

A Variational Approach to Upscaling Heterogeneous Media

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Heterogeneous Media

What makes a medium *heterogeneous*?

- Large relative variation in material properties
- Abrupt changes in material properties
- Large variation in spatial scales

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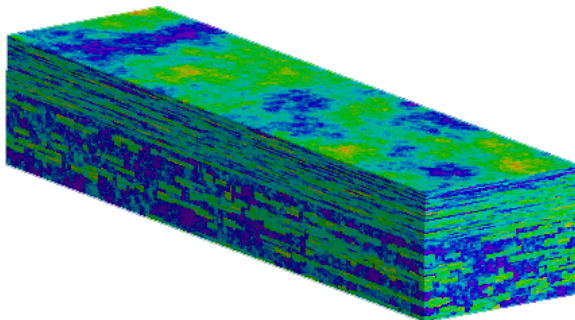
- Large relative variation in material properties
- Abrupt changes in material properties
- Large variation in spatial scales

Why do we care?

- Many natural media are heterogeneous
- Fine-scale variation affects macroscopic behavior
- Simulation of heterogeneous media must resolve variation

Subsurface Flow

Rate of flow through a reservoir depends on its composition



- Porosity & Permeability vary on scales from mm upwards
- Domain is $\sim 100\text{m} \times 50\text{m} \times 10\text{m}$

Darcy's Law

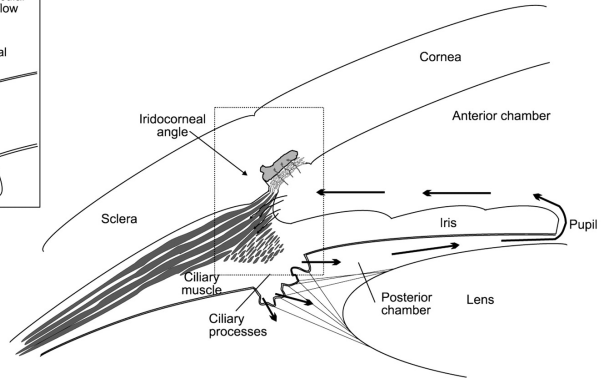
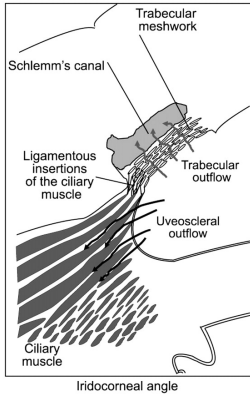
Model hydraulic head, h , of a fluid confined in a porous media

$$Q = -\mathcal{K}\nabla h$$

$$S_s \frac{\partial h}{\partial t} + \nabla \cdot Q = q$$

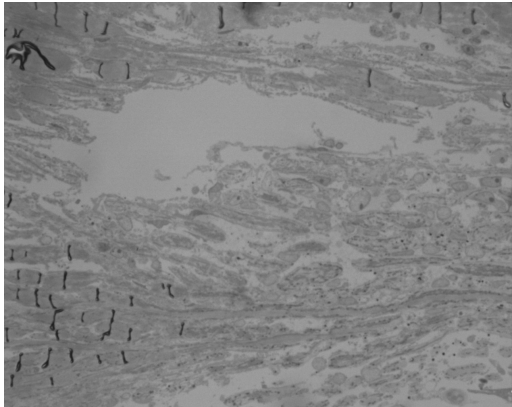
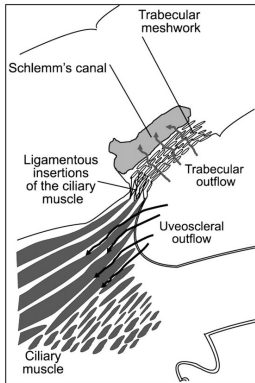
- Q denotes the Darcy-law flux
- q represents external sources or sinks of fluid
- Material properties
 - ▶ S_s = specific storage
 - ▶ \mathcal{K} = hydraulic conductivity

Ocular Flow



A. Llobet et al, *News Physiol. Sci.* **18**, pp 205-209, 2003.

Trabecular Meshwork



(left) A. Llobet et al, *News Physiol. Sci.* **18**, pp 205-209, 2003.

(right) Courtesy W.D. Stamer, U of Arizona & J.J. Heys, Arizona State U

Cardiac Bidomain Equations

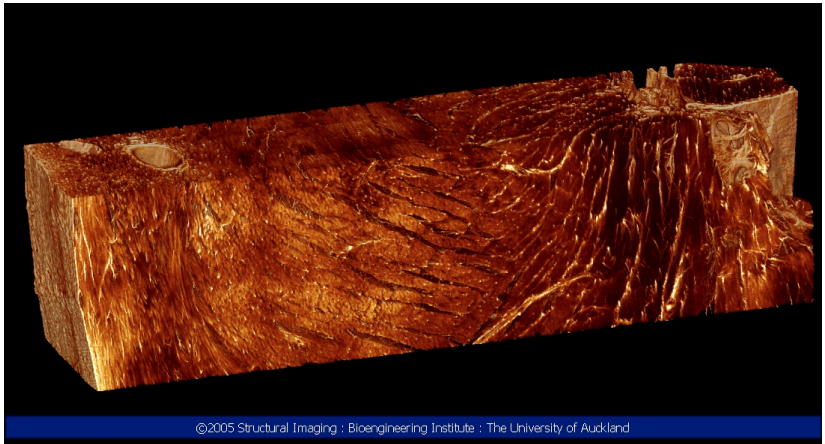
Model intra- and extra-cellular potentials, ϕ_i and ϕ_e , in cardiac tissue:

$$V_m = \phi_i - \phi_e$$

$$A_m C_m \frac{\partial V_m}{\partial t} - \nabla \cdot (\sigma_i \nabla V_m) = \nabla \cdot (\sigma_i \nabla \phi_e) - A_m I_{ion} \\ - \nabla \cdot ((\sigma_i + \sigma_e) \nabla \phi_e) = \nabla \cdot (\sigma_i \nabla V_m) + i_e(t)$$

- A_m is surface-to-volume ratio of the cell membrane
- C_m is the membrane capacitance per unit area
- I_{ion} represents ionic currents
- $i_e(t)$ represents **extracellular current injections**
- Material properties
 - ▶ σ_i = intracellular conductivity
 - ▶ σ_e = extracellular conductivity

Cardiac Tissue



Sample of rat left ventricular wall, dimensions are approximately $3.6 \times 0.8 \times 0.8$ mm.

