

Lecture 4a

Using the SAMR approach for elliptic and parabolic problems

*Course Block-structured Adaptive Mesh Refinement Methods for Conservation Laws
Theory, Implementation and Application*

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Outline

Adaptive geometric multigrid methods

- Linear iterative methods for Poisson-type problems

- Multi-level algorithms

- Multigrid algorithms on SAMR data structures

- Example

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Comments on parabolic problems

Poisson equation

$$\begin{aligned}\Delta q(\mathbf{x}) &= \psi(\mathbf{x}), \quad \mathbf{x} \in \Omega \subset \mathbb{R}^d, \quad q \in C^2(\Omega), \quad \psi \in C^0(\Omega) \\ q &= \psi^\Gamma(\mathbf{x}), \quad \mathbf{x} \in \partial\Omega\end{aligned}$$

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Discrete Poisson equation in 2D:

$$\frac{Q_{j+1,k} - 2Q_{jk} + Q_{j-1,k}}{\Delta x_1^2} + \frac{Q_{j,k+1} - 2Q_{jk} + Q_{j,k-1}}{\Delta x_2^2} = \psi_{jk}$$

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Operator

$$\mathcal{A}(Q_{\Delta x_1, \Delta x_2}) = \begin{bmatrix} & & \frac{1}{\Delta x_2^2} & & \\ \frac{1}{\Delta x_1^2} & - \left(\frac{2}{\Delta x_1^2} + \frac{2}{\Delta x_2^2} \right) & & \frac{1}{\Delta x_2^2} & \\ & & \frac{1}{\Delta x_2^2} & & \end{bmatrix} Q(x_{1,j}, x_{2,k}) = \psi_{jk}$$

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$$Q_{jk} = \frac{1}{\sigma} \left[(Q_{j+1,k} + Q_{j-1,k})\Delta x_2^2 + (Q_{j,k+1} + Q_{j,k-1})\Delta x_1^2 - \Delta x_1^2 \Delta x_2^2 \psi_{jk} \right]$$

$$\text{with } \sigma = \frac{2\Delta x_1^2 + 2\Delta x_2^2}{\Delta x_1^2 \Delta x_2^2}$$

Iterative methods

Jacobi iteration

$$Q_{jk}^{m+1} = \frac{1}{\sigma} \left[(Q_{j+1,k}^m + Q_{j-1,k}^m) \Delta x_2^2 + (Q_{j,k+1}^m + Q_{j,k-1}^m) \Delta x_1^2 - \Delta x_1^2 \Delta x_2^2 \psi_{jk} \right]$$

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Efficient parallelization / patch-wise application not possible!

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Checker-board or Red-Black Gauss Seidel iteration

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 for $j + k \bmod 2 = 0$
- $$Q_{jk}^{m+1} = \frac{1}{\sigma} \left[(Q_{j+1,k}^{m+1} + Q_{j-1,k}^{m+1}) \Delta x_2^2 + (Q_{j,k+1}^{m+1} + Q_{j,k-1}^{m+1}) \Delta x_1^2 - \Delta x_1^2 \Delta x_2^2 \psi_{jk} \right]$$
 for $j + k \bmod 2 = 1$

Iterative methods

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for $j + k \bmod 2 = 1$

Gauss-Seidel methods require $\sim 1/2$ of iterations than Jacobi method, however, iteration count still proportional to number of unknowns [Hackbusch, 1994]

Smoothing vs. solving

ν iterations with iterative linear solver

$$Q^{m+\nu} = \mathcal{S}(Q^m, \psi, \nu)$$

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Defect after m iterations

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Defect after $m + \nu$ iterations

$$d^{m+\nu} = \psi - \mathcal{A}(Q^{m+\nu}) = \psi - \mathcal{A}(Q^m + v_\nu^m) = d^m - \mathcal{A}(v_\nu^m)$$

with correction

$$v_\nu^m = \mathcal{S}(\vec{0}, d^m, \nu)$$

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Neglecting the sub-iterations in the smoother we write

$$Q^{n+1} = Q^n + v = Q^n + \mathcal{S}(d^n)$$

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Observation: Oscillations are damped faster on coarser grid.

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Observation: Oscillations are damped faster on coarser grid.

