

GMRES with Multigrid

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STOKES EQUATIONS

$$-\nabla \cdot (\nu \epsilon(\mathbf{u})) + \nabla p = \mathbf{f} \quad (1)$$

$$\nabla \cdot \mathbf{u} = 0 \quad (2)$$

→ assuming ν to be constant simplifies (1) to

$$-\nu \nabla^2 \mathbf{u} + \nabla p = \mathbf{f}$$

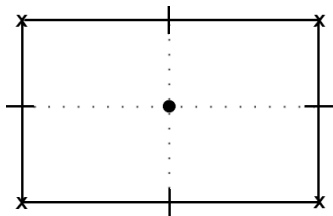
Weak form: Find $u \in H_0^1(\Omega)$ and $p \in L^2(\Omega)/\mathbb{R}$ such that

$$\begin{aligned} a(\mathbf{u}, \mathbf{v}) + b(\mathbf{v}, p) &= (\mathbf{f}, \mathbf{v}) \quad \forall \mathbf{v} \in \mathbf{H}_0^1(\Omega) \\ b(\mathbf{u}, q) &= 0 \quad \forall q \in L^2(\Omega)/\mathbb{R} \end{aligned}$$

$$\Rightarrow \begin{bmatrix} A & B \\ B^T & 0 \end{bmatrix} \begin{bmatrix} \mathbf{u} \\ p \end{bmatrix} = \begin{bmatrix} \mathbf{f} \\ 0 \end{bmatrix}$$