M. Trew, B. Smail, and A. Pullan, preprint 3/7/2005.

Elliptic Model Problem

A simpler model still displays same sensitivity to heterogeneity:

$$-\nabla \cdot (\mathcal{K} \nabla h) = q$$

- Implicit time stepping adds lower-order term
- Main terms in operator-splitting approach
- Assume $\mathcal{K} = \mathcal{K}(\mathbf{x})$, possibly tensor-valued

Develop approach for model problem, then extend to particular applications

Simulation Challenges

Even for model problem, simulation can be difficult

- If $\mathcal{K}(\mathbf{x})$ varies on a fine-enough scale, simulation may be intractable

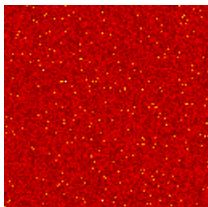
Example: 1 km \times 1 km \times 1 km reservoir, sediment varies on mm-scale requires 10^{18} Degrees of Freedom

Two approaches:

- Average conductivity to scale where simulation is possible
- Take variation in $\mathcal{K}(\mathbf{x})$ into account in discretization

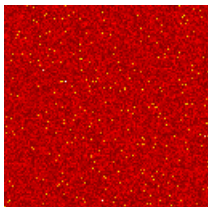
Effective Media

Given heterogeneous conductivity in a region, can we replace it by a homogeneous one without changing overall flow?



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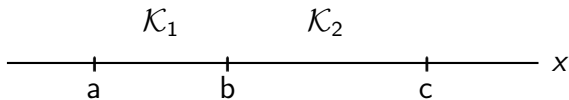


In general,

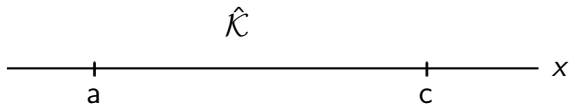
- depends on medium
- depends on flow conditions
- no single average always works

Effective Conductivity in One Dimension

Is it possible to replace a heterogeneous cell,

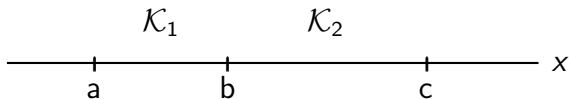


with an effective (*homogenized, or equivalent*) cell,

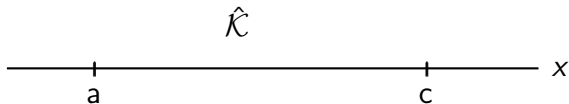


Effective Conductivity in One Dimension

Is it possible to replace a heterogeneous cell,



with an effective (*homogenized, or equivalent*) cell,



that doesn't perturb the solution outside $a < x < c$?

$$\hat{h}(a) = h(a),$$

$$\hat{h}(c) = h(c)$$

$$\hat{Q}(a) = Q(a),$$

$$\hat{Q}(c) = Q(c)$$

Harmonic Averages

One-dimensional model problem:

$$-\frac{\partial}{\partial x} \mathcal{K} \frac{\partial}{\partial x} h(x) = 0.$$

For constant \mathcal{K}_1 on $[a, b]$, integrating in x gives

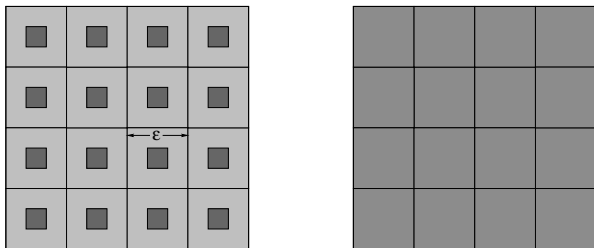
$$\begin{bmatrix} h(b) \\ Q(b) \end{bmatrix} = \begin{bmatrix} 1 & -\frac{b-a}{\mathcal{K}_1} \\ 0 & 1 \end{bmatrix} \begin{bmatrix} h(a) \\ Q(a) \end{bmatrix} = M_a^b \begin{bmatrix} h(a) \\ Q(a) \end{bmatrix}.$$

For a heterogeneous media, then

$$\begin{bmatrix} h(c) \\ Q(c) \end{bmatrix} = M_b^c M_a^b \begin{bmatrix} h(a) \\ Q(a) \end{bmatrix} = \hat{M}_a^c \begin{bmatrix} h(a) \\ Q(a) \end{bmatrix}$$

If $\hat{M}_a^c = M_b^c M_a^b$, then $\hat{\mathcal{K}} = (c - a) \left(\frac{b-a}{\mathcal{K}_1} + \frac{c-b}{\mathcal{K}_2} \right)^{-1}$.

Asymptotic Analysis



Let $\mathcal{K} = \mathcal{K}\left(\frac{\mathbf{x}}{\epsilon}\right)$, and consider

$$-\nabla \cdot \left(\mathcal{K} \left(\frac{\mathbf{x}}{\epsilon} \right) \nabla h_\epsilon \right) = q(\mathbf{x}).$$

A two-scale asymptotic analysis gives behavior as $\epsilon \rightarrow 0$.

Homogenization

Effective conductivity depends on unit cell, Y , relative to $\frac{\mathbf{x}}{\varepsilon}$.

Define

$$a_\varepsilon(u, v) = \int_Y \left(\mathcal{K} \left(\frac{\mathbf{x}}{\varepsilon} \right) \nabla_\varepsilon u \right) \cdot \nabla_\varepsilon v,$$

then

$$\xi^T \hat{\mathcal{K}} \xi^T = \min_{\phi \in H_p^1(Y)} a_\varepsilon(p_\xi + \nabla \phi, p_\xi + \nabla \phi),$$

where

- $\xi = \nabla p_\xi$ is constant
- $H_p^1(Y)$ is the Sobolev space, $H^1(Y)$, with periodic boundary conditions

Explicit Averages

- In one dimension, answer was harmonic average
- In d dimensions, theory limited to periodic media

How bad are simple, explicit averages at approximating effective conductivities?

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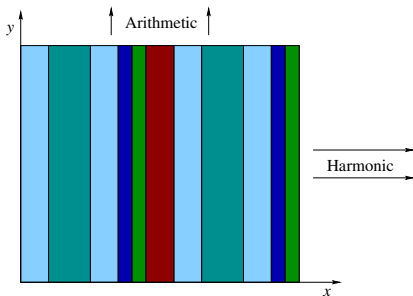
Arbitrarily.

Explicit Averages

- In one dimension, answer was harmonic average
- In d dimensions, theory limited to periodic media

How bad are simple, explicit averages at approximating effective conductivities?