Coarse grid correction:

$$Q^{n+1} = Q^n + v = Q^n + \mathcal{P}\mathcal{S}\mathcal{R}(d^n)$$

where \mathcal{R} is suitable restriction operator and \mathcal{P} a suitable prolongation operator

Two-grid correction method

Relaxation on current grid:

$$\bar{Q} = \mathcal{S}(Q^n, \psi, \nu)$$

$$Q^{n+1} = \bar{Q} + \mathcal{PS}(\vec{0}, \cdot, \mu) \mathcal{R}(\psi - \mathcal{A}(\bar{Q}))$$

Two-grid correction method

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Algorithm:

$$\begin{aligned}\bar{Q} &:= \mathcal{S}(Q^n, \psi, \nu) \\ d &:= \psi - \mathcal{A}(\bar{Q})\end{aligned}$$

$$\begin{aligned}d_c &:= \mathcal{R}(d) \\ v_c &:= \mathcal{S}(\vec{0}, d_c, \mu) \\ v &:= \mathcal{P}(v_c) \\ Q^{n+1} &:= \bar{Q} + v\end{aligned}$$

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$$Q^{n+1} := \bar{Q} + v$$

with smoothing:

$$d := \psi - \mathcal{A}(Q)$$

$$v := \mathcal{S}(0, d, \nu)$$

$$r := d - \mathcal{A}(v)$$

$$d_c := \mathcal{R}(r)$$

$$v_c := \mathcal{S}(0, d_c, \mu)$$

$$v := v + \mathcal{P}(v_c)$$

$$Q^{n+1} := Q + v$$

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with pre- and post-iteration:

$$d := \psi - \mathcal{A}(Q)$$

$$v := \mathcal{S}(0, d, \nu_1)$$

$$r := d - \mathcal{A}(v)$$

$$d_c := \mathcal{R}(r)$$

$$v_c := \mathcal{S}(0, d_c, \mu)$$

$$v := v + \mathcal{P}(v_c)$$

$$d := d - \mathcal{A}(v)$$

$$r := \mathcal{S}(0, d, \nu_2)$$

$$Q^{n+1} := Q + v + r$$

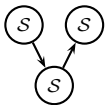
[Hackbusch, 1985]

Multi-level methods and cycles

V-cycle

$$\gamma = 1$$

2-grid

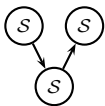


Multi-level methods and cycles

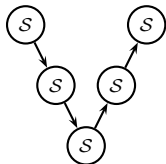
V-cycle

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2-grid



3-grid

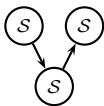


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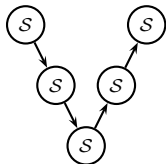
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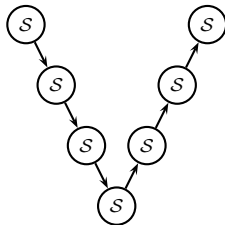
2-grid



3-grid



4-grid

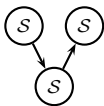


Multi-level methods and cycles

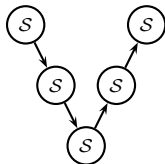
V-cycle

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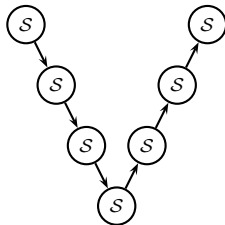
2-grid



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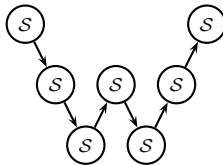


4-grid



W-cycle

$$\gamma = 2$$

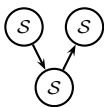


Multi-level methods and cycles

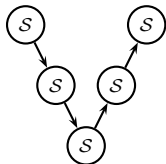
V-cycle

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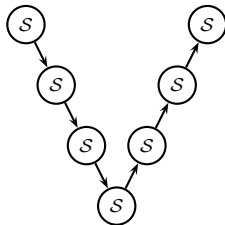
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3-grid

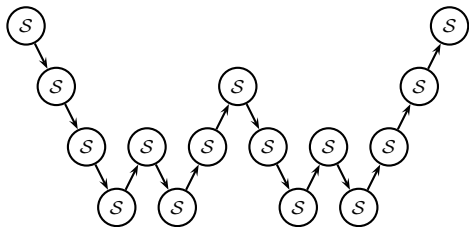
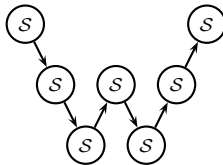


4-grid



W-cycle

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[Hackbusch, 1985] [Wesseling, 1992] ...

Stencil modification at coarse-fine boundaries in 1D

1D Example: Cell j , $\psi - \nabla \cdot \nabla q = 0$

$$d_j^l = \psi_j - \frac{1}{\Delta x_l} \left(\frac{1}{\Delta x_l} (Q_{j+1}^l - Q_j^l) - \frac{1}{\Delta x_l} (Q_j^l - Q_{j-1}^l) \right)$$

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H is approximation to *derivative* of Q^l .

