

# Toward black-box adaptive domain decomposition methods

**Frédéric Nataf**

Laboratory J.L. Lions (LJLL), CNRS, Alpines Inria and Univ. Paris VI

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joint work with

**Victorita Dolean** (Univ. Nice Sophia-Antipolis)

**Patrice Hauret** (Michelin, Clermont-Ferrand)

**Frédéric Hecht** (LJLL)

**Pierre Jolivet** (LJLL)

**Clemens Pechstein** (Johannes Kepler Univ., Linz)

**Robert Scheichl** (Univ. Bath)

**Nicole Spillane** (LJLL)

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- 1 Some Applications
- 2 An abstract 2-level Schwarz: the GenEO algorithm
- 3 Conclusion

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- 2 An abstract 2-level Schwarz: the GenEO algorithm
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Large Sparse discretized system with  
**strongly heterogeneous coefficients**  
(high contrast, nonlinear, multiscale)

E.g. Darcy pressure equation,  
 $P^1$ -finite elements:

$$AU = F$$

$$\text{cond}(A) \sim \frac{\alpha_{\max}}{\alpha_{\min}} h^{-2}$$

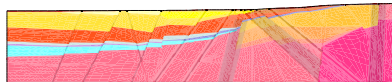
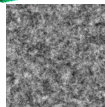
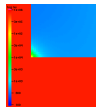
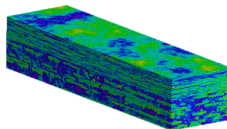
**Goal:**

iterative solvers

robust in **size** and **heterogeneities**

**Applications:**

flow in heterogeneous /  
stochastic / layered media  
structural mechanics  
time dependent waves  
etc.





$$-\operatorname{div}(\sigma(\mathbf{u})) = \mathbf{f},$$

where

$$\sigma_{ij}(\mathbf{u}) = 2\mu\varepsilon_{ij}(\mathbf{u}) + \lambda\delta_{ij}\operatorname{div}(\mathbf{u}), \quad \varepsilon_{ij}(\mathbf{u}) = \frac{1}{2} \left( \frac{\partial u_i}{\partial x_j} + \frac{\partial u_j}{\partial x_i} \right),$$
$$\mathbf{f} = (0, g)^T = (0, 10)^T, \quad \mu = \frac{E}{2(1+\nu)}, \quad \lambda = \frac{E\nu}{(1+\nu)(1-2\nu)}.$$

After a FEM (finite element method) discretization:

$$\Rightarrow \mathbf{AU} = \mathbf{F}$$

where  $\mathbf{A}$  is a large sparse highly heterogeneous matrix.

# Black box solvers (solve(MAT,RHS,SOL))

	Direct Solvers	Iterative Solvers
Pros	Robustness	Naturally
Cons	Difficult to	Robustness

Domain Decomposition Methods (DDM): Hybrid solver → should be naturally parallel and robust

General form:

$Au = f$ , solved with PCG for a preconditioner  $M^{-1}$ .

What's classical: Robustness with respect to problem size (scalability)

What's New here: Provable Robustness in the SPD case with respect to:

- Coefficients jumps
- System of PDEs

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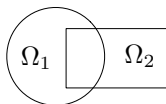
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# The First Domain Decomposition Method

## The original Schwarz Method (H.A. Schwarz, 1870)

$$\begin{aligned} -\Delta(u) &= f \quad \text{in } \Omega \\ u &= 0 \quad \text{on } \partial\Omega. \end{aligned}$$



Schwarz Method :  $(u_1^n, u_2^n) \rightarrow (u_1^{n+1}, u_2^{n+1})$  with

$$\begin{aligned} -\Delta(u_1^{n+1}) &= f \quad \text{in } \Omega_1 \\ u_1^{n+1} &= 0 \quad \text{on } \partial\Omega_1 \cap \partial\Omega \\ u_1^{n+1} &= u_2^n \quad \text{on } \partial\Omega_1 \cap \overline{\Omega_2}. \end{aligned}$$

$$\begin{aligned} -\Delta(u_2^{n+1}) &= f \quad \text{in } \Omega_2 \\ u_2^{n+1} &= 0 \quad \text{on } \partial\Omega_2 \cap \partial\Omega \\ u_2^{n+1} &= u_1^{n+1} \quad \text{on } \partial\Omega_2 \cap \overline{\Omega_1}. \end{aligned}$$

Parallel algorithm, converges but very slowly, overlapping subdomains only.