Arbitrarily. Depending on flow conditions:



Weak Forms

Consider solution of

$$-\nabla \cdot \mathcal{K}(\mathbf{x}) \nabla p(\mathbf{x}) = q(\mathbf{x})$$

Weak Forms

Consider solution of

$$(-\nabla \cdot \mathcal{K}(\mathbf{x})\nabla p(\mathbf{x})) \varphi(\mathbf{x}) = q(\mathbf{x})\varphi(\mathbf{x})$$

Weak Forms

Consider solution of

$$\int_{\Omega} (-\nabla \cdot \mathcal{K}(\mathbf{x}) \nabla p(\mathbf{x})) \varphi(\mathbf{x}) = \int_{\Omega} q(\mathbf{x}) \varphi(\mathbf{x})$$

Weak Forms

Consider solution of

$$\int_{\Omega} (\mathcal{K}(\mathbf{x}) \nabla p(\mathbf{x})) \cdot \nabla \varphi(\mathbf{x}) = \int_{\Omega} q(\mathbf{x}) \varphi(\mathbf{x}) + \text{BCs}$$

Weak Forms

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$$\int_{\Omega} (\mathcal{K}(\mathbf{x}) \nabla p(\mathbf{x})) \cdot \nabla \varphi(\mathbf{x}) = \int_{\Omega} q(\mathbf{x}) \varphi(\mathbf{x}) + \text{BCs}$$

Define

$$a(u, v) = \int_{\Omega} (\mathcal{K}(\mathbf{x}) \nabla u(\mathbf{x})) \cdot \nabla v(\mathbf{x})$$

Properties of $a(u, v)$:

- Defined for u (and v) such that $\int_{\Omega} \nabla u \cdot \nabla u < \infty$
- Positive Definite: $a(u, u) > 0$ for $u \neq 0$
- Symmetric: $a(u, v) = a(v, u)$,

Weak form defines an inner product and a norm on $H^1(\Omega)$

Subspace Minimization

Let p be the solution of

$$a(p, \varphi) = \int_{\Omega} q(\mathbf{x})\varphi(\mathbf{x}) + \text{BCs for all } \varphi \in H^1(\Omega).$$

Given a subspace, $\mathcal{V} \subset H^1(\Omega)$, best solution in \mathcal{V} is

$$p_{\mathcal{V}} = \underset{v \in \mathcal{V}}{\operatorname{argmin}} a(p - v, p - v)$$

Minimizer must satisfy

$$a(p_{\mathcal{V}}, \varphi) = \int_{\Omega} q(\mathbf{x})\varphi(\mathbf{x}) + \text{BCs for all } \varphi \in \mathcal{V}$$

Basis Functions

Suppose $\mathcal{V} = \text{span}\{\phi_j(\mathbf{x})\}_{j=1}^n$, then $p_{\mathcal{V}}(\mathbf{x}) = \sum_{j=1}^n p_j \phi_j(\mathbf{x})$.
Then,

$$\sum_{j=1}^n p_j a(\phi_j, \phi_i) = \int_{\Omega} q(\mathbf{x}) \phi_i(\mathbf{x}) + \text{BCs}_i = q_i \text{ for all } i.$$

Writing $\mathbf{p} = (p_1, p_2, \dots, p_n)^T$ and $\mathbf{q} = (q_1, q_2, \dots, q_n)^T$, then

$$A\mathbf{p} = \mathbf{q},$$

where $A_{ij} = a(\phi_j, \phi_i)$.

Classical Finite Elements

Want to choose basis, $\{\phi_j\}_{j=1}^n$, so that

- $p_{\mathcal{V}}$ is a good approximation to p
- A and \mathbf{q} are easy to calculate
- $A\mathbf{p} = \mathbf{q}$ is easy to solve

Classical Finite Elements

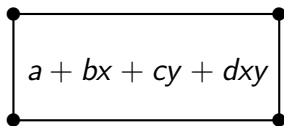
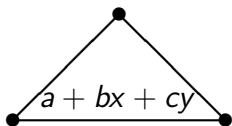
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Typical choices:

- Piecewise linears on triangles and tetrahedra
- Piecewise bilinears on quadrilaterals
- Piecewise trilinears on hexahedra

Local bases on polyhedra, with as many degrees of freedom as nodes



Approximation Properties

- Take $\{\hat{\phi}_j\}_{j=1}^{\infty}$ to be an $a(\cdot, \cdot)$ -orthogonal basis for H^1
- $\{\hat{\phi}_j\}_{j=1}^n$ is a basis for $\mathcal{V} \subset H^1$

Writing $p = \sum_{j=1}^{\infty} \hat{p}_j \hat{\phi}_j$, $p_{\mathcal{V}} = \sum_{j=1}^n \hat{p}_j \hat{\phi}_j$

$$a(p - p_{\mathcal{V}}, p - p_{\mathcal{V}}) = \sum_{j=n+1}^{\infty} \hat{p}_j^2 a(\hat{\phi}_j, \hat{\phi}_j)$$

Want the projection of p onto \mathcal{V}^{\perp} to be small in the $a(\cdot, \cdot)$ -norm

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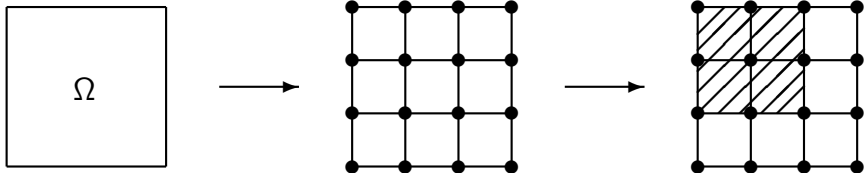
For a general q (+ BCs), $\hat{p}_j = \frac{\int_{\Omega} q \hat{\phi}_j}{a(\hat{\phi}_j, \hat{\phi}_j)}$

- Important to capture modes where $\frac{\int_{\Omega} q \hat{\phi}_j}{a(\hat{\phi}_j, \hat{\phi}_j)}$ is large
- Important to capture functions where $\frac{a(\varphi, \varphi)}{\langle \varphi, \varphi \rangle}$ is small

Multiscale Finite Element Method

Compute nodal basis of modes where $\frac{a(\varphi, \varphi)}{\langle \varphi, \varphi \rangle}$ is small

- Given Ω , partition into elements on scale for computation
- For each node, choose non-zero support over neighboring elements



T. Hou and X. Wu, *J. Comput. Phys.*, **134**, pp. 169–189, 1997.

T. Hou, X. Wu, and Z. Cai, *Math. Comp.*, **68**, pp. 913–943, 1999.

Multiscale Finite Element Method

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- Nodal basis implies $\phi_i(\mathbf{x}_j) = \delta_{ij}$
- Take $\phi_i(\mathbf{x}) = 0$ on boundary of its support

Can $\phi_i = \operatorname{argmin} \left\{ \frac{a(\varphi, \varphi)}{\langle \varphi, \varphi \rangle} : \varphi(\mathbf{x}_j) = \delta_{ij}, \varphi(\mathbf{x}) = 0 \text{ on } \partial\Omega_i \right\}$?

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I don't know.