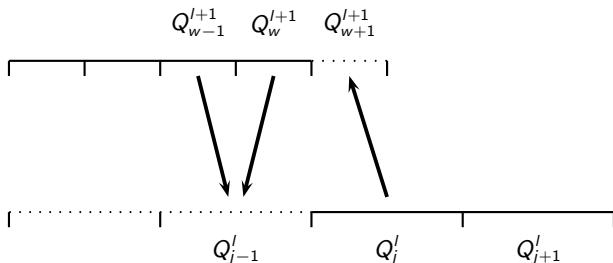
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Consider 2-level situation with $r_{l+1} = 2$:



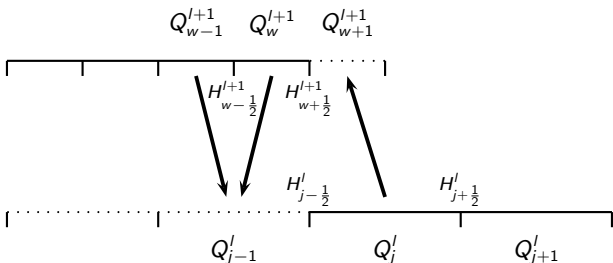
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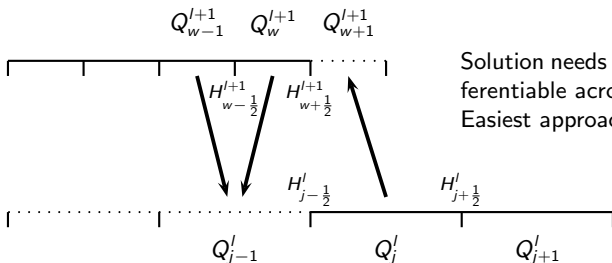
Stencil modification at coarse-fine boundaries in 1D

1D Example: Cell j , $\psi - \nabla \cdot \nabla q = 0$

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H is approximation to *derivative* of Q^l .

Consider 2-level situation with $r_{l+1} = 2$:



Solution needs to be continuously differentiable across interface.

Easiest approach: $H_{w+\frac{1}{2}}^{l+1} \equiv H_{j-\frac{1}{2}}^l$

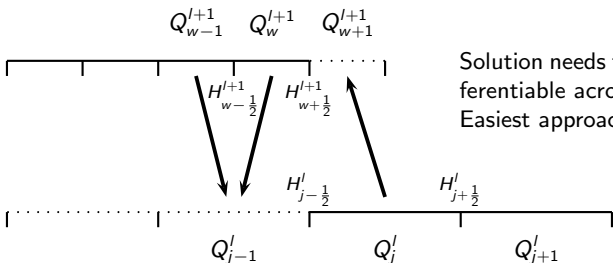
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H is approximation to derivative of Q^l .

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No specific modification necessary for 1D vertex-based stencils, cf. [Bastian, 1996]

Stencil modification at coarse-fine boundaries in 1D II

$$\text{Set } H_{w+\frac{1}{2}}^{l+1} = H_{\mathcal{I}}.$$

Stencil modification at coarse-fine boundaries in 1D II

Set $H_{w+\frac{1}{2}}^{l+1} = H_{\mathcal{I}}$. Inserting Q gives

$$\frac{Q_{w+1}^{l+1} - Q_w^{l+1}}{\Delta x_{l+1}} = \frac{Q_j^l - Q_w^{l+1}}{\frac{3}{2}\Delta x_{l+1}}$$

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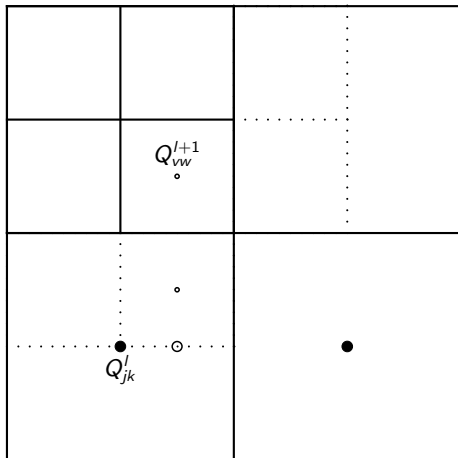
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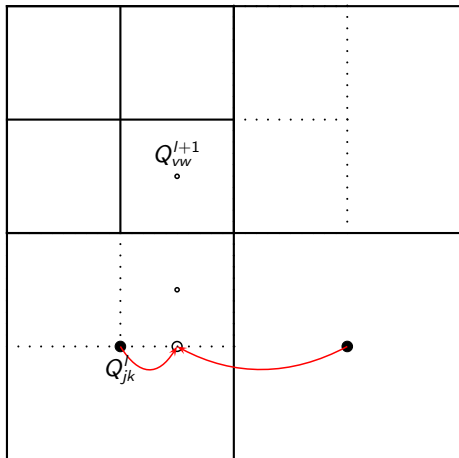
$$\check{d}_j^l = \psi_j - \frac{1}{\Delta x_l} \left(\frac{1}{\Delta x_l} (Q_{j+1}^l - Q_j^l) - \frac{2}{3\Delta x_{l+1}} (Q_j^l - Q_w^{l+1}) \right)$$