The parallel version is called **Jacobi Schwarz method (JSM)**.



# Jacobi and Schwarz (I): Algebraic point of view

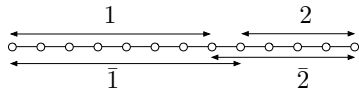
The set of indices is partitioned into two sets  $\mathcal{N}_1$  and  $\mathcal{N}_2$ :

$$\begin{pmatrix} A_{11} & A_{12} \\ A_{21} & A_{22} \end{pmatrix} \begin{pmatrix} x_1 \\ x_2 \end{pmatrix} = \begin{pmatrix} b_1 \\ b_2 \end{pmatrix}$$

The **block-Jacobi algorithm** reads:

$$\begin{pmatrix} A_{11} & 0 \\ 0 & A_{22} \end{pmatrix} \begin{pmatrix} x_1^{k+1} \\ x_2^{k+1} \end{pmatrix} = - \begin{pmatrix} 0 & A_{12} \\ A_{21} & 0 \end{pmatrix} \begin{pmatrix} x_1^k \\ x_2^k \end{pmatrix} + \begin{pmatrix} b_1 \\ b_2 \end{pmatrix}$$

It corresponds to solving a Dirichlet boundary value problem in each subdomain with Dirichlet data taken from the other one at the previous step  $\iff$  **Schwarz method with minimal overlap**



**Figure :** Domain decomposition with minimal overlap

# Jacobi and Schwarz (II): Larger overlap

Let  $\delta$  be a non negative integer

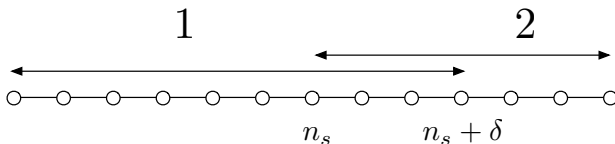


Figure : Domain decomposition with overlap

$$A = \begin{pmatrix} n_s & & \\ & \text{---} & \\ & & n_s + 1 \end{pmatrix} \text{ or } \begin{pmatrix} & & n_s + \delta \\ & \text{---} & \\ n_s & & \end{pmatrix}$$

Figure : Matrix decomposition Without or with overlap

# Strong and Weak scalability

How to evaluate the efficiency of a domain decomposition?

## Strong scalability (Amdahl)

"How the solution time varies with the number of processors for a fixed *total* problem size"

## Weak scalability (Gustafson)

"How the solution time varies with the number of processors for a fixed problem size *per processor*."

## Not achieved with the one level method

Number of subdomains	8	16	32	64
ASM	18	35	66	128

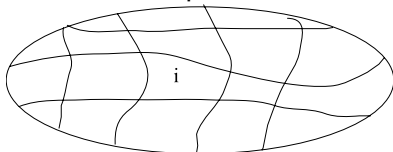
The iteration number increases linearly with the number of subdomains in one direction.

# How to achieve scalability

Stagnation corresponds to a few very low eigenvalues in the spectrum of the preconditioned problem. They are due to the lack of a global exchange of information in the preconditioner.

$$-\Delta u = f \text{ in } \Omega$$

$$u = 0 \text{ on } \partial\Omega$$



The mean value of the solution in domain  $i$  depends on the value of  $f$  on all subdomains.

A classical remedy consists in the introduction of a **coarse problem** that couples all subdomains. This is closely related to **deflation technique** classical in linear algebra (see Nabben and Vuik's papers in SIAM J. Sci. Comp, 200X) and multigrid techniques.

