

Resolution of PDE using domain decomposition methods on TeraFLOPic architectures

Pierre Jolivet

Laboratoire Jacques-Louis Lions^{*}
Laboratoire Jean Kuntzmann[‡]

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Third workshop on FreeFem++

Work supervised by F. Nataf^{*}, C. Prud'Homme[‡], F. Hecht^{*}

Outline

- 1 Introduction
- 2 Clusters
- 3 Domain decomposition methods
 - One-level methods
 - Two-level methods
 - Numerical results
- 4 Conclusion

Context

We want to solve large systems arising from the finite element method.

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- parallel direct solvers (MUMPS, SuperLU ..),
- parallel iterative solvers (Hypre ..),
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⇒ high-performance algorithms on massively parallel distributed memory multiprocessor architectures.

Within FreeFem++

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C++ plus load function within FreeFem++:

⇒ OpenMP (shared memory architectures), C for CUDA (GPGPU).

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FreeFem++ is working on the following parallel architectures (amongst others):

	N° of cores	Memory	Peak performance	Compilers
hpc1@LJLL	64@2.00 Ghz	252 Go	< 1 TFLOP/s	Intel
titane@CEA	12192* @2.93 Ghz	37 To	140 TFLOP/s	Intel
babel@IDRIS	40960@850 Mhz	20 To	139 TFLOP/s	IBM+GNU

* + 46080 CUDA cores

<http://www-ccrt.cea.fr>, Bruyères-le-Châtel, France.

<http://www.idris.fr>, Orsay, France.

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The original Schwarz method for a 2-way decomposition

Consider the following BVP in \mathbb{R}^d ($d = 2, 3$):

$$\begin{aligned}\nabla \cdot \kappa \nabla u &= F(u) && \text{in } \Omega \\ B(u) &= 0 && \text{on } \partial\Omega\end{aligned}$$

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$$\begin{array}{llll}\nabla \cdot \kappa \nabla u_1^{n+1} = F(u_1^{n+1}) & \text{in } \Omega_1 & \nabla \cdot \kappa \nabla u_2^{n+1} = F(u_2^{n+1}) & \text{in } \Omega_2 \\ B(u_1^{n+1}) = 0 & \text{on } \partial\Omega_1 \cap \partial\Omega & B(u_2^{n+1}) = 0 & \text{on } \partial\Omega_2 \cap \partial\Omega \\ u_1^{n+1} = u_2^n & \text{on } \partial\Omega_1 \cap \overline{\Omega_2} & u_2^{n+1} = u_1^n & \text{on } \partial\Omega_2 \cap \overline{\Omega_1}\end{array}$$

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- ② Highly heterogeneous coefficient κ in the BVP,
 \implies long plateaux in the convergence of the algorithm.

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Construction of E

- ① build a coarse mesh $\mathcal{T}_{\text{coarse}}$, then a new fespace and a new varf on $\mathcal{T}_{\text{coarse}}$ (available under `examples++-mpi/MPIGMRES*D.edp`).
- ② use the low-frequency modes of the *Dirichlet-to-Neumann* operator of the BVP at the interface of each neighboring domains (Nataf *et al.* , 2011).

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Two-level methods

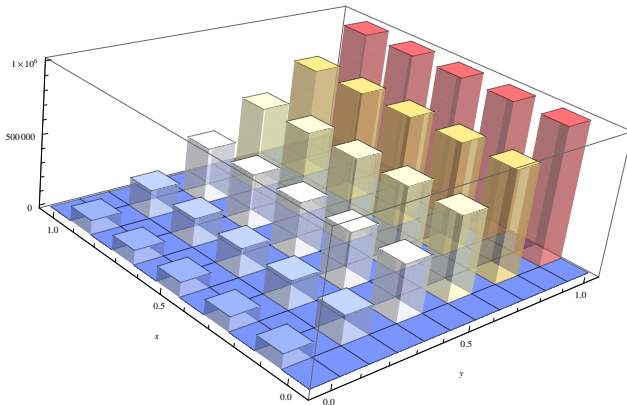
- ① solve the global coarse problem $Ex = b$ on one node,
- ② use the solution of the coarse problem on each local fine problem.

This is used to precondition a Krylov method (CG, GMRES) (Tang *et al.* , 2009).

