Domain decomposition techniques for hyperbolic equations on unstructured grids

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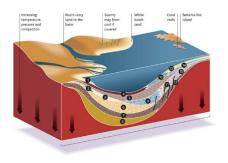


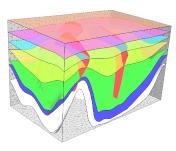




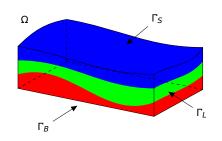
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Motivation - sedimentary basins





Mathematical model



$$\begin{cases} \nabla \cdot (\mu(\nabla \boldsymbol{u} + \nabla \boldsymbol{u}^T)) - \nabla \rho = -\rho \boldsymbol{g} & \text{in } \Omega \times (0, T] \\ \nabla \cdot \boldsymbol{u} = 0 & \text{in } \Omega \times (0, T] \\ \frac{\partial}{\partial t} \{\rho, \mu\} + \boldsymbol{u} \cdot \nabla \{\rho, \mu\} = 0 & \text{in } \Omega \times (0, T] \\ \rho = \rho_0, \quad \mu = \mu_0 & \text{in } \Omega \times \{0\} \\ \boldsymbol{u} = \bar{\boldsymbol{u}} & \text{on } \Gamma \end{cases}$$

Stokes system discretization - 1

bilinear forms

$$a(\mathbf{u}, \mathbf{v}) = \int_{\Omega} \mu \nabla \mathbf{u} : \nabla \mathbf{v} \qquad \forall \mathbf{u}, \mathbf{v} \in \mathbf{H}^{1}$$

$$b(p, \mathbf{v}) = \int_{\Omega} p \nabla \cdot \mathbf{v} \qquad \forall p \in L_{0}^{2}, \quad \forall \mathbf{v} \in \mathbf{H}^{1}$$

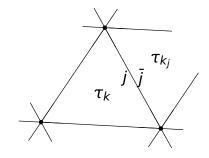
$$f(\mathbf{v}) = -\int_{\Omega} \rho \mathbf{g} \cdot \mathbf{v} \qquad \forall \mathbf{v} \in \mathbf{H}^{1}$$

weak formulation

$$\begin{cases} a(\mathbf{u}, \mathbf{v}) + b(p, \mathbf{v}) = f(\mathbf{v}) & \forall \mathbf{v} \in \mathbf{X} \\ b(q, \mathbf{u}) = 0 & \forall q \in \mathbf{S} \end{cases}$$

Stokes system discretization - 2

- tetrahedral grid $\mathcal{T}_h(\Omega)$, n_e elements and n_p points
- element τ_k , $\bigcup_k \tau_k = \mathcal{T}_h$
- $\blacksquare \ X = \{ \nu_h \in H^1 : \nu_h|_{\tau_k} \in \mathbb{P}^1_b \}$
- $S = \{q_h \in L_0^2 : v_h|_{\tau_k} \in \mathbb{P}^1\}$



Characteristic function

$$\lambda_i(\mathbf{x}) = \begin{cases} 1 & \text{if } \mathbf{x} \in \Omega_i \\ 0 & \text{if } \mathbf{x} \notin \Omega_i \end{cases}$$

$$\rho = \sum_{i=1}^{s} \lambda_i \rho_i, \quad \mu = \sum_{i=1}^{s} \lambda_i \mu_i$$

i = -, -, - (s components)

evolution equation for λ

$$\frac{\partial \boldsymbol{\lambda}}{\partial t} + \boldsymbol{u}^n \cdot \nabla \boldsymbol{\lambda} = 0$$

Characteristic function discretization - 1

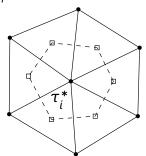
- multi-fluid support
- robust
- efficient
- automatic topology changes

Characteristic function discretization - 2

finite volume explicit method

$$\boldsymbol{\lambda}_h^{n+1} = \boldsymbol{\lambda}_h^n - \Delta t^n \boldsymbol{u}^n \cdot \nabla \boldsymbol{\lambda}_h^n$$

- dual mesh $\mathcal{T}_h^*(\Omega)$ with n_p elements
- element τ_i^* , $\bigcup_i \tau_i = \mathcal{T}_h^*$
- $\lambda_h \in V_0^*, V_0^* = \{ \varphi_h \in L^2 : \varphi_h|_{\tau_i^*} \in \mathbb{P}^0 \}$