Hou et al. ignore the denominator

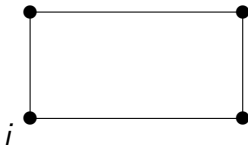
- define ϕ_i piecewise on each element
- fix boundary conditions and solve $a(\phi_i, \varphi) = 0$ on interior

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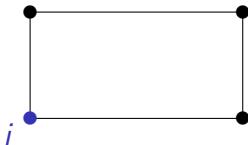
Artificial Boundary Conditions

Consider the element adjacent to node i ,



Artificial Boundary Conditions

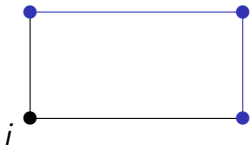
Consider the element adjacent to node i ,



- Fix $\phi_i(\mathbf{x}_i) = 1$

Artificial Boundary Conditions

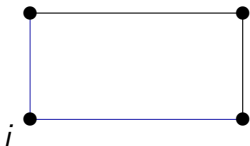
Consider the element adjacent to node i ,



- Fix $\phi_i(\mathbf{x}_i) = 1$
- Set $\phi_i(\mathbf{x}) = 0$ on $\partial\Omega_i$

Artificial Boundary Conditions

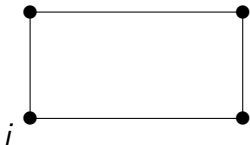
Consider the element adjacent to node i ,



- Fix $\phi_i(\mathbf{x}_i) = 1$
- Set $\phi_i(\mathbf{x}) = 0$ on $\partial\Omega_i$
- Impose boundary conditions on remaining edges

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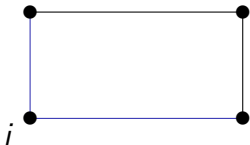
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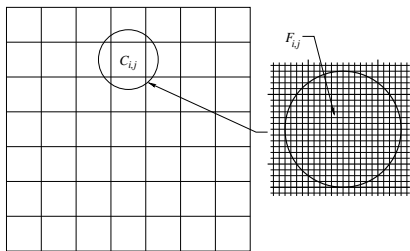
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- **Impose boundary conditions on remaining edges**
- Solve $a(\phi_i, \varphi) = 0$ in interior

Exact boundary conditions aren't known

- use linear
- solve one-dimensional problem along edge

Computational Cost of MSFEM

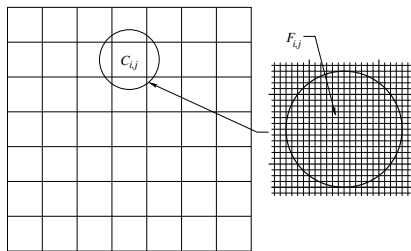
For each node of each element, need to compute basis function



- constant permeability tensor given on each fine-scale cell $F_{i,j}$
- choose computational scale, $C_{i,j}$
- solve for basis function of node (k, l) over $C_{i,j}$

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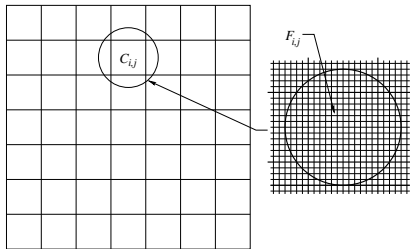
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We had three goals for our basis:

- good approximation
- **easy to calculate A and \mathbf{q}**
- easy to solve $A\mathbf{p} = \mathbf{q}$

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What is the cost of finding four basis functions over each element, compared to solving fine-scale equations?

Multigrid: Relaxation on $A\mathbf{x} = \mathbf{b}$

- Want to improve approximation, $\mathbf{x}^{(0)}$
- Introduce residual, $\mathbf{r}^{(0)} = \mathbf{b} - A\mathbf{x}^{(0)} = A(\mathbf{x} - \mathbf{x}^{(0)})$
- Take $\mathbf{x}^{(1)} = \mathbf{x}^{(0)} + D^{-1}\mathbf{r}^{(0)}$, for $D^{-1} \approx A^{-1}$

Error propagation form: $\mathbf{e}^{(1)} = (I - D^{-1}A) \mathbf{e}^{(0)}$

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Error propagation form: $\mathbf{e}^{(n)} = (I - D^{-1}A)^n \mathbf{e}^{(0)}$

Jacobi and Gauss-Seidel may be slow to converge, but their **failure is structured**

- Eigenvectors of small eigenvalues of $D^{-1}A$ are slow to change
- Can we use this to our advantage?

Multigrid: Subspace Correction

Dominant error after relaxation lies in a subspace

What if we could resolve this error by another process that acted only on the subspace?

Need

- complementary process
- way to combine its results with relaxation

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Want a map from the subspace to the whole space.

Interpolation!

Multigrid: Variational Coarsening

- Have $\mathbf{x}^{(1)}$, approximation after relaxation
- Let P be map from any subspace to whole space
- Corrected approximation will be $\mathbf{x}^{(2)} = \mathbf{x}^{(1)} + P\mathbf{x}_c$

What is the **best** \mathbf{x}_c for correction?

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What is the **best** \mathbf{x}_c for correction?

Symmetric and positive-definite matrix, A , defines an inner product and a norm:

$$\langle \mathbf{x}, \mathbf{y} \rangle_A = \mathbf{y}^T A \mathbf{x} \quad \text{and} \quad \|\mathbf{x}\|_A^2 = \mathbf{x}^T A \mathbf{x}$$

Best then means closest to the exact solution in norm:

$$\mathbf{y}^* = \underset{\mathbf{y}}{\operatorname{argmin}} \|\mathbf{x} - \mathbf{y}\|_A$$

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What is the **best** \mathbf{x}_c for correction?

Closest approximation to \mathbf{x} after correction given by

$$\mathbf{x}_c = \underset{\mathbf{y}_c}{\operatorname{argmin}} \|\mathbf{x} - (\mathbf{x}^{(1)} + P\mathbf{y}_c)\|_A$$

Best \mathbf{x}_c satisfies $(P^T A P)\mathbf{x}_c = P^T A(\mathbf{x} - \mathbf{x}^{(1)}) = P^T \mathbf{r}^{(1)}$

Multigrid: the V-Cycle

Multigrid Components

$$\text{Relax: } \mathbf{x}^{(1)} = \mathbf{x}^{(0)} + \mathbf{D}^{-1} \mathbf{r}^{(0)}$$

- Relaxation

- Use a relaxation process (such as Jacobi or Gauss-Seidel) to damp errors
- Remaining error satisfies $\mathbf{Ae}^{(1)} = \mathbf{r}^{(1)} = \mathbf{b} - \mathbf{Ax}^{(1)}$

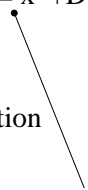
Multigrid: the V-Cycle

Multigrid Components

- Relaxation
- Restriction

$$\text{Relax: } \mathbf{x}^{(1)} = \mathbf{x}^{(0)} + \mathbf{D}^{-1} \mathbf{r}^{(0)}$$

Restriction



- Transfer residual to subspace
- Compute $P^T \mathbf{r}^{(1)}$

Multigrid: the V-Cycle

Multigrid Components

- Relaxation
- Restriction
- Subspace Correction

$$\text{Relax: } \mathbf{x}^{(1)} = \mathbf{x}^{(0)} + \mathbf{D}^{-1} \mathbf{r}^{(0)}$$

Restriction

$$\text{Solve: } \mathbf{P}^T \mathbf{A} \mathbf{P} \mathbf{x}_c = \mathbf{P}^T \mathbf{r}^{(1)}$$

