h = (E_a)_{a=1 \dots A}$
- $h_\ell =$ circumcircle diameter of T_ℓ
- $h = \max_\ell h_\ell =$ mesh step of \mathcal{T}_h



Properties

- $\bar{\Omega} = \bigcup_{\ell=1 \dots L} T_\ell$
- $T_\ell \cap T_m = \{\emptyset; 1 \text{ vertex}; 1 \text{ full edge}\} \rightarrow$ conformal mesh
- $\overset{\circ}{T}_\ell \neq \emptyset \rightarrow$ no flat triangle

Finite element P_k

For a given mesh \mathcal{T}_h , we define:

$$V_h := \{v_h \in C^0(\bar{\Omega}) : v_h|_{T_\ell} \in P_k(T_\ell), \ell = 1 \dots L\} \subset V \quad (\text{by construction})$$

where $P_k(T)$ is the space of polynomials of degree $\leq k$.

$$\begin{array}{l} \text{Find } u \in V \quad \text{such that } a(u, v) = b(v), \quad \forall v \in V \\ \text{Find } u_h \in V_h \quad \text{such that } a(u_h, v_h) = b(v_h), \quad \forall v_h \in V_h \end{array}$$

Exact problem

- Equivalence of formulation: u solution of (DF) $\Leftrightarrow u$ solution of (VF)
- Well-posedness: by Lax-Milgram Theorem

Approximate problem

- Well-posedness: by Lax-Milgram Theorem
- Convergence of the numerical solution:

Let $(\mathcal{T}_h)_h$, a regular family of meshes composed of elements P_k .

If $u \in H^{k+1}(\Omega)$ with $k \geq 1$, then there exist constants C_1 and C_2 such that, $\forall h$,

$$\|u - u_h\|_{L^2(\Omega)} \leq C_1 h^{k+1} |u|_{k+1, \Omega}$$

$$\|u - u_h\|_{H^1(\Omega)} \leq C_2 h^k |u|_{k+1, \Omega}$$

Recap on finite elements — Algebraic system [1/3]

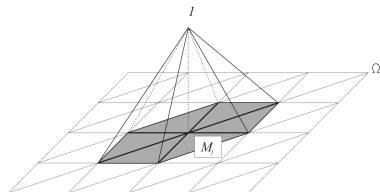
For the sake of simplicity, we consider P_1 finite elements.

Basis functions for V_h

- Lagrange functions $(\phi_i)_{i=1\dots N}$:

$$\phi_i \in V_h : \phi_i(M_j) = \delta_{ij}, \forall i, j$$

Property: $\text{supp}(\phi_i) = \bigcup_{\ell \text{ s.t. } M_i \subset T_\ell} T_\ell$



- The dimension of V_h is N . \rightarrow Number of vertices/nodes in the mesh
- The functions $(\phi_i)_{i=1\dots N}$ form a basis of V_h .
- Every solution $v_h \in V_h$ is characterized by the values $(v_h(M_i))_{i=1\dots N}$:

$$v_h(\mathbf{x}) = \sum_{i=1}^N v_h(M_i) \phi_i(\mathbf{x}), \quad \forall \mathbf{x} \in \Omega$$

Find $u_h \in V_h$ such that

$$\int_{\Omega} \nabla u_h \cdot \nabla v_h \, d\Omega + \int_{\Omega} u_h v_h \, d\Omega = \int_{\Omega} f v_h \, d\Omega + \int_{\partial\Omega} g v_h \, d\Omega, \quad \forall v_h \in V_h$$

Since $V_h = \text{span}(\phi_1, \dots, \phi_N)$, it is sufficient to use the basis functions as test functions:

$$\left| \begin{array}{l} \text{Find } u_h \in V_h \text{ such that} \\ \int_{\Omega} \nabla u_h \cdot \nabla \phi_i \, d\Omega + \int_{\Omega} u_h \phi_i \, d\Omega = \int_{\Omega} f \phi_i \, d\Omega + \int_{\partial\Omega} g \phi_i \, d\Omega, \quad i = 1 \dots N \end{array} \right.$$

The approximate solution can be written as

$$u_h(\mathbf{x}) = \sum_{j=1}^N \underbrace{u_h(M_j)}_{X_j} \phi_j(\mathbf{x})$$

