

# High-Performance Computing for Flows in Porous Media

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- ▶ Power wall
  - ▶ Power consumption is **the** limiting factor for exascale!
  - ▶ Tianhe 2: 33,8 PF (Linpack), 17.8 MW  $\Rightarrow$  500 MW
  - ▶ Clock rate stagnates but Moore's law is still valid
- ▶ Memory wall
  - ▶ Bandwidth insufficient to sustain peak performance
  - ▶ (Still) significant latency for main memory access
  - ▶ Memory access is energetically costly
  - ▶ Memory is hierarchical: register, cache, local, remote, ...
- ▶ ILP wall
  - ▶ Automatic extraction of instruction level parallelism stagnates
  - ▶ Revival of vectorization in form of SIMD instructions
- ▶ Other
  - ▶ Heterogeneous node architectures: CPU + coprocessors
  - ▶ Reduced reliability due to immense number of components?

**EXA-SCALE**  $\Rightarrow$   $10^6$  **processors**  $\times$   $10^3$  **threads**  $\times$   $10^9$  **FLOPS**

- ▶ Exploit three levels of parallelism:  
MPI + shared memory multithreading + SIMD instructions
- ▶ FLOPS are “for free”, memory access is expensive
- ▶ Matrix-based methods
  - ▶ **Robust preconditioners** available, e.g. **AMG**
  - ▶ Poor floating point performance, bound by memory bandwidth
- ▶ Matrix-free methods
  - ▶ Explicit time-stepping or certain solvers, e.g. CG, GMG
  - ▶ Low-order FE schemes:
    - ▶ Exploit problem structure: linear transformation, regular grids, constant coefficients, “stencil codes”
    - ▶ Vectorization over several elements
  - ▶ **High-order (DG) FE schemes**: naive complexity  $O(p^{\alpha d} h^{-d})$ 
    - ▶ Complexity reduction through **tensor-product approach**
    - ▶ Medium high order: Vectorization within one element
    - ▶ Very high order: Several threads to one element
- ▶ Other: mixed precision, increased asynchronicity

Porous Media Flow Problems and Their Discretization

Matrix-based Methods: Algebraic Multigrid

Matrix-free Methods: Sum Factorization for DG

More Results

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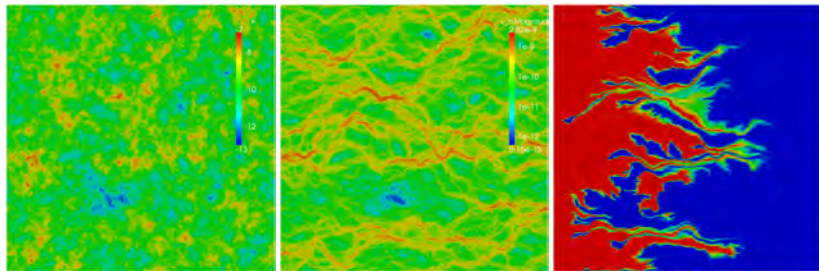
More Results

# Single Phase Flow and Transport

Parabolic/hyperbolic problem coupled to elliptic problem:

$$\nabla \cdot \mathbf{v} = f, \quad \mathbf{v} = -\frac{K}{\mu}(\nabla p - \rho \mathbf{g}),$$
$$\partial_t(\Phi c) + \nabla \cdot (c\mathbf{v} - D(\mathbf{v})\nabla c) = q.$$

Tensor  $K$ ,  $D$ , heterogeneity, convection-dominated effective macro-scale descriptions, uncertainty of parameters



Permeability  $K$

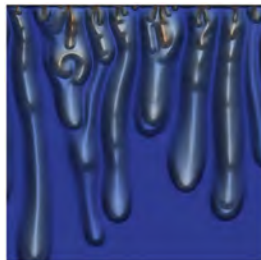
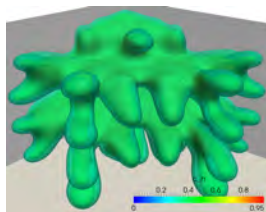
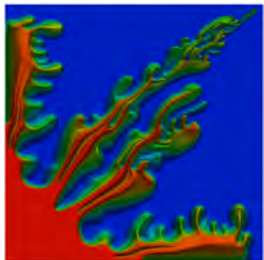
Velocity magnitude  $\|\mathbf{u}\|$

Concentration  $c$

$$\nabla \cdot \mathbf{v} = 0, \quad \mathbf{v} = -\frac{K}{\mu(c)}(\nabla p - \rho(c)g)$$

$$\partial_t(\Phi c) + \nabla \cdot (c\mathbf{v} - D(v)\nabla c) = 0$$

Density driven flow, miscible displacement  
Can exploit high resolution schemes



**Total fluid conservation:**

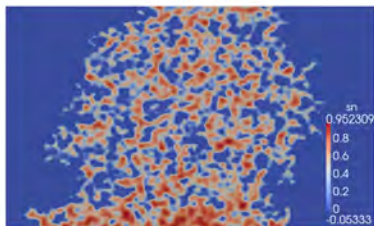
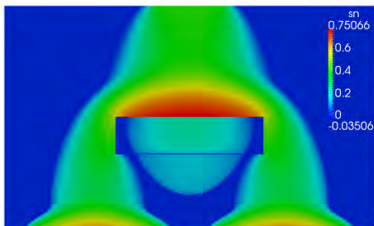
$$\nabla \cdot (v_a - \lambda_n K \nabla \phi_c) = q, \quad v_a = -\lambda_t K \nabla \phi_w.$$

**Conservation of non-wetting phase:**

$$\Phi \partial_t (1 - \psi(\phi_c)) + \nabla \cdot (f_n (1 - \psi(\phi_c)) v_a - \lambda_n (1 - \psi(\phi_c)) K \nabla \phi_c) = q_n.$$

Primary variables:  $\phi_w = p_w - \rho_w g d$ ,  $\phi_c = p_c - (\rho_n - \rho_w) g d$

$\psi(\phi_c)$ ,  $\lambda_n(\phi_c)$ ,  $\lambda_t(\phi_c)$ ,  $f_n(\phi_c)$  are nonlinear functions





## ▶ Advantages

- ▶ Potentially high order convergence
- ▶ High-order methods beneficial even with low regularity
- ▶ Element-wise conservative
- ▶ Full permeability tensors, discontinuous coefficients
- ▶ Handles elliptic, parabolic and hyperbolic equations equally well
- ▶ Unstructured grids, nonconforming refinement
- ▶ Contiguous memory access pattern
- ▶ Block matrix structure: less indirect access, BLAS level 3

## ▶ Disadvantages

- ▶ No monotonicity (for 2nd order problems)
- ▶ Free parameter to tune
- ▶ Higher number of DOFs compared to standard FE or FV
- ▶ Linear systems considered to be more difficult to solve