- Use subspace correction to eliminate dominating errors
- Best correction, \mathbf{x}_c , in terms of A -norm satisfies

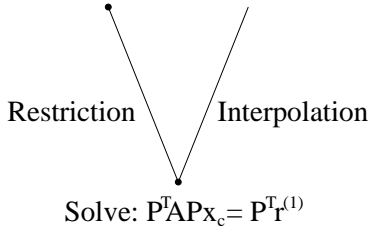
$$\mathbf{P}^T \mathbf{A} \mathbf{P} \mathbf{x}_c = \mathbf{P}^T \mathbf{r}^{(1)}$$

Multigrid: the V-Cycle

Multigrid Components

- Relaxation
 - Restriction
 - Subspace Correction
 - Interpolation
-
- Transfer correction to fine scale
 - Compute $\mathbf{x}^{(2)} = \mathbf{x}^{(1)} + P\mathbf{x}_c$

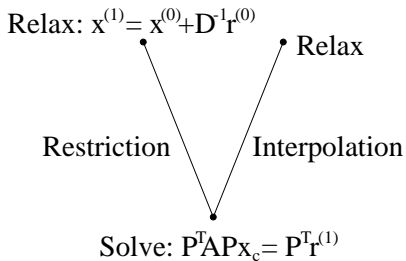
$$\text{Relax: } \mathbf{x}^{(1)} = \mathbf{x}^{(0)} + \mathbf{D}^{-1}\mathbf{r}^{(0)}$$



Multigrid: the V-Cycle

Multigrid Components

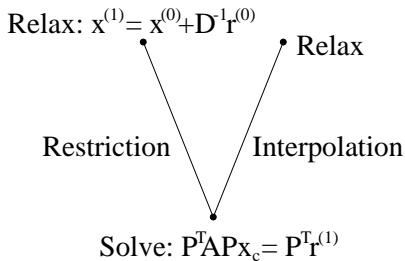
- Relaxation
- Restriction
- Subspace Correction
- Interpolation
- Relaxation
- Relax once again to damp errors introduced in subspace correction



Multigrid: the V-Cycle

Multigrid Components

- Relaxation
- Restriction
- Subspace Correction
- Interpolation
- Relaxation



Direct solution of coarse-grid problem isn't practical

Recursion!

Apply same methodology to solve coarse-grid problem

Multigrid: Operator-Induced Interpolation

Success of multigrid iteration depends on how well the range of P captures the slow-to-converge modes of relaxation

- For simple relaxation, **slow-to-converge modes** are close to **eigenvectors of A with small eigenvalues**
- Knowing structure of A (or continuum problem that generated it) allows effective choice of P

For $-\nabla \cdot \mathcal{K} \nabla p$, Black Box MG reduces error in the A -norm

- by a factor bounded less than 1 per iteration
- at a cost per iteration proportional to the size of A

MSFEM and Optimal Solvers

For scalar elliptic PDEs, discretized by standard finite elements, **multigrid is an optimal solver**.

- Error-reduction factor bounded independent of matrix size
- Iteration cost is bounded proportional to matrix size

In essence, solving a problem with $2n$ degrees of freedom takes twice as long as solving one with n degrees of freedom.

For MSFEM:

- Each basis function requires fine-scale solve over each element in its support
- Total cost is proportional to number of fine-scale nodes
- **Same as cost of solving fine-scale problem itself!**

Multigrid and Approximation

Optimal approximation properties rely on representing functions where $\frac{a(\varphi, \varphi)}{\langle \varphi, \varphi \rangle}$ is small

Operator-Induced Interpolation, P ,

- chosen based on discrete operator
- must accurately represent modes where $\frac{\mathbf{x}^T A \mathbf{x}}{\mathbf{x}^T \mathbf{x}}$ is small

Variational coarsening

- restricts A to range of interpolation
- explicitly constructs coarse-scale discrete model,
 $A_c = P^T A P$

Modes needed for good approximation properties are also needed for good multigrid performance

Implicit Basis Functions

Fine-scale finite-element discretization:

$$A_{ij} = \mathbf{e}_j^T \mathbf{A} \mathbf{e}_i = \int_{\Omega} (\mathcal{K}(\mathbf{x}) \nabla \phi_j) \cdot \nabla \phi_i$$

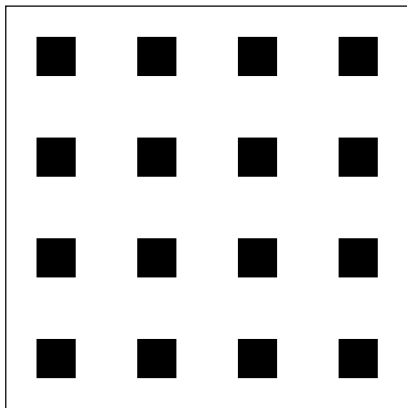
Variational coarsening gives coarse-grid operator,

$$\begin{aligned}(A_c)_{ij} &= (P^T A P)_{ij} = (P \hat{\mathbf{e}}_j)^T A (P \hat{\mathbf{e}}_i) \\ &= \int_{\Omega} \left(\mathcal{K}(\mathbf{x}) \nabla \left(\sum_k p_{kj} \phi_k \right) \right) \cdot \nabla \left(\sum_l p_{li} \phi_l \right) \\ &= \int_{\Omega} \left(\mathcal{K}(\mathbf{x}) \nabla \hat{\phi}_j \right) \cdot \nabla \hat{\phi}_i\end{aligned}$$

Variational coarsening **implicitly defines basis functions** on coarse scale, $\hat{\phi}_i = \sum_l p_{li} \phi_l$.

Multigrid Basis Functions

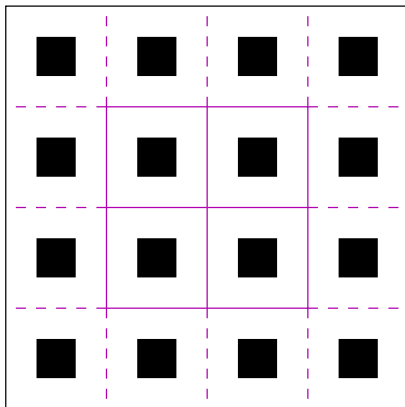
Variational multigrid defines a multiscale finite element basis



Periodic tiling of inclusion problem: $\mathcal{K} = 1000$ in inclusions,
 $\mathcal{K} = 1$ in background

Multigrid Basis Functions

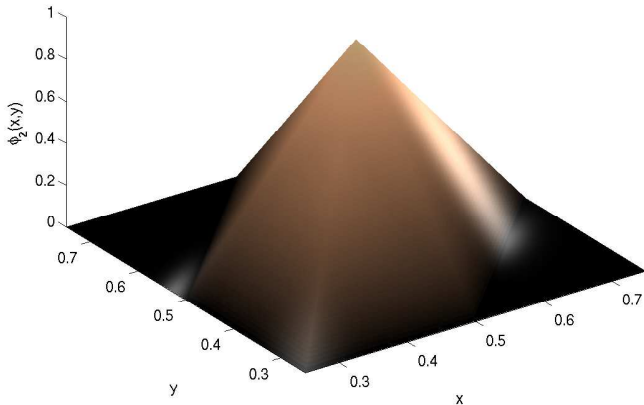
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Multigrid Basis Functions