Find $(X_j)_{j=1 \dots N} \in \mathbb{R}^N$ such that

$$\sum_{j=1}^N \left[\int_{\Omega} \nabla \phi_j \cdot \nabla \phi_i \, d\Omega + \int_{\Omega} \phi_j \phi_i \, d\Omega \right] X_j = \int_{\Omega} f \phi_i \, d\Omega + \int_{\partial\Omega} g \phi_i \, d\Omega, \quad i = 1 \dots N$$

Find $\mathbf{x} \in \mathbb{R}^N$ such that $\mathbf{Ax} = \mathbf{f}$

$$\text{with } \left\{ \begin{array}{ll} A_{ij} = \int_{\Omega} \nabla \phi_j \cdot \nabla \phi_i \, d\Omega + \int_{\Omega} \phi_j \phi_i \, d\Omega & (i, j = 1 \dots N) \\ f_i = \int_{\Omega} f \phi_i \, d\Omega + \int_{\partial\Omega} g \phi_i \, d\Omega & (i = 1 \dots N) \\ x_j = X_j & (j = 1 \dots N) \end{array} \right.$$

Properties

- \mathbf{A} is **symmetric positive definite**.

$$\begin{aligned} \text{For } \mathbf{y} \in \mathbb{R}^N \setminus \{0\} : (\mathbf{Ay} | \mathbf{y}) &= \sum_{i=1}^N \sum_{j=1}^N y_i a(\phi_i, \phi_j) y_j \\ &= a \left(\sum_{i=1}^N \phi_i y_i, \sum_{j=1}^N \phi_j y_j \right) && \text{(bilinearity of } a) \\ &= a(\mathbf{y}_h, \mathbf{y}_h) \geq \alpha_a \|\mathbf{y}_h\|^2 && \text{(coercivity of } a) \end{aligned}$$

- \mathbf{A} is (very) **sparse**.

$$\left. \begin{array}{l} A_{ij} \neq 0 \text{ if } \text{supp}(\phi_i) \cap \text{supp}(\phi_j) \neq \emptyset \\ \text{supp}(\phi_i) = \bigcup_{\ell \text{ s.t. } M_i \subset T_\ell} T_\ell \end{array} \right\} \Rightarrow A_{ij} \neq 0 \text{ if } \exists \ell \text{ such that } M_i, M_j \subset T_\ell$$

Algebraic systems resulting from finite element discretizations

Recap on finite elements

Sequential implementation

Parallel implementation

$$\boxed{\mathbf{Ax} = \mathbf{f}} \quad \text{with} \quad \begin{cases} A_{ij} = \int_{\Omega} \nabla \phi_j \cdot \nabla \phi_i \, d\Omega + \int_{\Omega} \phi_j \phi_i \, d\Omega & (i, j = 1, \dots, N) \\ f_i = \int_{\Omega} f \phi_i \, d\Omega + \int_{\partial\Omega} g \phi_i \, d\Omega & (i = 1, \dots, N) \end{cases}$$

Computation of matrix \mathbf{A}

The elements of \mathbf{A} can be rewritten as:

$$\begin{aligned} A_{ij} &= \int_{\Omega} (\nabla \phi_j \cdot \nabla \phi_i + \phi_j \phi_i) \, d\Omega \\ &= \sum_{\ell=1}^L \int_{T_{\ell}} (\nabla \phi_j \cdot \nabla \phi_i + \phi_j \phi_i) \, dT \\ &= \sum_{\substack{\ell \text{ such that} \\ M_i, M_j \subset T_{\ell}}} \int_{T_{\ell}} (\nabla \phi_j \cdot \nabla \phi_i + \phi_j \phi_i) \, dT \end{aligned}$$



For each element T_ℓ , we define three **local basis functions** $(\tau_I^\ell)_{I=1}^3$ such that:

$$\boxed{\phi_i|_{T_\ell} = \tau_I^\ell} \quad (I = 1, 2, 3)$$

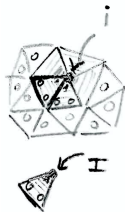
with $M_i \subset T_\ell$ and the corresponding indices $\text{LocalToGlobal}(\ell, I) = i$.