SIPG + weighted averages (*Di Pietro, Ern, Guermond '2008*):

$$\begin{aligned}
 a_h(p, w; c) &= \sum_{T \in \mathcal{T}_h} \int_T \left[ \frac{K}{\mu(c)} (\nabla p - \rho(c)g) \right] \cdot \nabla w - \sum_{F \in \mathcal{F}_h^{iD}} \int_F \nu_F \cdot \left\{ \frac{K}{\mu(c)} (\nabla p - \rho(c)g) \right\}_\omega \llbracket w \rrbracket \\
 &\quad - \sum_{F \in \mathcal{F}_h^{iD}} \int_F \theta \nu_F \cdot \left\{ \frac{K}{\mu(c)} \nabla w \right\}_\omega \llbracket p \rrbracket + \sum_{F \in \mathcal{F}_h^{iD}} \int_F \gamma_{F,f} \llbracket w \rrbracket \llbracket p \rrbracket \\
 l(w) &= \sum_{T \in \mathcal{T}_h} \int_T q_f w - \sum_{F \in \mathcal{F}_h^N} \int_F j_f w \quad \theta = 1 \text{ (SIPG)}, \theta = 0 \text{ (OBB)}, \theta = -1 \text{ (NIPG)} \\
 \delta_{Kn}^\pm &= \nu_F^t K^\pm \nu_F / \mu(c)^\pm, \quad \omega^- = \frac{\delta_{Kn}^+}{\delta_{Kn}^- + \delta_{Kn}^+}, \quad \omega^+ = \frac{\delta_{Kn}^-}{\delta_{Kn}^- + \delta_{Kn}^+}.
 \end{aligned}$$

Penalty term ( $\langle \cdot, \cdot \rangle$  harmonic average,  $p$  polynomial degree,  $d$  space dimension):

$$\gamma_{F,f} = \alpha \langle \delta_{Kn}^-, \delta_{Kn}^+ \rangle M, \quad M = p(p + d - 1) \frac{|F|}{\min(|T^-(F)|, |T^+(F)|)}$$

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More Results

- ▶ Wish to solve large  $Ax = b$
- ▶ Assume hierarchy of matrices

$$A_l \in \mathbb{R}^{l_l \times l_l} \quad l = 0, \dots, L, \quad A_L = A$$

$l_l = \{1, \dots, n_l\}$  denotes the index set on level  $l$

- ▶ Transfer matrices

$$R_l \in \mathbb{R}^{l_{l-1} \times l_l}, \quad P_l \in \mathbb{R}^{l_l \times l_{l-1}} \quad l = 1, \dots, L$$

- ▶ Geometric multigrid:  $A_l, R_l, P_l$  related to hierarchy of finite element spaces

$$V_l = V(\mathcal{T}_l) = \text{span}\{\varphi_{i,l} : i \in l_l\} \quad l = 0, \dots, L$$

constructed from a mesh hierarchy

$$\mathcal{T}_l \quad l = 0, \dots, L$$

- ▶ Algebraic multigrid: Only  $A$  and  $b$  are given

**Require:**  $x$  approximation of solution of linear system  $Ax = b$

**Ensure:**  $x$  improved approximation

$d_L \leftarrow b - Ax; v_L \leftarrow 0;$  ▷ Compute defect

**for**  $l \in \{L, \dots, 1\}$  **do**

Smooth( $v_l, d_l, \nu_1$ ); ▷ Pre-smoothing (updates  $v_l$ )

$d \leftarrow d_l - A_l v_l.$  ▷ Compute defect

$d_{l-1} \leftarrow R_l d;$  ▷ Restrict defect

$v_{l-1} \leftarrow 0;$  ▷ Prepare coarser level

**end for**

Solve  $A_0 v_0 = d_0;$  ▷ Exact coarsest level solve

**for**  $l \in \{1, \dots, L\}$  **do**

$v_l \leftarrow v_l + \xi P_l v_{l-1};$  ▷ Prolongate coarse grid correction

Smooth( $v_l, d_l, \nu_2$ ); ▷ Post-smoothing (updates  $v_l$ )

**end for**

$x \leftarrow x + v_L;$

## ▶ Geometric multigrid

- ▶  $\mathcal{T}_l, 0 \leq l \leq L$ : nested sequence of meshes
- ▶  $V_l = V(\mathcal{T}_l)$  nested sequence of finite element spaces
- ▶ Basis representation:  $\phi_{i,l-1} = \sum_{j \in I_l} (R_l)_{i,j} \phi_{j,l}$  gives  $R_l$
- ▶ Then  $A_{l-1} = R_l A_l P_l$  with  $P_l = R_l^T$  (Galerkin product)

## ▶ Ruge-Stüben algebraic multigrid

- ▶ Partition index set  $I_l = F_l \cup C_l$  and set  $I_{l-1} = C_l$
- ▶ Determine entries of prolongation matrix  $P_l \in \mathbb{R}^{I_l \times C_l}$
- ▶ Set  $A_{l-1} = R_l A_l P_l$  with  $R_l = P_l^T$
- ▶ Selection of  $C_l$  and interpolation factors depend on  $A_l$

## ▶ Agglomeration algebraic multigrid

- ▶ Partition index set  $I_l = \bigcup_{i=1}^{n_l-1} C_{i,l}$  (clustering)
- ▶ Partitioning depends on  $A_l$
- ▶ Set  $(R_l)_{i,j} = 1$  if  $j \in C_{i,l}$  and zero else (“piecewise constants”)
- ▶ Set  $A_{l-1} = R_l A_l P_l$  with  $P_l = R_l^T$
- ▶ Important: scale coarse grid correction by factor  $\xi$   
( $\xi \approx 2$  for regular  $2^d$  coarsening of diffusion problem)

- ▶ References
  - ▶ Developed independently in the mid 90s: *Vanek et al (1994)*, *Braess (1995)*, *Raw et al (1996)*
  - ▶ Theory and practice by Notay, e.g. *Notay, Napov, JCP, 2015*
  - ▶ Combinatorial algebraic multigrid: I. Koutis, G.L. Miller, and R. Peng. arXiv:1102.4842v4
- ▶ Our implementation follows *Raw et al (1996)*
  - ▶ Symmetric strength of connection criterion (for M-matrices): Edge  $(i, j) \in I_l \times I_l$  is called strong iff

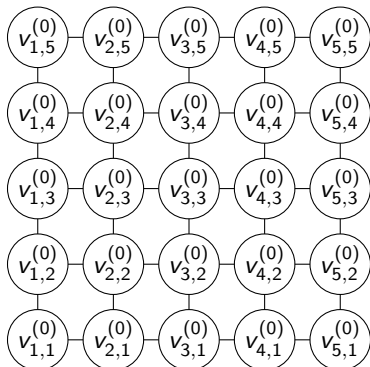
$$\frac{a_{ij}a_{ji}}{a_{ii}a_{jj}} \geq \delta \min(\eta_i, \eta_j)$$

with  $\eta_i = \max_j \frac{a_{ij}a_{ji}}{a_{ii}a_{jj}}$  and threshold  $0 < \delta < 1$ .