Stencil modification at coarse-fine boundaries: 2D



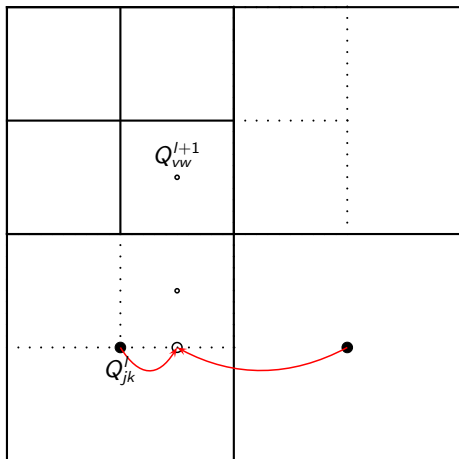
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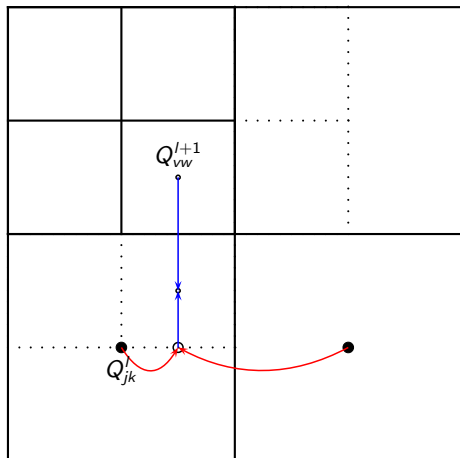
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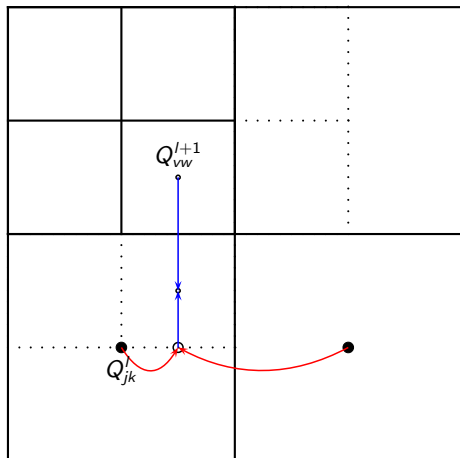
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In general:

$$Q_{v,w-1}^{l+1} = \left(1 - \frac{2}{r_{l+1} + 1} \right) Q_{vw}^{l+1} + \frac{2}{r_{l+1} + 1} \left((1-f) Q_{jk}^l + f Q_{j+1,k}^l \right)$$

with

$$f = \frac{x_{1,l+1}^v - x_{1,l}^j}{\Delta x_{1,l}}$$

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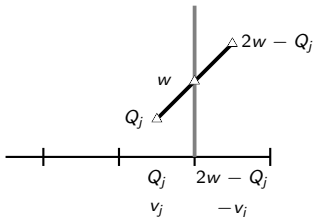
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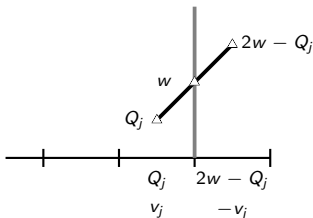
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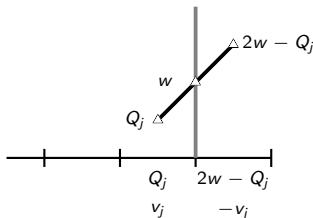
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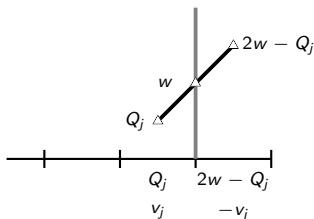
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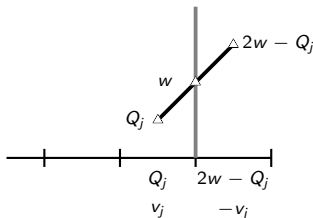
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- ▶ **Adaptation criteria**