$$\boxed{\mathbf{Ax} = \mathbf{f}} \quad \text{with} \quad \left| \begin{array}{l} A_{ij} = \int_{\Omega} \nabla \phi_j \cdot \nabla \phi_i \, d\Omega + \int_{\Omega} \phi_j \phi_i \, d\Omega \quad (i, j = 1, \dots, N) \\ f_i = \int_{\Omega} f \phi_i \, d\Omega + \int_{\partial\Omega} g \phi_i \, d\Omega \quad (i = 1, \dots, N) \end{array} \right.$$

Computation of matrix \mathbf{A}

The elements of \mathbf{A} can be rewritten as:

$$\begin{aligned} A_{ij} &= \int_{\Omega} (\nabla \phi_j \cdot \nabla \phi_i + \phi_j \phi_i) \, d\Omega \\ &= \sum_{\ell=1}^L \int_{\hat{T}_{\ell}} (\nabla \phi_j \cdot \nabla \phi_i + \phi_j \phi_i) \, dT \\ &= \sum_{\substack{\ell \text{ such that} \\ M_i, M_j \subset T_{\ell}}} \int_{\hat{T}_{\ell}} (\nabla \phi_j \cdot \nabla \phi_i + \phi_j \phi_i) \, dT \\ &= \sum_{\substack{\ell \text{ such that} \\ M_i, M_j \subset T_{\ell}}} \int_{\hat{T}_{\ell}} (\nabla \tau_J^{\ell} \cdot \nabla \tau_I^{\ell} + \tau_J^{\ell} \tau_I^{\ell}) \, dT \quad \text{with} \quad \left| \begin{array}{l} \text{LocalToGlobal}(\ell, I) = i \\ \text{LocalToGlobal}(\ell, J) = j \end{array} \right. \\ &= \sum_{\substack{\ell \text{ such that} \\ M_i, M_j \subset T_{\ell}}} A_{IJ}^{\ell} \quad \text{with} \quad A_{IJ}^{\ell} = \int_{\hat{T}_{\ell}} (\nabla \tau_J^{\ell} \cdot \nabla \tau_I^{\ell} + \tau_J^{\ell} \tau_I^{\ell}) \, dT \end{aligned}$$



The matrix $\mathbf{A}^{\ell} \in \mathbb{R}^{3 \times 3}$ is a local element-wise matrix corresponding to element T_{ℓ} .

Assembling of \mathbf{A}

```

Initialization:  $\mathbf{A} = 0$ ;
for  $\ell = 1, \dots, L$  do
  | Computation of local matrix  $\mathbf{A}^\ell$ ;
  | for  $I = 1, 2, 3$  do
  | | for  $J = 1, 2, 3$  do
  | | |  $i \leftarrow \text{LocalToGlobal}(\ell, I)$ ;
  | | |  $j \leftarrow \text{LocalToGlobal}(\ell, J)$ ;
  | | |  $A_{ij} \leftarrow A_{ij} + A_{IJ}^\ell$ ;
  | | | end
  | | end
  | end
end
  
```

Assembling of \mathbf{f} (*volume term*)

```

Initialization:  $\mathbf{f} = 0$ ;
for  $\ell = 1, \dots, L$  do
  | Computation of local vector  $\mathbf{f}^\ell$ ;
  | for  $I = 1, 2, 3$  do
  | |  $i \leftarrow \text{LocalToGlobal}(\ell, I)$ ;
  | |  $f_i \leftarrow f_i + f_I^\ell$ ;
  | | end
  | end
end
  
```

Parallelization strategy?

Sequential implementation — Iterative solution procedure

Computation of a matrix-vector product $\mathbf{y} = \mathbf{Az}$

We would like to compute

$$y_i = \sum_{j=1}^N A_{ij} z_j \quad (i = 1, \dots, N)$$

where

- y_i is a resulting quantity associated to node M_i
- z_j is a quantity associated to node M_j (e.g. solution, residual, ...)
- $A_{ij} \neq 0$ only if there is at least one triangle containing M_i and M_j
i.e. if (M_i, M_j) is an edge $\in \mathcal{E}_h$.

Matrix-vector product $\mathbf{y} = \mathbf{Az}$

```
for  $M_i \in \mathcal{M}_h$  do
  for  $M_j \in \mathcal{M}_h$  such that  $(M_i, M_j) \in \mathcal{E}_h$  do
     $y_i \leftarrow y_i + A_{ij} z_j$ ;
  end
end
end
```

Parallelization strategy?