# Adding a coarse space

We add a coarse space correction (*aka* second level)

Let  $V_H$  be the coarse space and  $Z$  be a basis,  $V_H = \text{span } Z$ , writing  $R_0 = Z^T$  we define the two level preconditioner as:

$$M_{ASM,2}^{-1} := R_0^T (R_0 A R_0^T)^{-1} R_0 + \sum_{i=1}^N R_i^T A_i^{-1} R_i.$$

The **Nicolaides approach** is to use the kernel of the operator as a coarse space, this is the constant vectors, in local form this writes:

$$Z := (R_i^T D_i R_i \mathbf{1})_{1 \leq i \leq N}$$

where  $D_i$  are chosen so that we have a partition of unity:

$$\sum_{i=1}^N R_i^T D_i R_i = Id.$$

# Theoretical convergence result

## Theorem (Widlund, Dryja)

Let  $M_{ASM,2}^{-1}$  be the two-level additive Schwarz method:

$$\kappa(M_{ASM,2}^{-1}A) \leq C \left(1 + \frac{H}{\delta}\right)$$

where  $\delta$  is the size of the overlap between the subdomains and  $H$  the subdomain size.

This does indeed work very well

Number of subdomains	8	16	32	64
ASM	18	35	66	128
ASM + Nicolaides	20	27	28	27

# Failure for Darcy equation with heterogeneities

$$\begin{aligned} -\nabla \cdot (\alpha(\mathbf{x}, y) \nabla u) &= 0 \quad \text{in } \Omega \subset \mathbb{R}^2, \\ u &= 0 \quad \text{on } \partial\Omega_D, \\ \frac{\partial u}{\partial n} &= 0 \quad \text{on } \partial\Omega_N. \end{aligned}$$



Decomposition



$\alpha(\mathbf{x}, y)$

Jump	1	10	$10^2$	$10^3$	$10^4$
ASM	39	45	60	72	73
ASM + Nicolaides	30	36	50	61	65

## Our approach

Fix the problem by an optimal and proven choice of a coarse space  $\mathcal{Z}$ .

- 1 Some Applications
- 2 An abstract 2-level Schwarz: the GenEO algorithm
  - Choice of the coarse space
  - Parallel implementation
- 3 Conclusion



## Strategy

Define an appropriate coarse space  $V_{H2} = \text{span}(Z_2)$  and use the framework previously introduced, writing  $R_0 = Z_2^T$  the two level preconditioner is:

$$P_{ASM2}^{-1} := R_0^T (R_0 A R_0^T)^{-1} R_0 + \sum_{i=1}^N R_i^T A_i^{-1} R_i.$$

## The coarse space must be

- Local (calculated on each subdomain)  $\rightarrow$  parallel
- Adaptive (calculated automatically)
- Easy and cheap to compute
- Robust (must lead to an algorithm whose convergence is proven not to depend on the partition nor the jumps in coefficients)

# Abstract eigenvalue problem

**Gen.EVP** per subdomain:

Find  $p_{j,k} \in V_{h|\Omega_j}$  and  $\lambda_{j,k} \geq 0$ :

$$a_{\Omega_j}(p_{j,k}, v) = \lambda_{j,k} a_{\Omega_j^o}(\Xi_j p_{j,k}, \Xi_j v) \quad \forall v \in V_{h|\Omega_j}$$

$$A_j p_{j,k} = \lambda_{j,k} X_j A_j^o X_j p_{j,k} \quad (X_j \dots \text{diagonal})$$

$a_D \dots$  restriction of  $a$  to  $D$

---

**In the two-level ASM:**

Choose first  $m_j$  eigenvectors per subdomain:

$$V_0 = \text{span} \{ \Xi_j p_{j,k} \}_{k=1, \dots, m_j}^{j=1, \dots, N}$$

This automatically includes Zero Energy Modes.

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# Comparison with existing works

**Galvis & Efendiev (SIAM 2010):**

$$\int_{\Omega_j} \kappa \nabla p_{j,k} \cdot \nabla v \, dx = \lambda_{j,k} \int_{\Omega_j} \kappa p_{j,k} v \, dx \quad \forall v \in V_{h|\Omega_j}$$

**Efendiev, Galvis, Lazarov & Willems (submitted):**

$$a_{\Omega_j}(p_{j,k}, v) = \lambda_{j,k} \sum_{i \in \text{neighb}(j)} a_{\Omega_j}(\xi_j \xi_i p_{j,k}, \xi_j \xi_i v) \quad \forall v \in V_{|\Omega_j}$$

$\xi_j \dots$  partition of unity, calculated adaptively (MS)

**Our gen.EVP:**

$$a_{\Omega_j}(p_{j,k}, v) = \lambda_{j,k} a_{\Omega_j^\circ}(\Xi_j p_{j,k}, \Xi_j v) \quad \forall v \in V_{h|\Omega_j}$$

both matrices typically singular  $\implies \lambda_{j,k} \in [0, \infty]$

Two technical assumptions.