Flux approximation

finite volume approximation on the dual grid

$$\lambda_{s}^{n+1} = (1+D^{n}) \lambda_{s}^{n} - \sum_{f} v_{f}^{n} \Phi(\widehat{\lambda}_{s,f}^{n}, \widehat{\lambda}_{s,\bar{f}}^{n})$$

$$D^{n} = \frac{\Delta t^{n}}{|\tau|} \oint_{\partial \tau} \mathbf{u} \cdot \mathbf{n} = \sum_{f} v_{f}^{n}$$

$$v_{f}^{n} = \frac{\Delta t^{n}}{|\tau|} \oint_{f} \mathbf{u} \cdot \mathbf{n}$$

$$\widehat{\lambda}_{s,f}^{n} = \lambda_{s}^{n} + \delta \lambda_{s,f}^{n}$$

$$\tau_{s}^{n} = \lambda_{s}^{n} + \delta \lambda_{s,f}^{n}$$

 $\delta\lambda^n_{s,f}$ comes from the constrined minimization problem

$$\begin{cases} \min_{\delta \lambda_{s,f}^n} \ \frac{1}{2} \sum_s \left(\lambda_s^n - \phi_{s,f}^n + \delta \lambda_{s,f}^n \right)^2 \\ \sum_s \delta \lambda_{s,f}^n = 0 \\ -\lambda_s^n < \delta \lambda_{s,f}^n < \min\left(\frac{1 + D^n - \nu_f^n |f|}{\nu_f^n |f|}, 1 - \lambda_s^n \right) \end{cases}$$

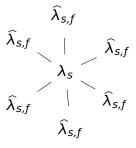
 $\delta\lambda^n_{s,f}$ as a best fit approx of LS 1

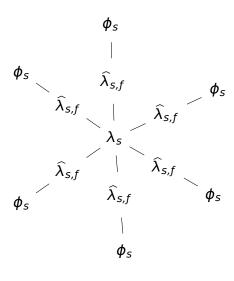
A. Cervone

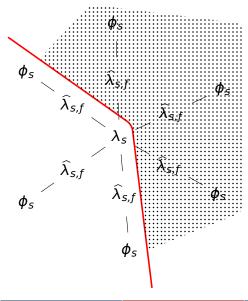
 $^{^{}m 1}$ Villa A., Formaggia L. Implicit tracking for multi-fluid simulations. JCP 229 (2010) 5788–5802

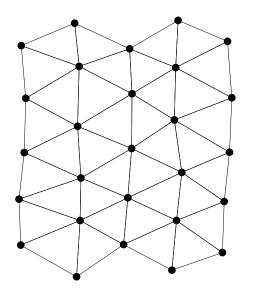


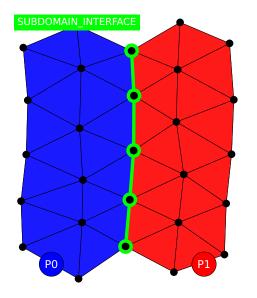
A. Cervone DD for hyperbolic eqns

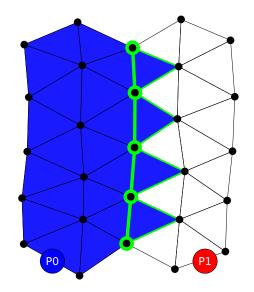


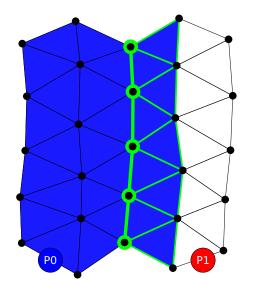


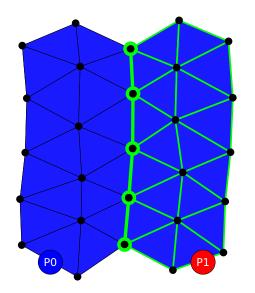












LifeV

- finite element library
- originally developed for life sciencies
- advanced FSI solvers
- heart modeling
- 1D models
- multiscale
- parallel
- based on Trilinos

LifeV maps

the MapEpetra object

- stores 2 Epetra_Maps
 - Unique: objects used to assembly
 - Repeated: objects used to share information
- handles import/exports between different maps
- operator=()
- operator+=()
- operator|=() for block support
- used to build all algebraic objects

Overlapping maps - 1

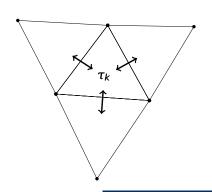
- based on connectivity
- each geometric entity knows its neighborhood