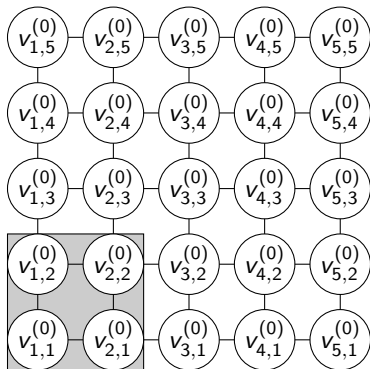
- ▶ Greedy region growing algorithm builds clusters of arbitrary size (aggressive coarsening)
- ▶ Fill-in minimization: roundness, merge small clusters
- ▶ Non-M-matrices: ignore positive off-diagonals
- ▶ Systems of PDEs: norm of blocks, guiding component

- ▶ Parallelization
  - ▶ Row-wise partitioning of the matrix with 1-overlap
  - ▶ Build clusters in parallel and exchange overlap
  - ▶ Build coarse system in parallel w/o communication
  - ▶  $P \gtrsim 5000$  requires successive coarsening
- ▶ Advantages
  - ▶ Quite robust for elliptic problems with heterogeneous coefficients
  - ▶ Sparsity pattern is very well preserved on coarser levels
  - ▶ Conjecture: AAMG V-Cycle preconditioner yields  $O(L)$  condition number
- ▶ Disadvantages
  - ▶ Needs to be used as preconditioner
  - ▶ Convergence for Poisson equation is generally not optimal (is cured by smoothed aggregation)

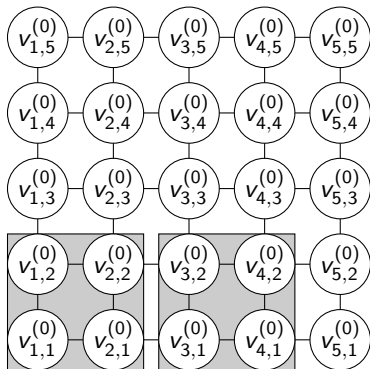




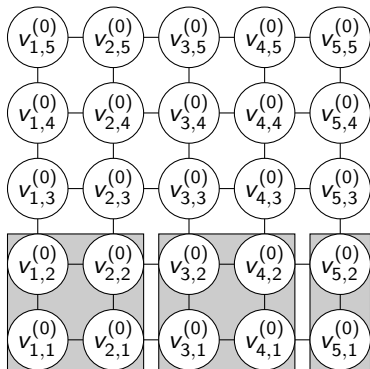
# Sequential Agglomeration



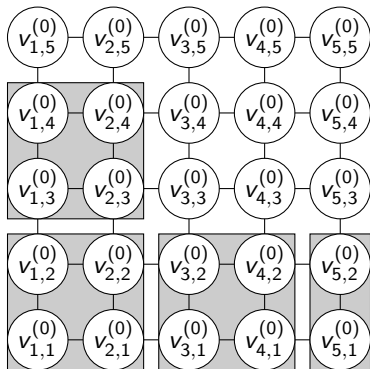
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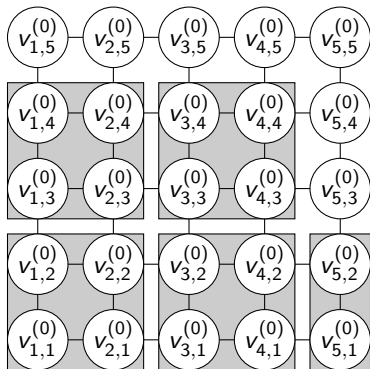
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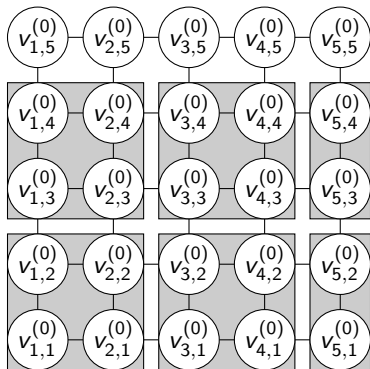
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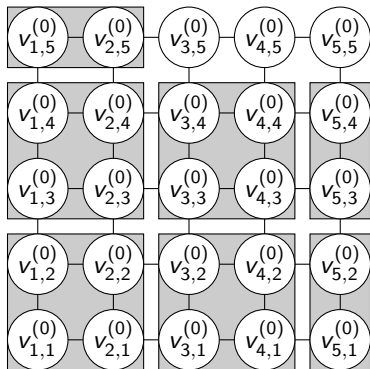
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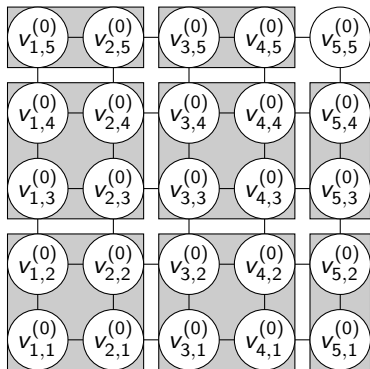


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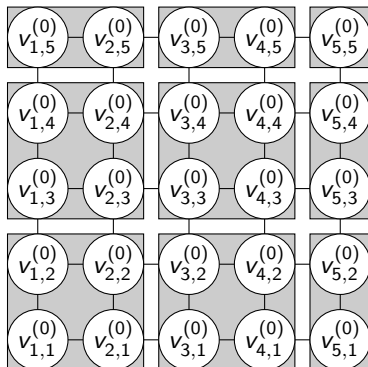




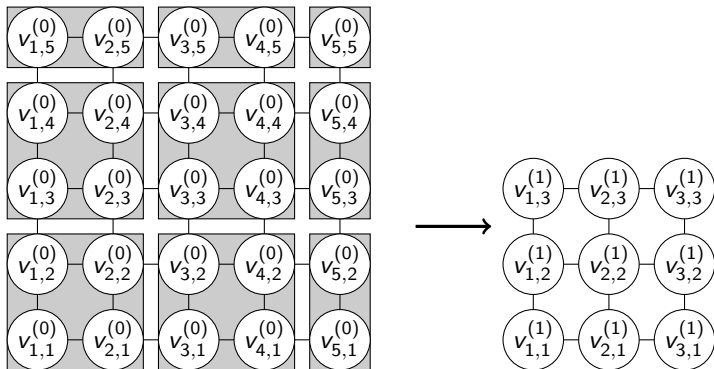
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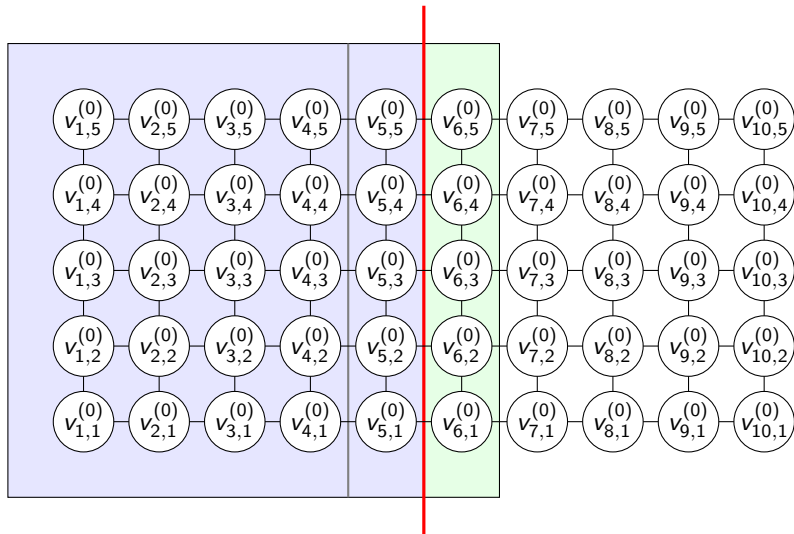