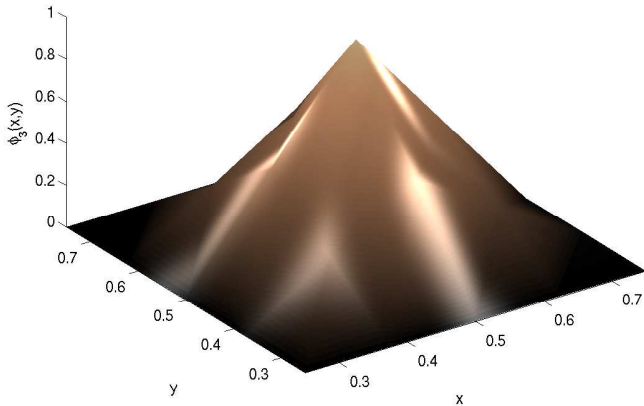
Variational multigrid defines a multiscale finite element basis



Bilinear basis function on coarse scale

Multigrid Basis Functions

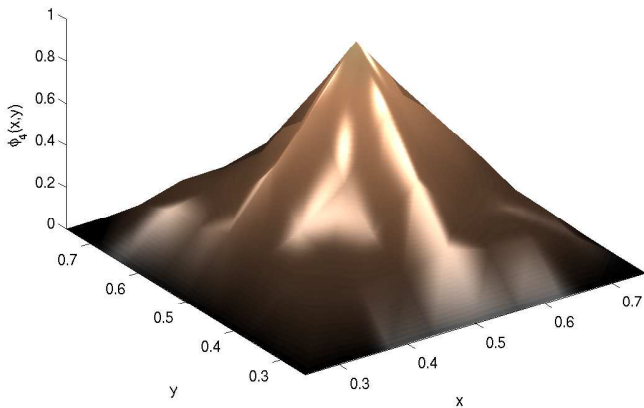
Variational multigrid defines a multiscale finite element basis



Basis function accounting for coarsest 2 scales

Multigrid Basis Functions

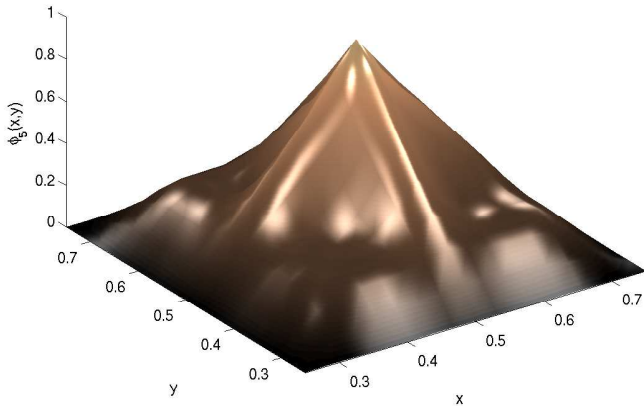
Variational multigrid defines a multiscale finite element basis



Basis function accounting for coarsest 3 scales

Multigrid Basis Functions

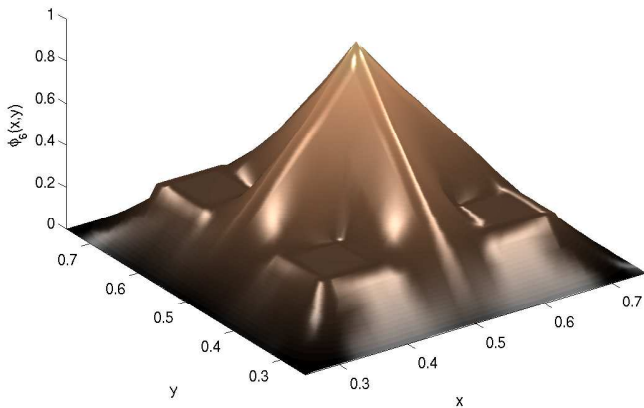
Variational multigrid defines a multiscale finite element basis



Basis function accounting for coarsest 4 scales

Multigrid Basis Functions

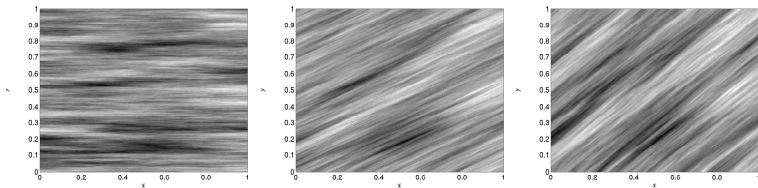
Variational multigrid defines a multiscale finite element basis



Basis function accounting for all scales

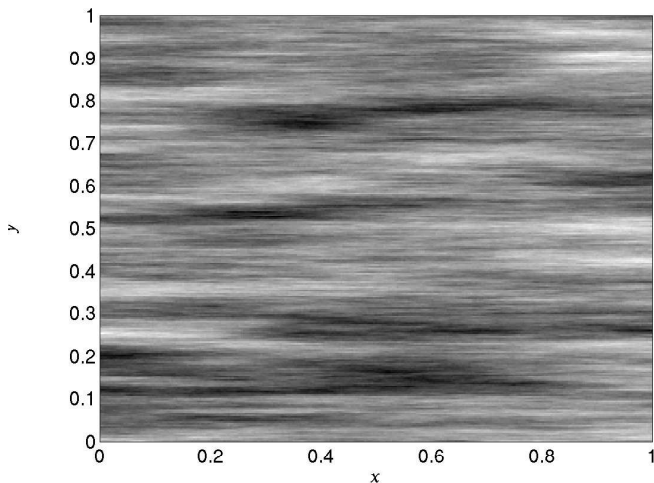
Geostatistical Media

- Principle axis of statistical anisotropy chosen
- Correlation length of 0.8 along axis, 0.04 across axis
- $\log_{10}(\mathcal{K})$ normally distributed with mean 0, variance 4



Multigrid Basis Functions

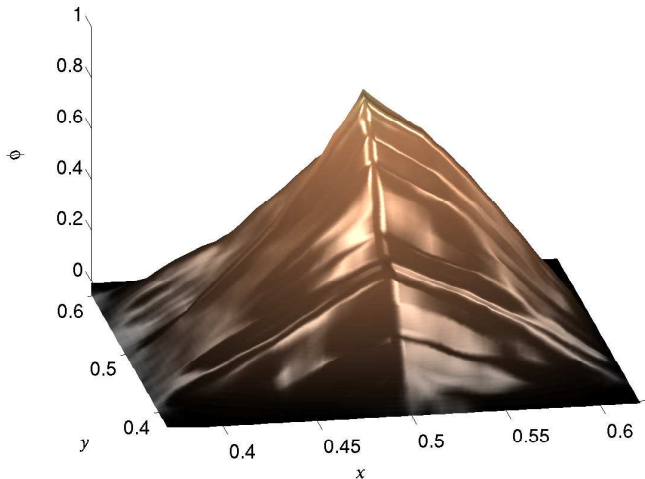
Variational multigrid defines a multiscale finite element basis