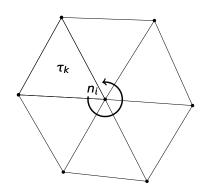




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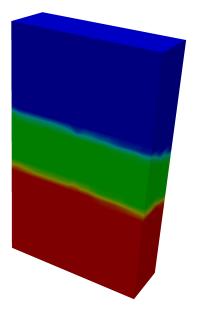


Overlapping maps - 2

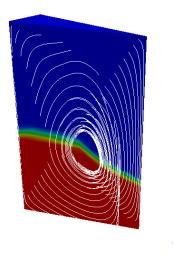
Algorithm

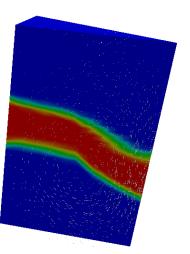
- build neighborhood information on the full mesh
- partition the mesh
- delete full mesh
- identify subdomain interface entities (check on neighbors)
- init searching set with subdomain interface entities
- for: level of overlap
 - add all neighbors not on current partition
 - replace searching set with added entities

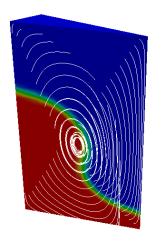
Simulation setup

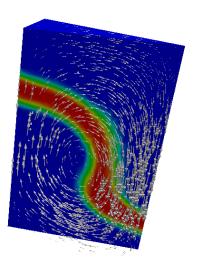


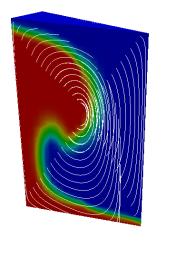
- $n_e \sim 200K$
- $n_p \sim 40K$
- dof ~ 750*K*
- $\mu = 3.0, \rho = 3.5$
- $\mu = 2.0, \rho = 0.2$
- $\mu = 0.1, \rho = 1$

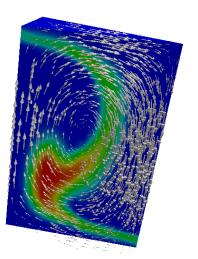


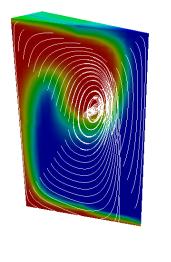


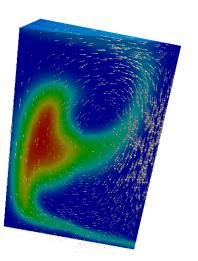


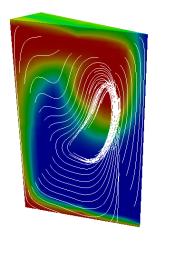


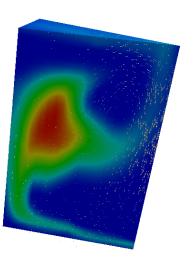


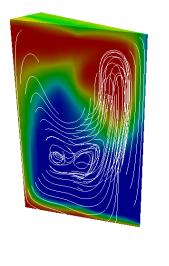


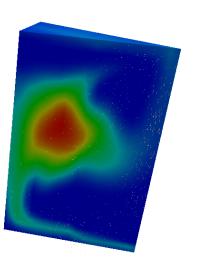










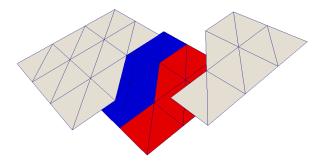


Performance analysis

- $\mu_1/\mu_2 \sim 1 \rightarrow CN = 7 \cdot 10^5$
- $\mu_1/\mu_2 \sim 10 \rightarrow CN = 2.5 \cdot 10^6$
- $\mu_1/\mu_2 \sim 100 \rightarrow CN = 2.5 \cdot 10^7$
- preconditioning computation
 - reuse?
 - Vanka smoothers
 - HYPRE
- mesh memory footpint

Assembly

- phisically replicated mesh
- resort to overlapping maps to avoid comm?
- computation intensive vs memory/comm intensive
- create graph for matrix init
- virtual subdomain interface nodes



ADR test





- 4.5M elements
- 2.2*M* points
- 4K interface elements
- 4K interface points

Performance

PROC/NODE	STANDARD	REPEATED	GAIN
2/2	11.52	11.1	3.6%
4/4	6.13	5.89	3.9%
8/8	3.11	3.01	3.2%
16/2	1.71	1.64	4.1%
16/4	1.63	1.62	0.0%
16/16	1.80	1.71	5.0%
64/8	0.47	0.42	10.6%
64/64	0.21	0.18	14.3%

thanks for the attention... go watch the poster!





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