Algebraic systems resulting from finite element discretizations

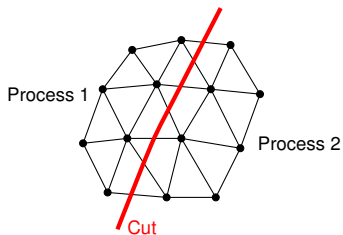
Recap on finite elements

Sequential implementation

Parallel implementation

Parallel implementation — Parallelization strategies

Partition by groups of vertices



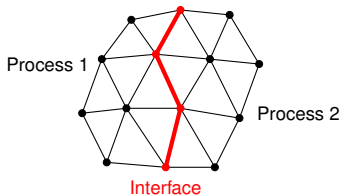
Parallel assembling

not natural

Parallel matrix-vector product

rather natural

Partition by groups of elements



Parallel assembling

rather natural

Parallel matrix-vector product

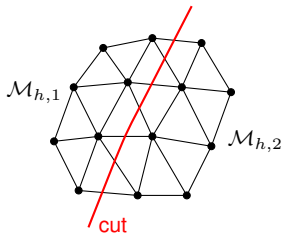
not natural

Parallel implementation — Strategy by groups of vertices [1/3]

The vertices/nodes are distributed between the different processes:

$$\mathcal{M}_h = \bigcup_{p=1}^P \mathcal{M}_{h,p} \quad \text{with } \mathcal{M}_{h,p} \cap \mathcal{M}_{h,q} = \emptyset \text{ if } p \neq q$$

where $\mathcal{M}_{h,p}$ is the group of vertices/nodes corresponding to process p .



Strategy for matrix-vector product $\mathbf{y} = \mathbf{A}\mathbf{z}$

- Each process p computes the part of \mathbf{y} corresponding to nodes $\mathcal{M}_{h,p}$:

$$y_i = \sum_j A_{ij} z_j \quad \text{with } i \in \mathcal{M}_{h,p}$$

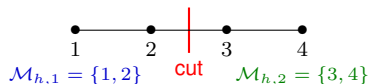
- Each process p stores the elements of \mathbf{z} and \mathbf{y} , and the lines of \mathbf{A} with indices $\in \mathcal{M}_{h,p}$.

A priori, no duplication of data, but computing \mathbf{y} requires communications.
The edges between nodes of $\mathcal{M}_{h,1}$ and $\mathcal{M}_{h,2}$ indicates the dependencies.

Parallel implementation — Strategy by groups of vertices [2/3]

Illustration in 1D

Configuration with 3 P1 elements and 4 nodes:



Matrix-vector product:

$$\begin{bmatrix} y_1 \\ y_2 \\ y_3 \\ y_4 \end{bmatrix} = \begin{bmatrix} \cdot & \cdot & 0 & 0 \\ \cdot & \cdot & \cdot & 0 \\ 0 & \cdot & \cdot & \cdot \\ 0 & 0 & \cdot & \cdot \end{bmatrix} \begin{bmatrix} z_1 \\ z_2 \\ z_3 \\ z_4 \end{bmatrix}$$

$$\begin{bmatrix} y_1 \\ y_2 \\ y_3 \\ y_4 \end{bmatrix} = \underbrace{\begin{bmatrix} \cdot & \cdot & 0 & 0 \\ \cdot & \cdot & \cdot & 0 \\ 0 & \cdot & \cdot & \cdot \\ 0 & 0 & \cdot & \cdot \end{bmatrix}}_{\substack{\text{Computed by process 1} \\ \mathbf{y}_1 = \mathbf{A}_1 \mathbf{z}}} \begin{bmatrix} z_1 \\ z_2 \\ z_3 \\ z_4 \end{bmatrix} + \underbrace{\begin{bmatrix} \cdot & \cdot & 0 & 0 \\ \cdot & \cdot & \cdot & 0 \\ 0 & \cdot & \cdot & \cdot \\ 0 & 0 & \cdot & \cdot \end{bmatrix}}_{\substack{\text{Computed by process 2} \\ \mathbf{y}_2 = \mathbf{A}_2 \mathbf{z}}} \begin{bmatrix} z_1 \\ z_2 \\ z_3 \\ z_4 \end{bmatrix}$$