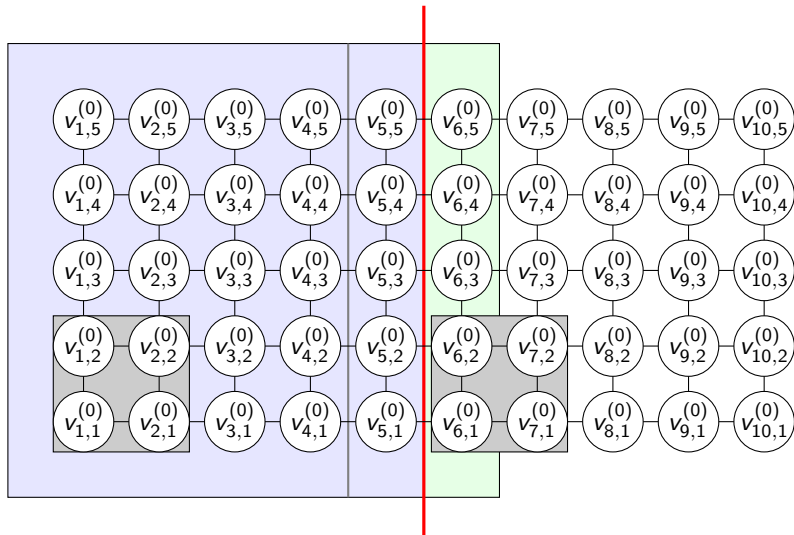
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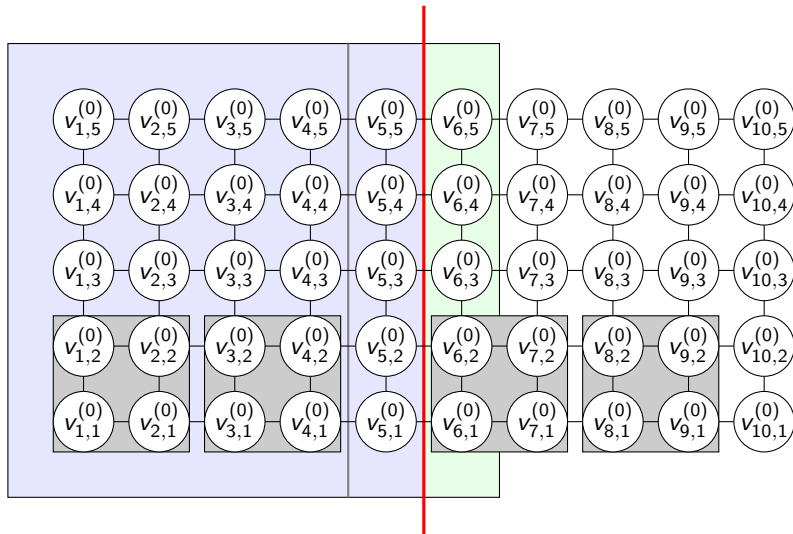


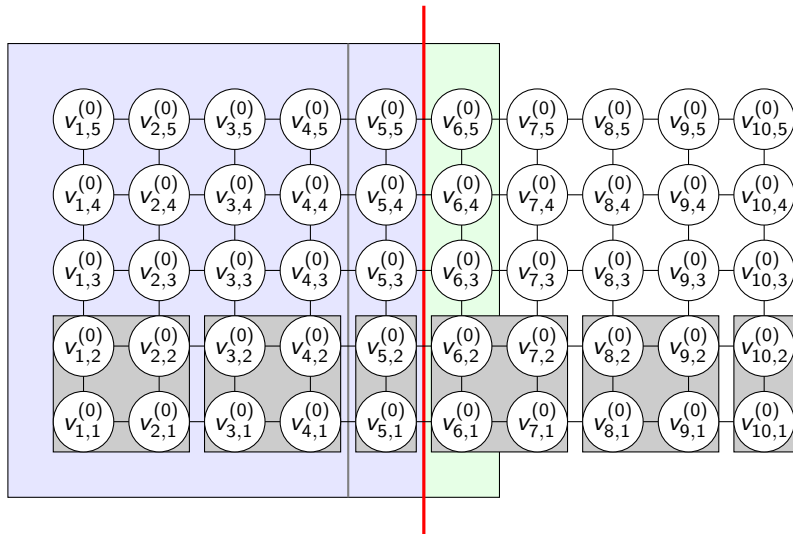
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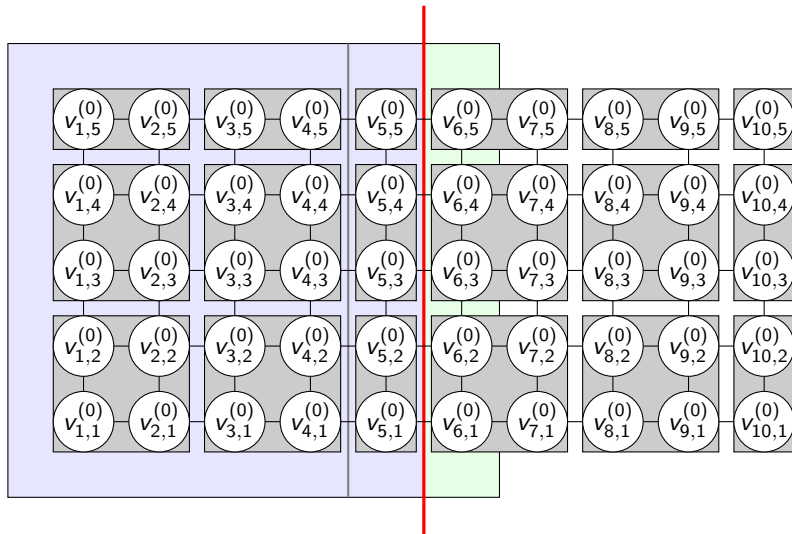




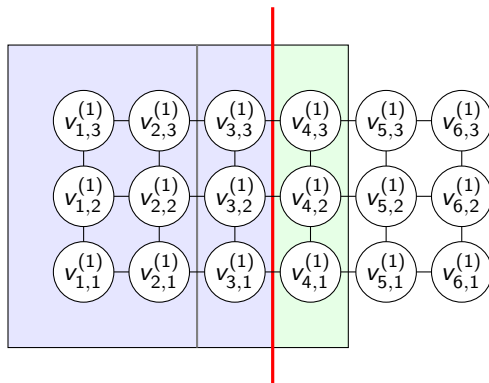












- ▶ AAMG and results presented is work of **Markus Blatt**
- ▶ **BlueGene/P** at Jülich Supercomputing Center
- ▶  $P \cdot 80^3$  degrees of freedom ( $5120^3$  finest mesh), CCFV
- ▶ **Poisson problem**,  $10^{-8}$  reduction
- ▶ AMG used as preconditioner in BiCGStab (2 V-Cycles!)
- ▶ **Scaling starts at 1**

procs	1/h	lev.	TB	TS	NIT	TIT	TT
1	80	5	19.86	31.91	8	3.989	51.77
8	160	6	27.7	46.4	10	4.64	74.2
64	320	7	74.1	49.3	10	4.93	123
512	640	8	76.91	60.2	12	5.017	137.1
4096	1280	10	81.31	64.45	13	4.958	145.8
32768	2560	11	92.75	65.55	13	5.042	158.3
262144	5120	12	188.5	67.66	13	5.205	256.2