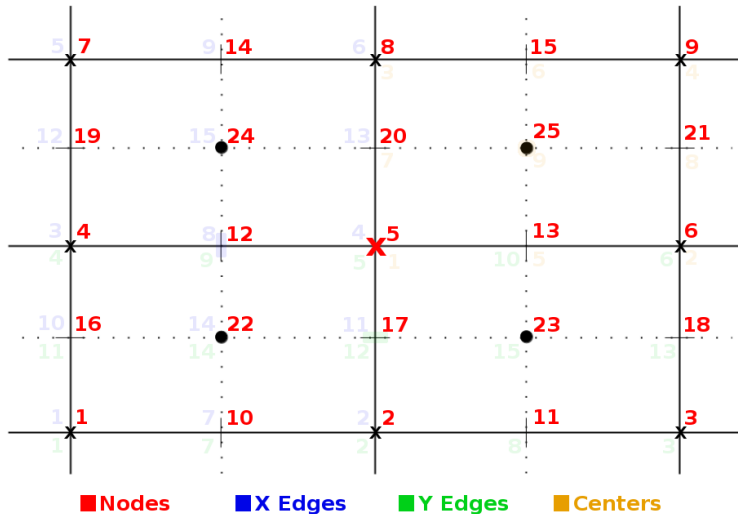
TAYLOR-HOOD ELEMENTS

Q2 elements for velocity - Q1 elements for pressure

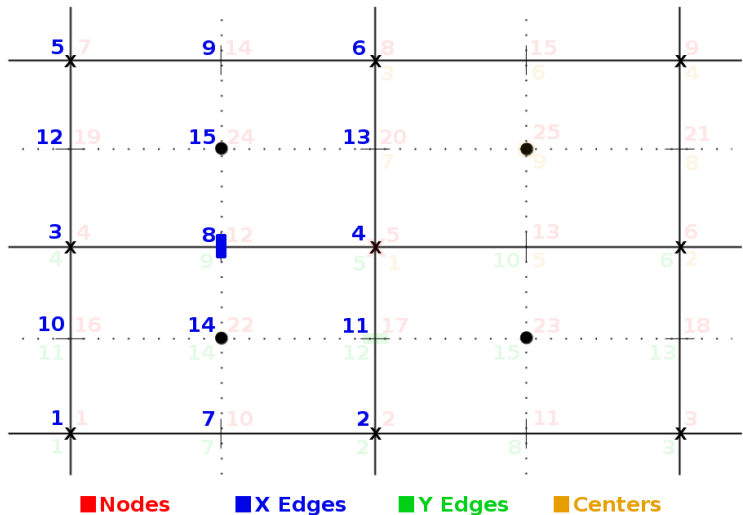


- known to be inf-sup stable
- equally spaced, regular structure

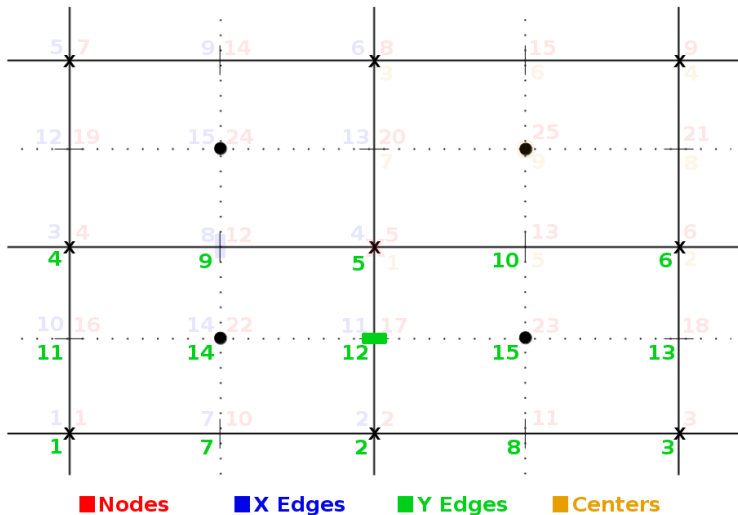
STENCIL STRUCTURE



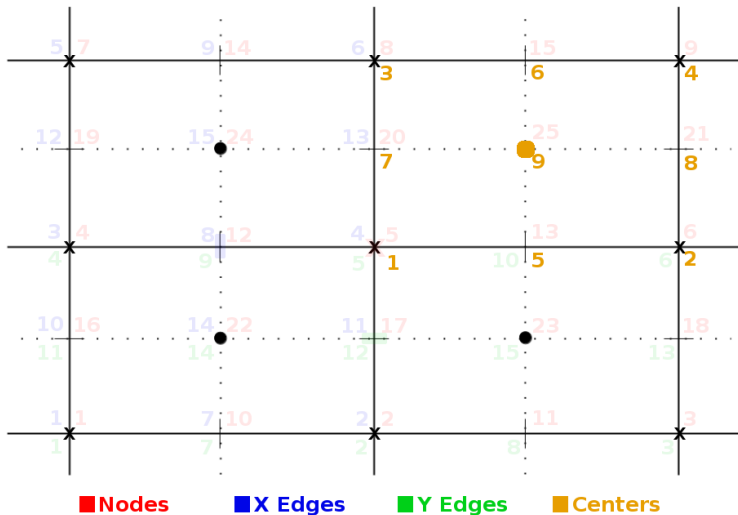
STENCIL STRUCTURE



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STENCIL EXTRACTION

Number of non-zeros per row (Laplacian):

	nodes	x-edges	y-edges	centers	Total
nodes	9	6	6	4	25
x-edges	6	3	4	2	15
y-edges	6	4	3	2	15
centers	4	2	2	1	9

→ Natural consequence: localised stencil-based calculations

→ Stencils independent from each other \Rightarrow parallelisable

GMRES

Initial system

$$A\mathbf{x} = \mathbf{b}$$

GMRES minimises 2-norm of residual

$$\|\mathbf{r}_m\|_2 = \|\mathbf{b} - A\mathbf{x}_m\|_2$$

over all vectors in the *Krylov-subspace*

$$\mathbf{x}_0 + \mathcal{K}_m(A, \mathbf{r}_0) = \mathbf{x}_0 + \text{span}\{\mathbf{r}_0, A\mathbf{r}_0, \dots, A^{m-1}\mathbf{r}_0\}.$$

(Right-)preconditioning:

$$AM^{-1}\mathbf{u} = \mathbf{b}$$

$$\mathbf{u} = M\mathbf{x}$$

→ V-Cycle with rediscratisation on coarser grids

"Ideal" Braess-Sarazin update:

$$\begin{bmatrix} \mathbf{u} \\ p \end{bmatrix}^{new} = \begin{bmatrix} \mathbf{u} \\ p \end{bmatrix}^{old} + \omega_{BS} \begin{bmatrix} tD & B \\ B^T & 0 \end{bmatrix}^{-1} \left(\begin{bmatrix} f \\ g \end{bmatrix} - A \begin{bmatrix} \mathbf{u} \\ p \end{bmatrix}^{old} \right)$$