Permeability field for 0 degrees

Multigrid Basis Functions

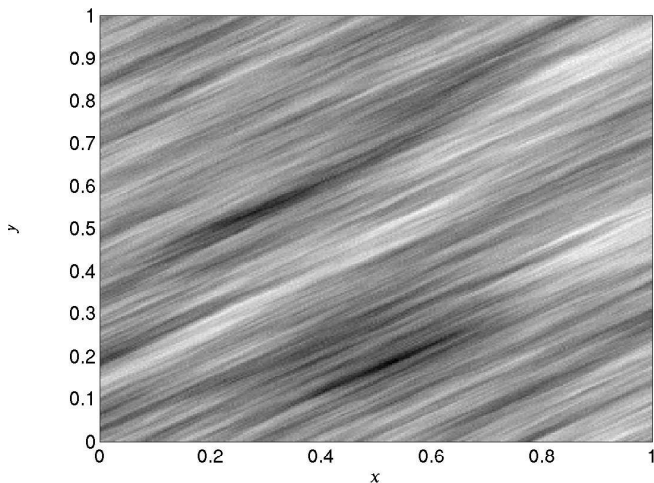
Variational multigrid defines a multiscale finite element basis



Basis for node at $(\frac{1}{2}, \frac{1}{2})$ for 0 degrees

Multigrid Basis Functions

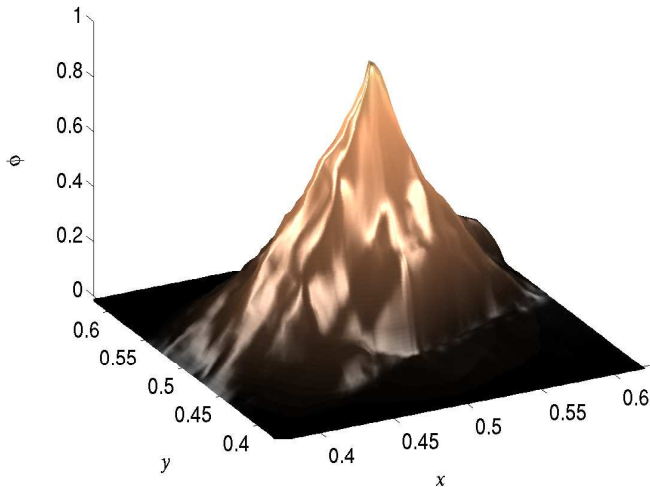
Variational multigrid defines a multiscale finite element basis



Permeability field for 30 degrees

Multigrid Basis Functions

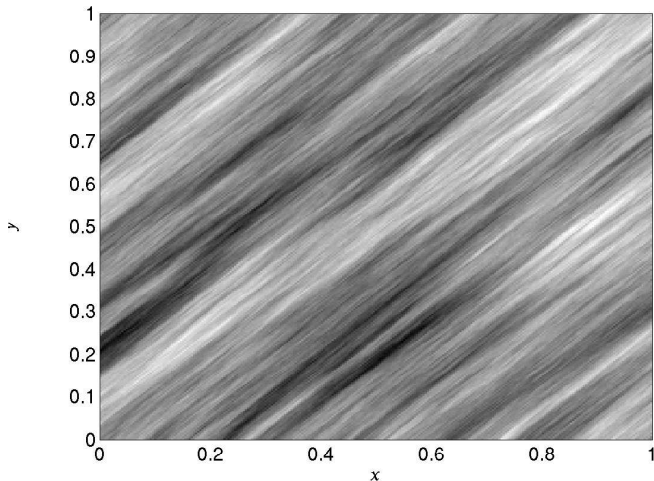
Variational multigrid defines a multiscale finite element basis



Basis for node at $(\frac{1}{2}, \frac{1}{2})$ for 30 degrees

Multigrid Basis Functions

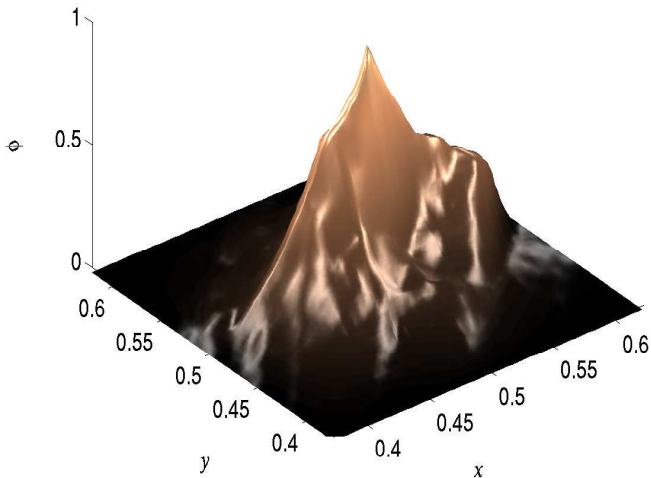
Variational multigrid defines a multiscale finite element basis



Permeability field for 45 degrees

Multigrid Basis Functions

Variational multigrid defines a multiscale finite element basis



Basis for node at $(\frac{1}{2}, \frac{1}{2})$ for 45 degrees

Implicit Upscaling

Multigrid coarse-scale operators represent consistently upscaled models

- Equivalent to finite element discretization with implicit basis functions
- Accurately represent small-Rayleigh quotient modes
- Require no fine-scale solution to form coarse-scale model
- Are easily solved using multigrid

Implicit Upscaling

Multigrid coarse-scale operators represent consistently upscaled models

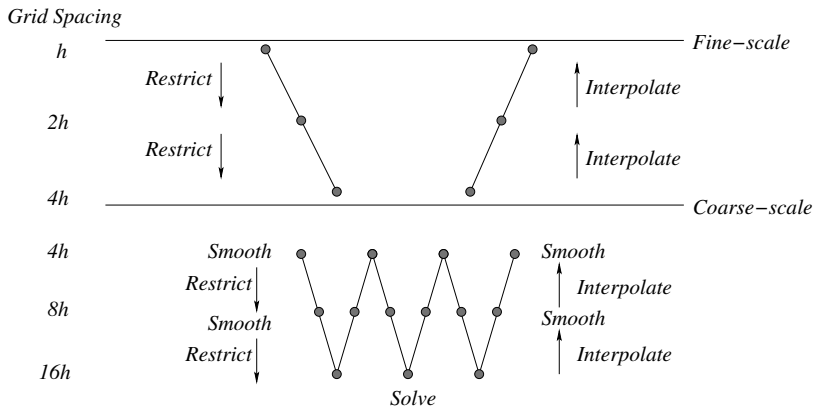
- Equivalent to finite element discretization with implicit basis functions
- Accurately represent small-Rayleigh quotient modes
- Require no fine-scale solution to form coarse-scale model
- Are easily solved using multigrid