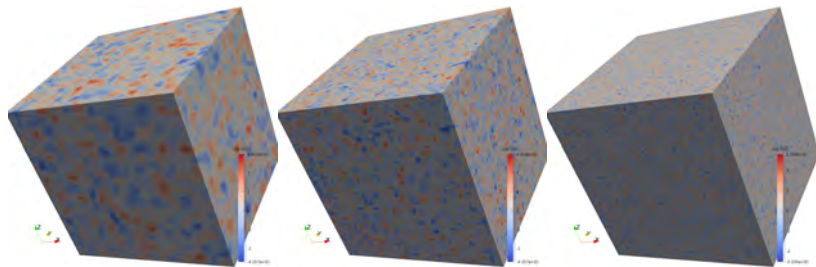
# Multiplicative AMG-DG Preconditioner

**Require:**  $x_F$  approximation of solution of DG system  $A_F x_F = b_F$

**Ensure:**  $x_F$  improved approximation

- $d_F \leftarrow b_F - A_F x_F; v_F \leftarrow 0;$  ▷ Compute defect
- Smooth( $v_F, d_F, \nu_1$ ); ▷ Pre-smoothing on DG system
- $d_C \leftarrow R_F(d_F - A_F v_F);$  ▷ Restrict defect to subspace
- $v_C \leftarrow 0;$  ▷ Prepare coarse level
- AMG( $v_C, d_C$ ); ▷ Apply AMG V-Cycle
- $v_F \leftarrow v_F + R_F^T v_C$  ▷ Add coarse grid correction
- Smooth( $v_F, d_F, \nu_2$ ); ▷ Post-smoothing on DG system
- $x_F \leftarrow x_F + v_F;$  ▷ Provide result

- ▶  $R_F$  obtained from embedding  $V_C \subset V_F = V_{DG}$
- ▶  $V_C$  may be  $Q_1$  conforming or  $P_0$
- ▶ Method is not purely algebraic
- ▶ Smoothing step takes majority of computing time
- ▶ Smoothing step may be implemented matrix-free
- ▶ See *B., Blatt, Scheichl, NLAA, 2012*

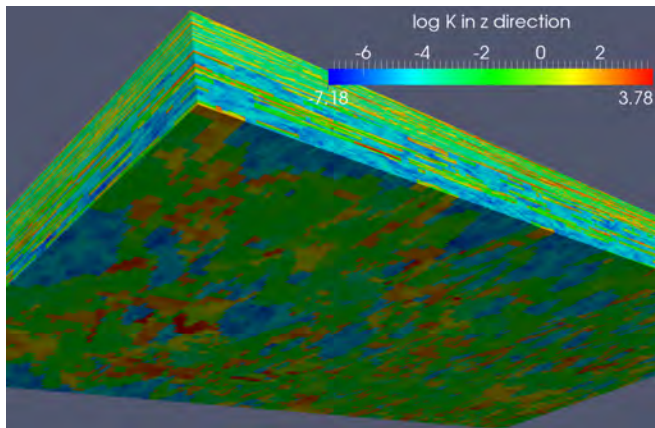


- ▶ Gaussian random field with Gaussian covariance function
- ▶ Picture: correlation length  $2h$ , variance 5
- ▶ 10 orders of magnitude variation
- ▶ Problem gets more difficult with refinement

# AMG Weak Scaling Results: Hybrid DG-AMG

Problem	$P$	TIT	NIT	TIT	NIT
		Q1		DG1/Q1	
CL=20h VAR=1	1	0.24	10	3.90	9
	8	0.32	14	4.95	12
	64	0.35	16	5.19	15
	512	0.37	17	5.24	18
	4096	0.61	19	5.88	18
		CCFV		DG1/P0	
CL=2h VAR=5	1	0.08	18	2.41	32
	8	0.12	25	3.11	49
	64	0.14	25	3.30	55
	512	0.15	40	3.36	74
	4096	0.75	50	3.92	94

- ▶  $60^3 = 216000$  cells per core,  $10^{-8}$  reduction
- ▶ Xeon CPU E5-2630 v3 @ 2.40GHz "Haswell", 8 cores
- ▶ 416 nodes, 2 CPUs per node



- ▶ Domain  $1200 \times 2200 \times 170 \text{ ft}^3$
- ▶  $60 \times 220 \times 85$  mesh, 1.1 Mio cells
- ▶ Anisotropic mesh, anisotropic diagonal permeability tensor
- ▶ 11 orders of magnitude variation

- ▶  $-\nabla \cdot (K\nabla u) = 0$ , Dirichlet on sides, Neumann top/bottom
- ▶  $10^{-6}$  reduction from zero initial value

### Strong Scaling

Scheme	What	# cores			
		1	4	16	80
CCFV	NIT	23	24	27	27
	TIT	0.27891	0.08955	0.02714	0.00907
	Overall Speedup		3.0	8.8	26.1
DG1/P0	NIT	35	38	37	42
	TIT	10.058	3.026	1.001	0.289
	Overall Speedup		3.1	9.5	29.0

### Iteration numbers for $h$ refinement, $P = 80$

# cells	$1.1 \cdot 10^6$	$8.8 \cdot 10^6$	$70.4 \cdot 10^6$
CCFV	27	33	39
DG1/P0	42	54	66

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# Exploiting Tensor Product Structure

Tensor product basis:  $\phi_i(x) = \theta_{i_d}(x_d) \dots \theta_{i_1}(x_1)$

Tensor product quadrature:  $\Xi_j = (\xi_{j_1}, \dots, \xi_{j_d})^T$

where  $i = (i_1, \dots, i_d) \in I^d$  and  $j = (j_1, \dots, j_d) \in J^d$

Evaluate finite element function on ref elem at quadrature points:

$$\begin{aligned} u_h(\Xi_i) &= \sum_{j \in J^d} c_j \phi_j(\Xi_i) && (i \in I^d) \\ &= \sum_{j_d \in J} \dots \sum_{j_1 \in J} \theta_{j_d}(\xi_{i_d}) \dots \theta_{j_1}(\xi_{i_1}) c_{j_1, \dots, j_d} && (\text{tensor product struct.}) \\ &= \sum_{j_d \in J} \dots \sum_{j_1 \in J} A_{i_d, j_d}^{(d)} \dots A_{i_1, j_1}^{(1)} c_{j_1, \dots, j_d} && (A^{(k)} \in \mathbb{R}^{m \times n}) \end{aligned}$$

Evaluation at all quadrature points is matrix-vector product:

$$y = Ax = (A^{(d)} \otimes \dots \otimes A^{(1)}) x, \quad A_{(i_1, \dots, i_d), (j_1, \dots, j_d)} = \prod_{k=1}^d A_{i_k, j_k}^{(k)}$$