Test case in \mathbb{R}^2

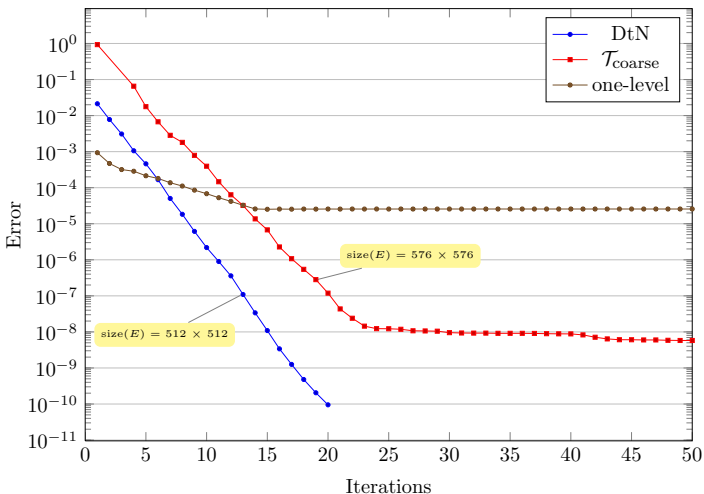
$$\begin{aligned}\nabla \cdot \kappa \nabla u &= 1 & \text{in } \Omega \\ u &= 0 & \text{on } \partial\Omega\end{aligned}$$

with a *skyscraper* viscosity $\kappa(x, y)$:



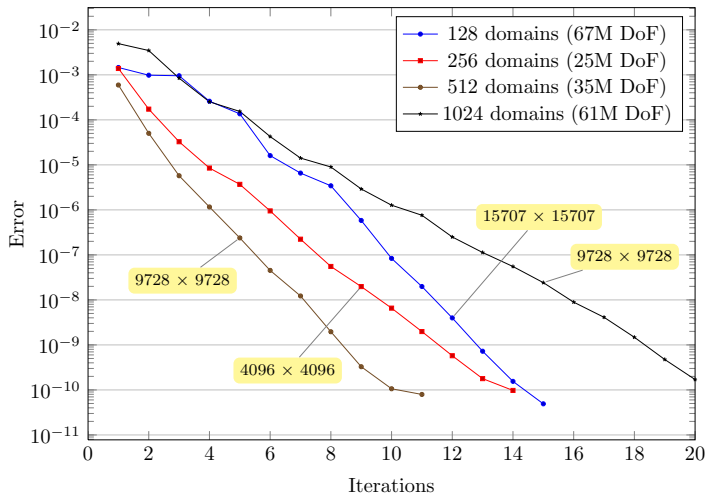
Convergence curves

2D test case running on 64 processors using \mathbb{P}_2 FE ($\varepsilon = 10^{-10}$, 450k DoF)



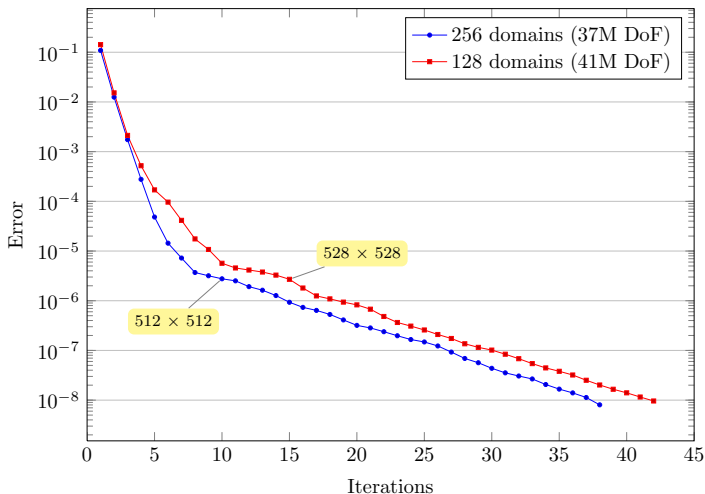
Some more 2D results

2D test case running on various number of processors using \mathbb{P}_1 FE ($\varepsilon = 10^{-10}$)



And some 3D results

3D test case running on various number of processors using \mathbb{P}_2 FE ($\varepsilon = 10^{-8}$)



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Final words

Concerning FreeFem++:

- ① efficient parallel algorithms (memory and CPU overhead with the FreeFem++ - MPI bindings),
- ② not very user-friendly interface at the moment (for DDM),
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Concerning domain decomposition methods:

- ① add another level,
- ② do parallel computations on more complex systems (e.g. elasticity).

Thanks for your attention.

NATAF, F., XIANG, H., DOLEAN, V., & SPILLANE, N. 2011.

A coarse space construction based on local Dirichlet to Neumann maps, to appear.

SIAM Journal on Scientific Computing.

TANG, J.M., NABBEN, R., VUIK, C., & ERLANGGA, Y.A. 2009.

Comparison of two-level preconditioners derived from deflation, domain decomposition and multigrid methods.

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