→ factorised system

$$\begin{bmatrix} tD & 0 \\ B^T & S \end{bmatrix} \begin{bmatrix} I & \frac{1}{t}D^{-1}B \\ 0 & I \end{bmatrix} \begin{bmatrix} \delta \mathbf{u} \\ \delta p \end{bmatrix} = \begin{bmatrix} \mathbf{r}_u \\ \mathbf{r}_p \end{bmatrix}$$

with Schur complement

$$S = -\frac{1}{t}B^T D^{-1}B$$

→ compute S exactly

→ approximately solve for solution using

$$S\delta p = r_p - \frac{1}{t}B^T D^{-1}\mathbf{r}_u \quad (3)$$

$$\delta\mathbf{u} = \frac{1}{t}D^{-1}(\mathbf{r}_u - B\delta p). \quad (4)$$

with one sweep of weighted Jacobi on (3), plugging solution into (4).

→ parameter choice: $t = 1.1$, $\omega = 0.7$
(based on small experiments)

FINE-SCALE PARALLELISM

GMRES: matrix-vector and vector-vector calculations
→ easy parallelisation

Multigrid: matrix-vector and vector-vector calculations
→ easy parallelisation
interpolation/restriction
→ easy parallelisation

Braess-Sarazin: matrix-vector and vector-vector calculations
→ easy parallelisation

Solve with S : weighted Jacobi
→ easy parallelisation

Calculate $\delta \mathbf{u}$, δp : matrix-vector and vector-vector calculations
→ easy parallelisation

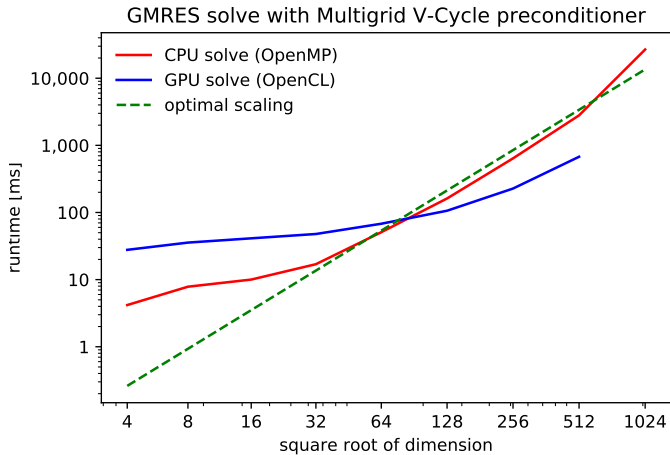
GPU CONSIDERATIONS

- ▶ fine-grain parallelism perfect for GPU's
- ▶ implemented using OpenCL
- ▶ additional cost of communication
 - moving initial data to GPU
 - moving final solution back to CPU
- ▶ Galerkin coarse-grid operators expensive
- ▶ direct solver on GPU?

REMARKS ON NUMERICAL EXPERIMENTS

- ▶ VCycle preconditioner
possible solve on coarsest grid:
 1. approximate: Braess-Sarazin smoother
 2. exact: UMFPACK library on CPU
- ▶ GPU/CPU calculations identical
(up to floating-point error)
- ▶ CPU: Intel Xeon E5; GPU: NVIDIA GeForce TITAN X

SPEEDUP COMPARISON



SUMMARY

- ▶ structured grid
 - stencil formulation
- ▶ low memory cost
- ▶ many independent calculations
 - well parallelisable
- ▶ OpenCL speeds things up by a factor of up to 4 on larger meshes
- ▶ on small meshes CPU performs better
 - overhead of moving to GPU dominant

FUTURE WORK

- ▶ CUDA version
- ▶ Vanka relaxation
Slow in serial (Braess-Sarazin better choice), potential benefits in parallel
- ▶ efficient Galerkin coarsening on GPU
- ▶ direct solve on GPU?

Thank you!