Same structure obtained for BLF evaluation:  $a(u_h, \phi_i)$ ,  $i \in I^d$

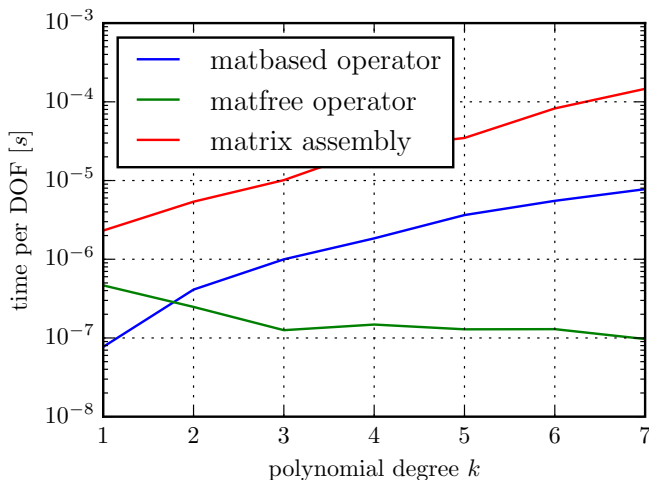
# Sum Factorization

Extract common factors:

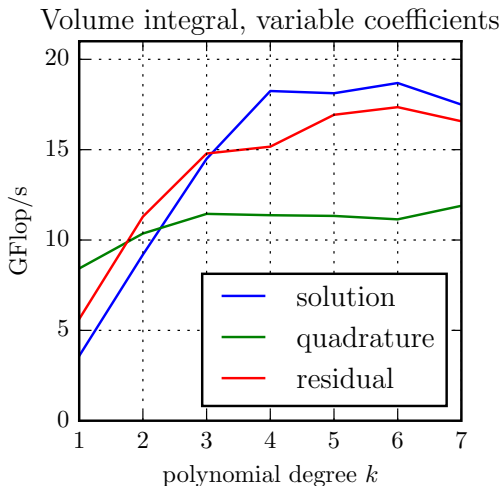
$$\begin{aligned}y_{i_1, \dots, i_d} &= \sum_{j_d \in J} \dots \sum_{j_1 \in J} A_{i_d, j_d}^{(d)} \dots A_{i_1, j_1}^{(1)} x_{j_1, \dots, j_d}^{(0)} \\ &= \sum_{j_d \in J} \dots \sum_{j_2 \in J} A_{i_d, j_d}^{(d)} \dots A_{i_2, j_2}^{(2)} \underbrace{\left( \sum_{j_1 \in J} A_{i_1, j_1}^{(1)} x_{j_1, \dots, j_d}^{(0)} \right)}_{x_{i_1, j_2, \dots, j_d}^{(1)}} \\ &= \sum_{j_d \in J} \dots \sum_{j_3 \in J} A_{i_d, j_d}^{(d)} \dots A_{i_3, j_3}^{(3)} \underbrace{\left( \sum_{j_2 \in J} A_{i_2, j_2}^{(2)} x_{i_1, j_2, \dots, j_d}^{(1)} \right)}_{x_{i_2, i_1, j_3, \dots, j_d}^{(2)}} \\ &= \dots = \sum_{j_d \in J} A_{i_d, j_d}^{(d)} x_{i_{d-1}, \dots, i_1, j_d}^{(d-1)}\end{aligned}$$

Do this for all  $(i_1, \dots, i_d)$  simultaneously

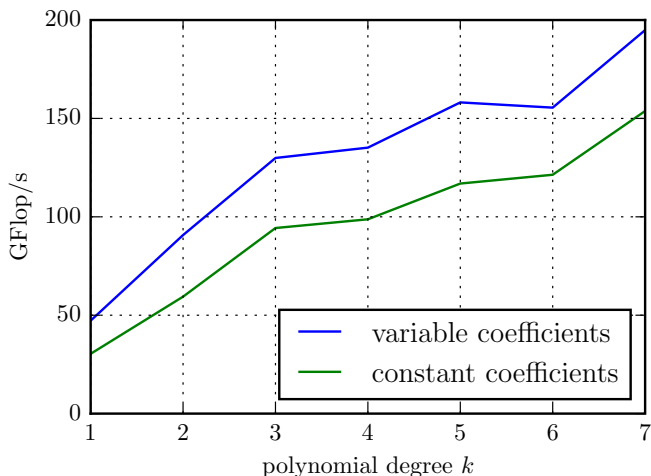
- ▶ Sum factorization algorithm by *Buis, Dyksen (1996)*:
- ▶ For  $k = 1, \dots, d$  do („dim by dim approach”)
  - ▶ compute **matrix-matrix product**  $X^{(k)} = A^{(k)}X^{(k-1)}$
  - ▶ “rotate”  $X^{(k)}$  to make  $j_{k+1}$  the row index (transposition in 2d)
- ▶ Complexity reduction ( $n = m = p + 1$ ):
  - ▶  $O(p^{2d}) \rightarrow O(p^{d+1})$  for operator application (mat-vec)
  - ▶  $O(p^{3d}) \rightarrow O(p^{2d+1})$  for Jacobian assembly
- ▶ Small matrices, e.g.  $(4 \times 4) \cdot (4 \times 16)$  for  $\mathbb{Q}_3$  in 3d
- ▶ FE operator evaluation, by e.g. *Melenk, Gerdes, Schwab (2001)*, *Kronbichler, Korman (2012)*:
  1. compute  $\hat{u}_h, \nabla \hat{u}_h$  at quadrature points,  $O(p^{d+1})$
  2. compute coefficients, geometry factors at quad. points,  $O(p^d)$
  3. compute residuals  $O(p^{d+1})$
- ▶ Applicable to nonlinear problems
- ▶ Applicable to non-affine transformations
- ▶ In DG face terms are dominating:  $O(p^d)$  complexity observed



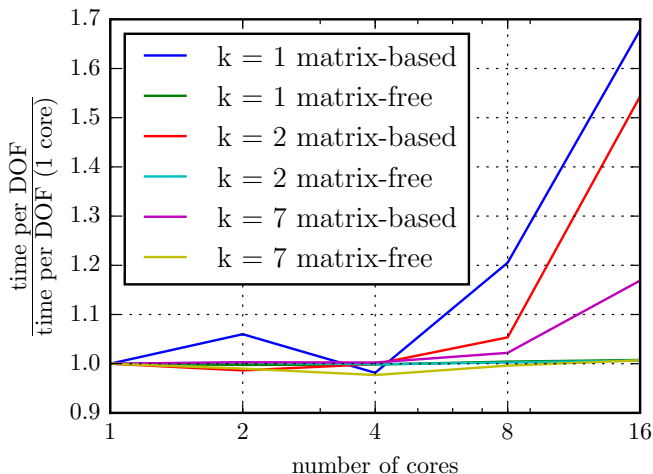
- ▶ **All results by Steffen Müthing**
- ▶ SIPG for 3d stationary convection-diffusion operator
- ▶ Matrix-based uses blocked SELL-C- $\sigma$  matrix format
- ▶ Matrix-assembly is sum-factorized!