Theorem (Spillane, Dolean, Hauret, N., Pechstein, Scheichl)

If for all  $j$ :  $0 < \lambda_{j,m_{j+1}} < \infty$ :

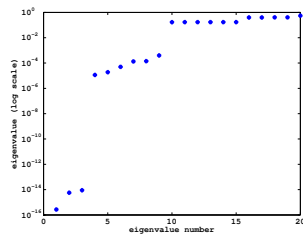
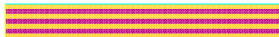
$$\kappa(M_{ASM,2}^{-1}A) \leq (1 + k_0) \left[ 2 + k_0 (2k_0 + 1) \max_{j=1}^N \left( 1 + \frac{1}{\lambda_{j,m_{j+1}}} \right) \right]$$

Possible criterion for picking  $m_j$ : (used in our Numerics)

$$\lambda_{j,m_{j+1}} < \frac{\delta_j}{H_j}$$

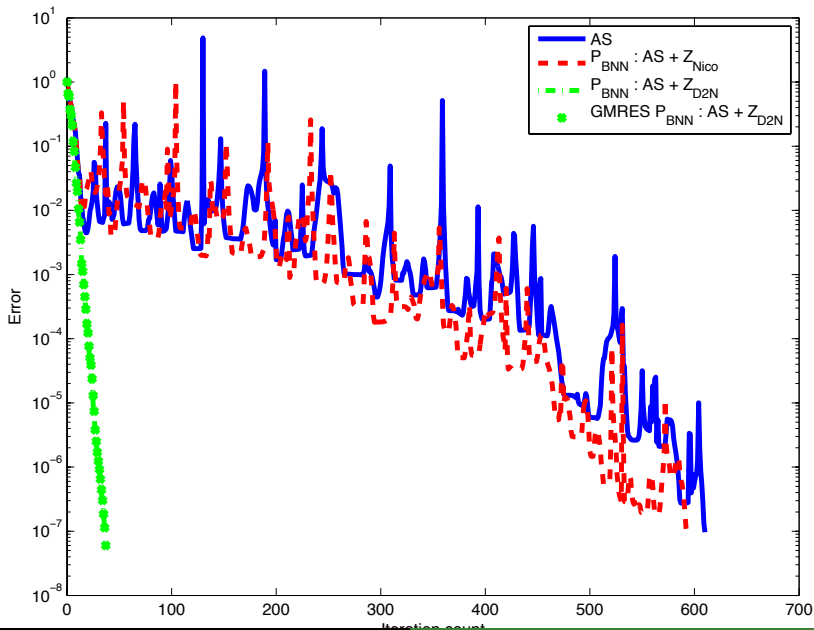
$H_j \dots$  subdomain diameter,  $\delta_j \dots$  overlap

# Eigenvalues and eigenvectors



Logarithmic scale

# Convergence



# Numerical results via a Domain Specific Language

FreeFem++ (<http://www.freefem.org/ff++>), with:

- Metis Karypis and Kumar 1998
- SCOTCH Chevalier and Pellegrini 2008
- UMFPACK Davis 2004
- ARPACK Lehoucq et al. 1998
- MPI Snir et al. 1995
- Intel MKL
- PARDISO Schenk et al. 2004
- MUMPS Amestoy et al. 1998
- PaStiX Hénou et al. 2005
- PETSC

Why use a DS(E)L instead of C/C++/Fortran/... ?

- performances close to low-level language implementation,
- hard to beat something as simple as:

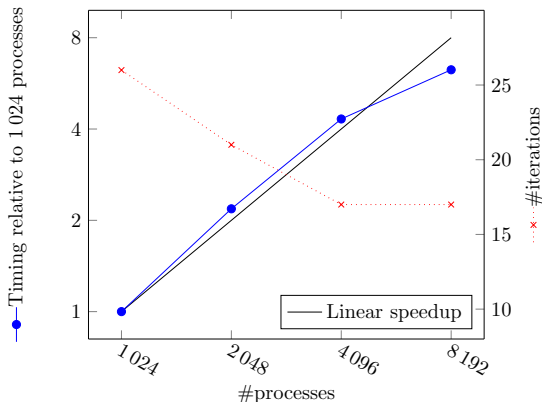
```
varf a(u, v) = int3d(mesh)([dx(u), dy(u), dz(u)]' * [dx(v), dy(v), dz(v)])  
              + int3d(mesh)(f * v) + on(boundary_mesh)(u = 0)
```