Computing y_2 (on proc 1) requires z_1 and z_2 (on proc 1) and z_3 (on proc p_2)

Parallel algorithms

Parallel assembly of \mathbf{A} **On each process** $p = 1, \dots, P$:Assemble matrix \mathbf{A}_p corresponding to lines of \mathbf{A} with indices $i \in \mathcal{M}_{h,p}$;Parallel matrix-vector product $\mathbf{y} = \mathbf{A}\mathbf{z}$ **On each process** $p = 1, \dots, P$:**for** q such that $\mathcal{M}_{h,p} \cap \mathcal{M}_{h,q} \neq \emptyset$ **do** Send values $\{z_i\}_{i \in \mathcal{M}_{h,p}}$ s.t. $\exists(i, j) \in \mathcal{E}_h$ with $j \in \mathcal{M}_{h,q}$ to process q ; Recv values $\{z_j\}_{j \in \mathcal{M}_{h,q}}$ s.t. $\exists(i, j) \in \mathcal{E}_h$ with $i \in \mathcal{M}_{h,p}$ from process q ;**end****for** $i \in \mathcal{M}_{h,p}$ **do** **for** j such that (i, j) is an edge **do** $y_i \leftarrow y_i + A_{p,ij}z_j$; **end****end**

Temporary storage, to store nodal values corresponding to neighboring process

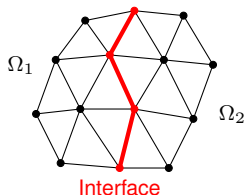
Parallel implementation — Strategy by groups of elements [1/3]

The elements are distributed between the different processes:

$$\bar{\Omega} = \bigcup_{p=1}^P \bar{\Omega}_p$$

where $\bar{\Omega}_p$ is the group of elements corresponding to process p .

If $\mathcal{M}_{h,p}$ is the set of nodes/vertices of Ω_p , then $\mathcal{M}_{h,p} \cap \mathcal{M}_{h,q} \neq \emptyset$ if $\bar{\Omega}_p \cap \bar{\Omega}_q \neq \emptyset$.



Strategy for matrix-vector product $\mathbf{y} = \mathbf{A}\mathbf{z}$

- ▶ Each process p performs the operations corresponding to the elements of $\bar{\Omega}_p$.
- ▶ Each process p stores elements of \mathbf{z} and \mathbf{y} and the lines of \mathbf{A} corresponding to vertices/nodes $\mathcal{M}_{h,p}$ (i.e. both interior and interface nodes).

Duplication of data corresponding to interface nodes

Parallel algorithms

Parallel assembly of \mathbf{A}

On each process $p = 1, \dots, P$:

Assemble matrix \mathbf{A}_p corresponding to elements of \mathcal{T}_p ;

Parallel matrix-vector product $\mathbf{y} = \mathbf{A}\mathbf{z}$

On each process $p = 1, \dots, P$:

for $i \in \mathcal{M}_{h,p}$ **do**

for $j \in \mathcal{M}_{h,p}$ *such that* $(i, j) \in \mathcal{E}_h$ **do**

$y_i \leftarrow y_i + A_{p,ij}z_j$;

end

end

for q *such that* $\mathcal{M}_{h,p} \cap \mathcal{M}_{h,q} \neq \emptyset$ **do**

 Send/Recv values $\{y_i\}$ for the interface nodes $\mathcal{M}_{h,p} \cap \mathcal{M}_{h,q}$;

 Accumulate these values to compute the total sums;

end

Summary

► Finite element scheme

- Exact/Approximate variational formulation of an elliptic problem
- P_1 finite elements — Convergence rate: h^2 in L^2 -norm and h^1 in H^1 -norm
- Linear system $\mathbf{Ax} = \mathbf{f}$:
 - $\mathbf{A} \in \mathbb{R}^{N \times N}$ is symmetric, positive definite, sparse
 - $\mathbf{x} \in \mathbb{R}^N$ contains the nodal values of the solution
 - N is the number of nodes/vertices

► Implementation

- Main loops:
 - Loop over the elements for the matrix assembly
 - Loop over the unknowns/nodes/vertices for solving the linear system
- Strategy for parallel implementation:
 - Partitioning by groups of nodes
 - Partitioning by groups of elements