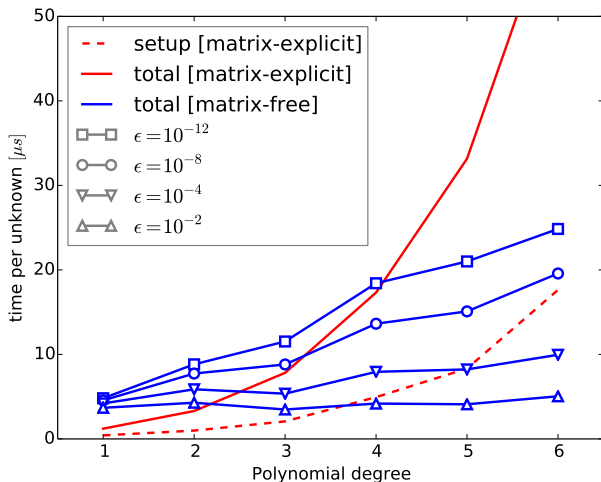
- ▶ SIPG for 3d stationary convection-diffusion operator
- ▶ Xeon E5-2698v3 (Haswell), Peak  $\approx 40$  GFlop/s per core
- ▶ Variable diffusion tensor and velocity field



- ▶ SIPG for 3d stationary convection-diffusion operator
- ▶ Xeon E5-2698v3 16 cores (Haswell), Peak  $\approx 650$  GFlop/s
- ▶ Variable diffusion tensor and velocity field



- ▶ Weak scaling: Constant #DOFs per core
- ▶ Xeon E5-2698v3 16 cores (Haswell)
- ▶ Matrix-based uses blocked SELL-C- $\sigma$  matrix format



- ▶ All results by Eike Müller
- ▶ SIPG for 3d stationary convection-diffusion equation



Porous Media Flow Problems and Their Discretization

Matrix-based Methods: Algebraic Multigrid

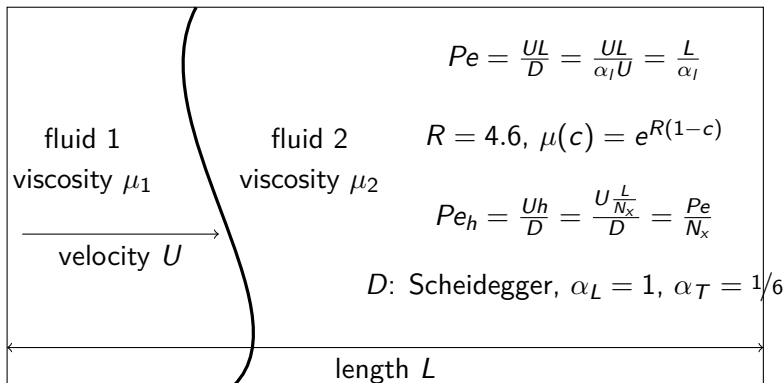
Matrix-free Methods: Sum Factorization for DG

More Results

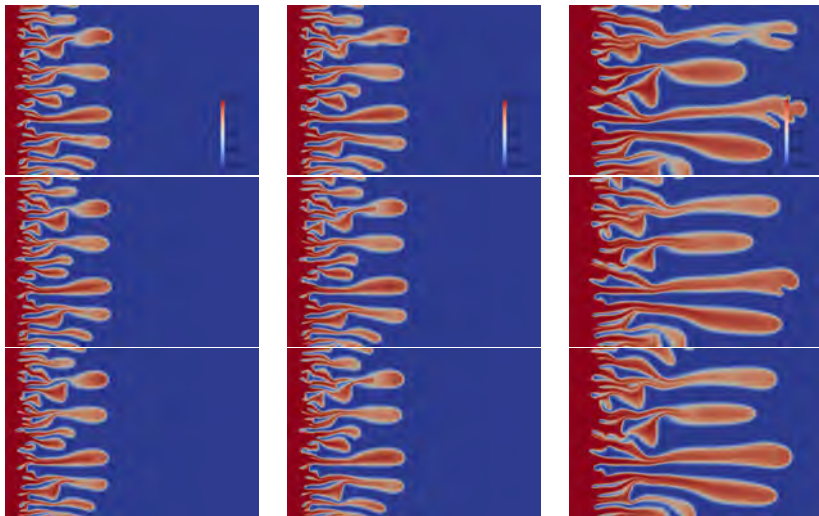
$$\nabla \cdot \mathbf{v} = 0, \quad \mathbf{v} = -\frac{K}{\mu(c)}(\nabla p - \rho \mathbf{g})$$

$$\partial_t(\Phi c) + \nabla \cdot (c\mathbf{v} - D(\mathbf{v})\nabla c) = 0$$

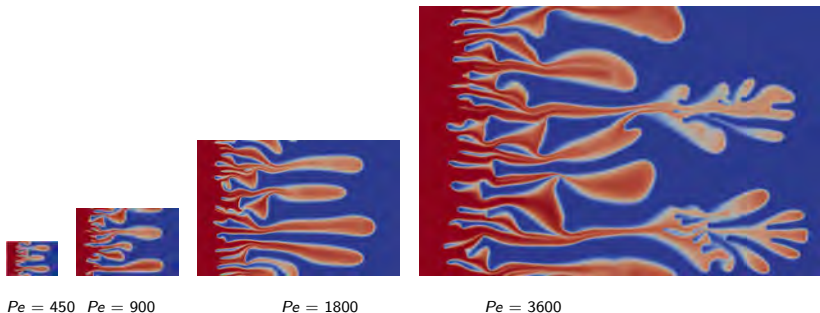
- ▶ Application e.g. in polymer flooding
- ▶ Unstable for less viscous fluid displacing more viscous fluid
- ▶ Decoupled implicit pressure / implicit concentration algorithm



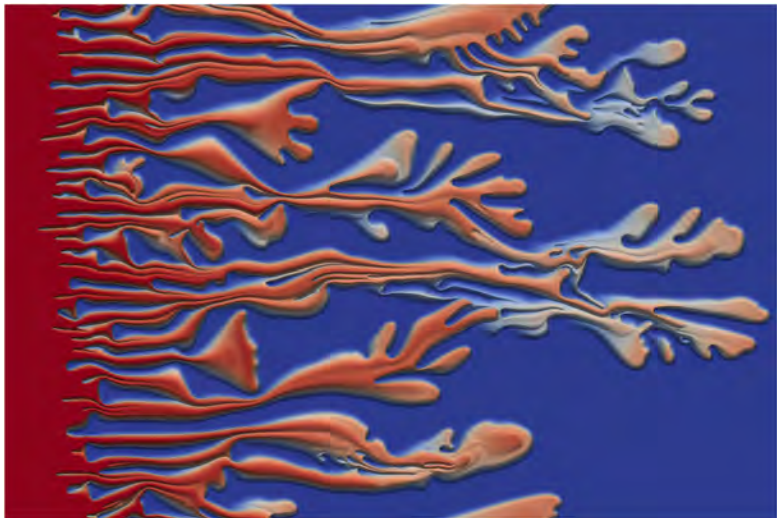
- ▶ Unstable flow for  $R > 0$
- ▶ Need  $Pe_h \approx 1$  to avoid unphysical oscillations
- ▶ Aspect ratio  $A = 3 : 2$ , periodic in  $y$



Top to bottom:  $Q_1/Q_1, Pe_h = 2$ ,  $Q_1/Q_1, Pe_h = 1$ ,  $Q_2/Q_2, Pe_h = 2$   
(All DG methods)