- Much More in F. Hecht's talk tomorrow



# Strong scalability in two dimensions heterogeneous elasticity (P. Jolivet with Frefeem ++)

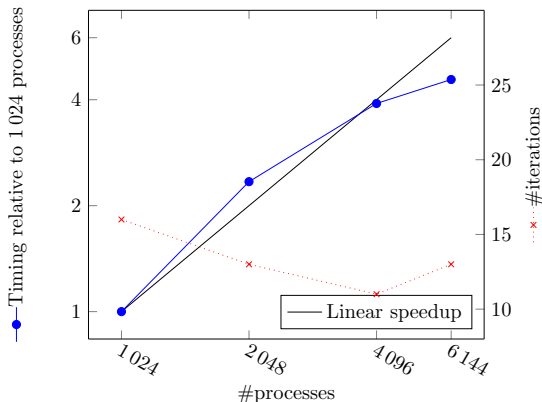
## Elasticity problem with heterogeneous coefficients



Speed-up for a 1.2 billion unknowns 2D problem. Direct solvers in the subdomains. Peak performance wall-clock time: 26s.

# Strong scalability in three dimensions heterogeneous elasticity

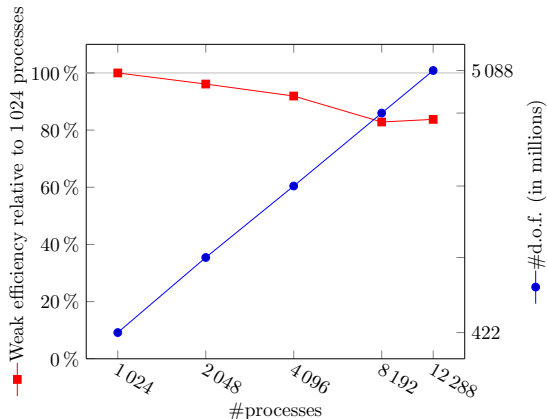
## Elasticity problem with heterogeneous coefficients



Speed-up for a 160 million unknowns 3D problem. Direct solvers in subdomains. Peak performance wall-clock time: 36s.

# Weak scalability in two dimensions

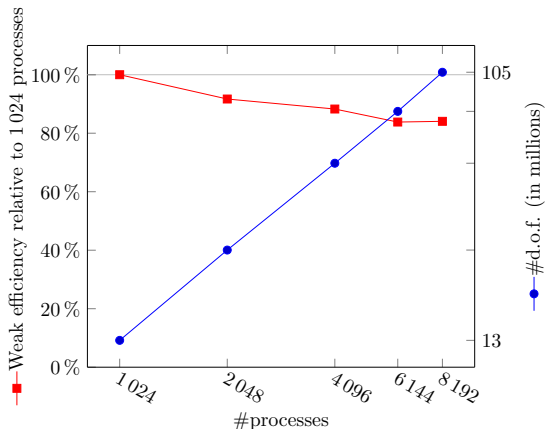
## Darcy problems with heterogeneous coefficients



Efficiency for a 2D problem. Direct solvers in the subdomains.  
Final size: 22 billion unknowns. Wall-clock time:  $\simeq$  200s.

# Weak scalability in three dimensions

## Darcy problems with heterogeneous coefficients



Efficiency for a 3D problem. Direct solvers in the subdomains.  
Final size: 2 billion unknowns. Wall-clock time:  $\simeq$  200s.

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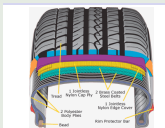
# Conclusion

## Summary

- Using generalized eigenvalue problems and projection preconditioning we are able to achieve a targeted convergence rate.
- Works for ASM, BNN and FETI methods
- This process can be implemented in a black box algorithm.

## Future work

- Build the coarse space on the fly.



- Multigrid like three (or more) level methods

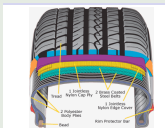
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**THANK YOU FOR YOUR ATTENTION!**