Algorithm:

- Form fine-scale discrete model
- Use operator-induced variation coarsening to create coarse-scale models
- Restrict sources and boundary conditions to chosen computational scale
- Solve model on chosen scale
- Interpolate solution to fine scale

The Multilevel Upscaling Algorithm

From a multigrid point of view, this is just not smoothing on scales finer than the coarse (computational) scale



Adaptivity

MLUPS framework is a natural setting for adaptivity

Variational multigrid approach

- creates a hierarchy of models at different scales
- naturally restricts A -norm to coarse scales
- allows for coarse-scale error estimation
- allows for local improvement on scales finer than chosen coarse scale

Nonlinear multigrid (FAS) framework gives flexible framework
for error estimation and control

Test problems

Two-dimensional geostatistical media

- Chosen axis of statistical anisotropy
- Correlation lengths of 0.8 along axis, 0.04 across axis
- $\log_{10}(\mathcal{K})$ normally distributed with mean 0, variance of 4

Boundary Conditions

- mean uniform flow driven by imposed Dirichlet boundaries
- $h(0, y) = 1, h(1, y) = 0$
- Homogeneous Neumann boundaries on top and bottom

Test problems

\mathcal{K} chosen to be piecewise constant on 256×256 mesh

Four algorithms:

- Bilinear finite elements on 256×256 mesh
- MSFEM with coarse scale of 8×8 elements
- MLUPS with coarse scale of 8×8 elements
- MLUPSa with coarse scale of 8×8 elements
 - ▶ MLUPSa is MLUPS with relaxation on all finer scales in final interpolation

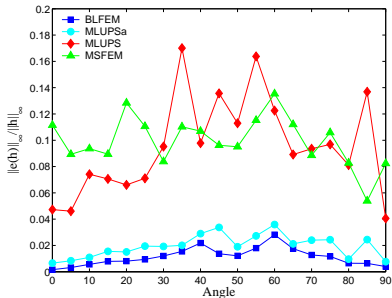
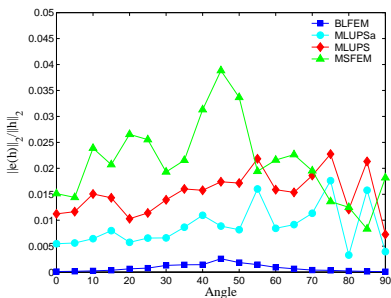
Accuracy measured versus solution of problem on 2048×2048 grid.

Errors in Fine-Scale Pressures

Errors are measured in discrete vector norms:

$$\|e(h)\|_2 = \left(\frac{1}{N} \sum_{i=1}^N e(h)_i^2 \right)^{\frac{1}{2}}, \quad \|e(h)\|_\infty = \max_i |e(h)_i|,$$

evaluated at each node on the 2048×2048 mesh.

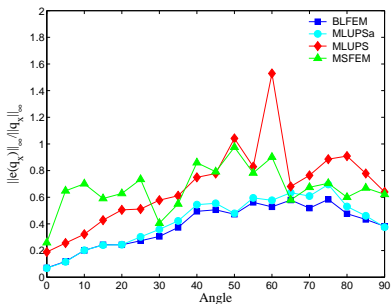
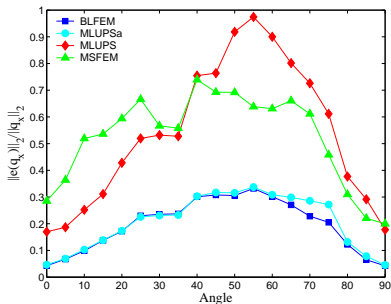


Errors in Fine-Scale Flux

Errors measured component-wise in discrete vector norms:

$$\|e(\mathbf{Q} \cdot \hat{\mathbf{x}})\|_2 = \left(\frac{1}{N} \sum_{i=1}^N e(\mathbf{Q} \cdot \hat{\mathbf{x}})_i^2 \right)^{\frac{1}{2}}, \quad \|e(\mathbf{Q} \cdot \hat{\mathbf{x}})\|_\infty = \max_i |e(\mathbf{Q} \cdot \hat{\mathbf{x}})_i|,$$

evaluated at cell-centers of the 2048×2048 mesh.

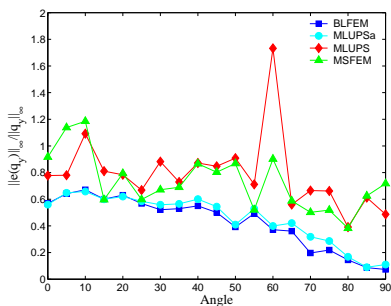
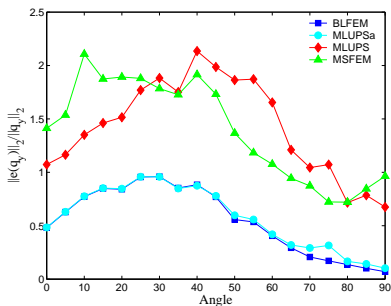


Errors in Fine-Scale Flux

Errors measured component-wise in discrete vector norms:

$$\|e(\mathbf{Q} \cdot \hat{\mathbf{y}})\|_2 = \left(\frac{1}{N} \sum_{i=1}^N e(\mathbf{Q} \cdot \hat{\mathbf{y}})_i^2 \right)^{\frac{1}{2}}, \quad \|e(\mathbf{Q} \cdot \hat{\mathbf{y}})\|_\infty = \max_i |e(\mathbf{Q} \cdot \hat{\mathbf{y}})_i|,$$

evaluated at cell-centers of the 2048×2048 mesh.



Summary

- Accurate simulation relies on resolving heterogeneities in media
- Coefficient upscaling only valid in special cases
- Variational principles allow accurate upscaling of model
- MSFEM approach accurate, but expensive
- Operator-induced multigrid also captures necessary modes
- Multilevel Upscaling (MLUPS) approach accurate, 15 times cheaper than MSFEM

S.P. MacLachlan & J.D. Moulton, *Water Resour. Res.*, **42**, 2006

<http://www.cs.umn.edu/~maclach/research/multiscale.pdf>

Ongoing Research

- Relationship between coefficient and model upscaling
- Accounting for mass conservation in variational framework
- Removing artificial boundary conditions from MSFEM
- Relationship between MSFEM and MLUPS
- 3D, time-dependent, nonlinear
- Stochastic coefficients, stochastic FEM
- Local error estimation and adaptivity

S.P. MacLachlan & J.D. Moulton, *Water Resour. Res.*, **42**, 2006

<http://www.cs.umn.edu/~maclach/research/multiscale.pdf>