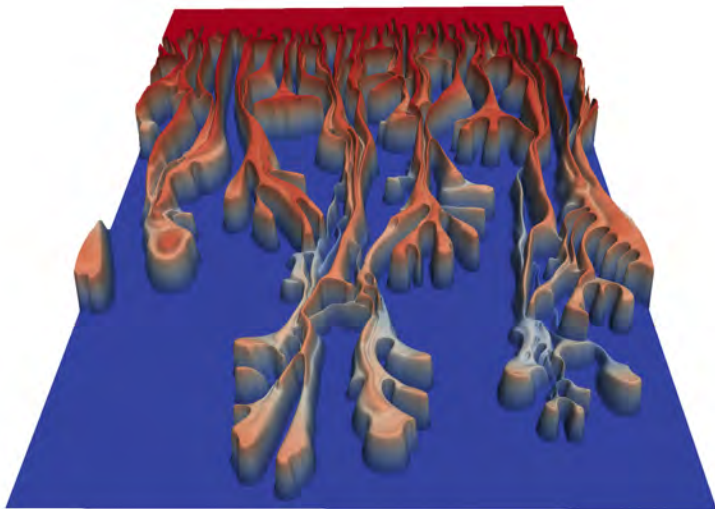


- ▶ Increasing  $Pe$  corresponds to increasing  $L$  and time
- ▶ Initially, fingers compete and merge, leading to spikes running on top of existing fingers
- ▶ Tip splitting starts at  $\approx Pe = 2000$
- ▶ Interaction of splitted fingers becomes increasingly complex



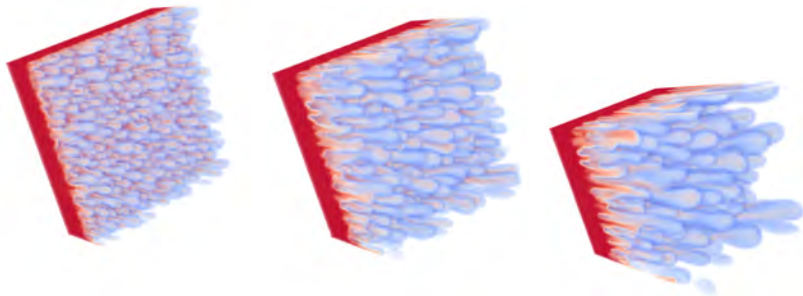
128 × Xeon E5-2630 v3 8 core,  $155 \cdot 10^6$  DOF, Alexander-2, 17000 time steps, 32h

# Pe = 7200 Warped View



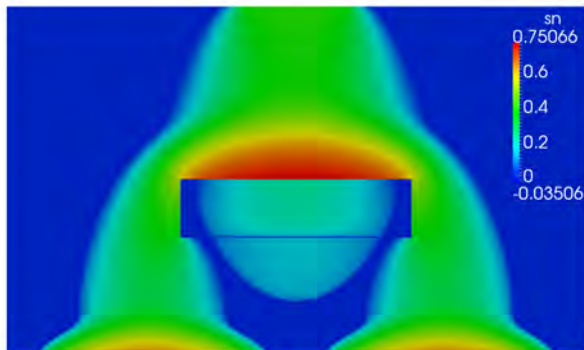
128 × Xeon E5-2630 v3 8 core,  $155 \cdot 10^6$  DOF, Alexander-2, 17000 time steps, 32h

# First 3d Results for $Pe = 300$



8 × Xeon E5-2680 v3 10 core,  $243 \cdot 10^6$  DOF, Alexander-2



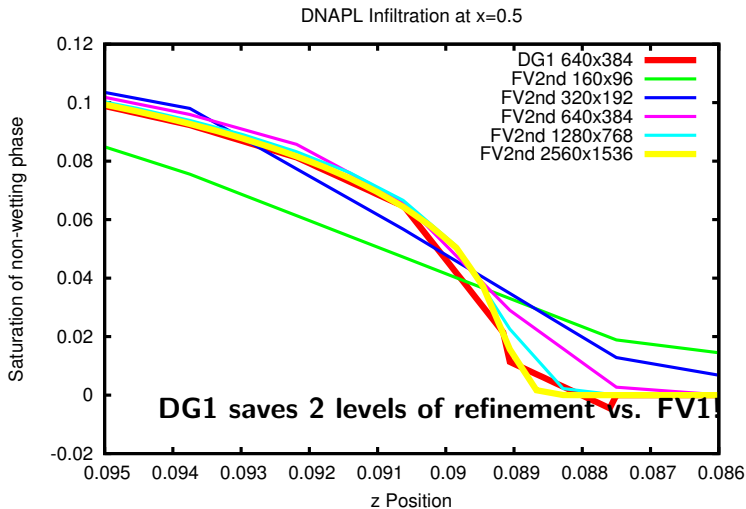


- ▶  $k_{r\alpha}$ : quadratic
- ▶  $p_c$ : Brooks-Corey,  $\lambda = 2$   
 $p_e^{bulk} = 755$ ,  
 $p_e^{lense} = 1163$

- ▶ Fully implicit fully coupled
- ▶ DG1 / Alexander(2)
- ▶ CCFV0: full upw., impl. Euler
- ▶ CCFV1: central, Alex.(2)

# At Free Boundary

DG1 640x384 vs. FV1



# Weak Scalability for DNAPL Infiltration

$T = [0, 3600s]$ , DG1/Alexander(2) vs. FV1/Alexander(2)

	$P$	$h^{-1}$	DOF	TS	ANL	ALIN	TT	LTIT
DG1	1	40	$7.7 \cdot 10^3$	30	11.4	1.8	156.8	0.144
	4	80	$3.1 \cdot 10^4$	71	11.0	3.4	466.1	0.052
	16	160	$1.2 \cdot 10^5$	125	12.0	5.7	1187.3	0.062
	<b>64</b>	<b>320</b>	<b><math>4.9 \cdot 10^5</math></b>	<b>263</b>	<b>11.4</b>	<b>9.1</b>	<b>3250.5</b>	<b>0.067</b>
	256	640	$2.0 \cdot 10^6$	563	10.7	15.3	11233.0	0.082
	1024	1280	$7.9 \cdot 10^6$	1369	9.2	17.0	41917.4	0.112
FV1	1	160	$3.1 \cdot 10^4$	60	7.7	1.8	191.2	0.043
	4	320	$1.2 \cdot 10^5$	120	8.2	2.2	592.1	0.069
	16	640	$4.9 \cdot 10^5$	246	8.4	2.6	1453.7	0.078
	<b>64</b>	<b>1280</b>	<b><math>2.0 \cdot 10^6</math></b>	<b>491</b>	<b>8.9</b>	<b>3.3</b>	<b>6496.1</b>	<b>0.151</b>
	256	2560	$7.9 \cdot 10^6$	1021	9.7	4.5	12774.9	0.156

**Accuracy(DG1) = Accuracy(FV1) on 2 times refined mesh**

- ▶ Agglomeration Algebraic Multigrid is reasonably robust and scalable
- ▶ Sum factorization has reduced complexity at increased performance
- ▶ Hybrid AMG-DG preconditioner can overcome memory wall and improve scalability
- ▶ All methods presented in this talk have been implemented in the DUNE software framework ([www.dune-project.org](http://www.dune-project.org)) as part of the **EXA-DUNE** project
- ▶ DUNE is also the basis of the *Open Porous Media Initiative* driven by StatOil (<http://opm-project.org>)

**Thank you for your attention!**