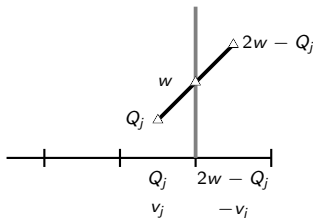
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- ▶ Looping instead of time steps and check of convergence



Additive geometric multigrid algorithm

AdvanceLevelMG(l) - Correction Scheme

Set ghost cells of Q^l

Calculate defect d^l from Q^l, ψ^l

$$d^l := \psi^l - \mathcal{A}(Q^l)$$

If ($l < l_{max}$)

 Calculate updated defect r^{l+1} from v^{l+1}, d^{l+1}

$$r^{l+1} := d^{l+1} - \mathcal{A}(v^{l+1})$$

 Restrict d^{l+1} onto d^l

$$d^l := \mathcal{R}_l^{l+1}(r^{l+1})$$

Do ν_1 smoothing steps to get correction v^l

$$v^l := \mathcal{S}(0, d^l, \nu_1)$$

If ($l > l_{min}$)

 Do $\gamma > 1$ times

 AdvanceLevelMG($l - 1$)

 Set ghost cells of v^{l-1}

 Prolongate and add v^{l-1} to v^l

$$v^l := v^l + \mathcal{P}_l^{l-1}(v^{l-1})$$

 If ($\nu_2 > 0$)

 Set ghost cells of v^l

 Update defect d^l according to v^l

$$d^l := d^l - \mathcal{A}(v^l)$$

 Do ν_2 post-smoothing steps to get r^l

$$r^l := \mathcal{S}(v^l, d^l, \nu_2)$$

 Add additional correction r^l to v^l

$$v^l := v^l + r^l$$

Add correction v^l to Q^l

$$Q^l := Q^l + v^l$$

Additive Geometric Multiplicative Multigrid Algorithm

Start - Start iteration on level l_{max}

For $l = l_{max}$ Downto $l_{min} + 1$ Do

 Restrict Q^l onto Q^{l-1}

 Regrid(0)

 AdvanceLevelMG(l_{max})

See also: [Trottenberg et al., 2001], [Canu and Ritzdorf, 1994]

Vertex-based: [Brandt, 1977], [Briggs et al., 2001]

Example

On $\Omega = [0, 10] \times [0, 10]$ use hat function

$$\psi = \begin{cases} -A_n \cos\left(\frac{\pi r}{2R_n}\right), & r < R_n \\ 0 & \text{elsewhere} \end{cases}$$

with $r = \sqrt{(x_1 - X_n)^2 + (x_2 - Y_n)^2}$
to define three sources with

n	A_n	R_n	X_n	Y_n
1	0.3	0.3	6.5	8.0
2	0.2	0.3	2.0	7.0
3	-0.1	0.4	7.0	3.0

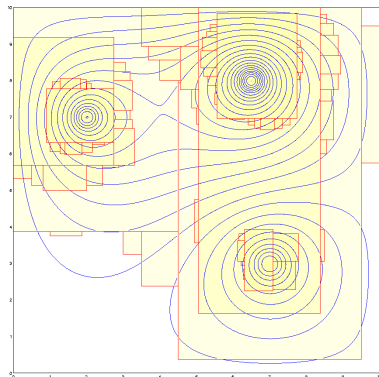
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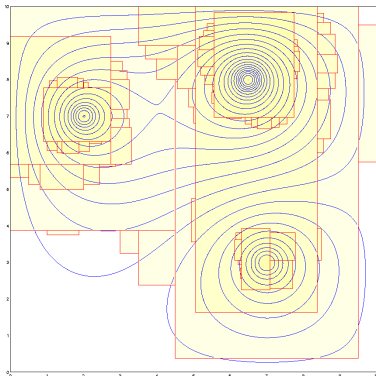
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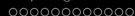


	128×128	1024×1024	1024×1024
l_{max}	3	0	0
l_{min}	-4	-7	-4
ν_1	5	5	5
ν_2	5	5	5
V-Cycles	15	16	341
Time [sec]	9.4	27.7	563

Stop at $\|d^l\|_{max} < 10^{-7}$ for $l \geq 0$, $\gamma = 1$, $r_l = 2$

Some comments on parabolic problems

- ▶ Consequences of time step refinement
- ▶ Level-wise elliptic solves vs. global solve
- ▶ If time step refinement is used an elliptic flux correction is unavoidable.
- ▶ The correction is explained in Bell, J. (2004). Block-structured adaptive mesh refinement. Lecture 2. Available at <https://ccse.lbl.gov/people/jbb/shortcourse/lecture2.pdf>.



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