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#### Overview

#### Model Problems

- Relaxation Methods
- Error convergence
- Multiple grids
- Performance
- Theoretical Considerations

Differential Equation in 1D

-u''(x) + au(x) = f(x)

for 0 < x < 1, a > 0Boundary: u(0) = u(1) = 0

 Partition continuous problem into n subintervals by "sampling" it at the grid points x<sub>i</sub> = jh, with h = 1/n

Grid Ω<sup>h</sup>:



Second order finite difference approximation

$$\frac{-v_{j-1} + 2v_j - v_{j+1}}{h^2} + av_j = f(x_j) \text{ with } 1 \le j \le n-1$$

 $v_0 = v_n = 0$ ; with **v** being the approximate solution to **u** Written in Matrix-Vector form

$$\frac{1}{h^{2}}\begin{bmatrix} 2+ah^{2} & -1 \\ -1 & 2+ah^{2} & -1 \\ & & \ddots \\ & & -1 & 2+ah^{2} \end{bmatrix} \cdot \begin{bmatrix} v_{1} \\ \vdots \\ \vdots \\ v_{n-1} \end{bmatrix} = \begin{bmatrix} f(x_{1}) \\ \vdots \\ \vdots \\ f(x_{n-1}) \end{bmatrix} = \begin{bmatrix} f_{1} \\ \vdots \\ \vdots \\ f_{n-1} \end{bmatrix}$$

Written compactly: Av = f

(Elliptic) Partial Differential Equation

-u<sub>xx</sub> - u<sub>yy</sub> + au = f(x,y)
for 0 < x,y < 1 , a > 0
Boundary: "Frame = 0"

 Sampled with a two-dimensional grid (n-1,m-1 interior grid points)



Sampling results in difference approximation

$$\frac{-v_{i-1,j} + 2v_{ij} - v_{i+1,j}}{h_x^2} + \frac{-v_{i,j-1} + 2v_{ij} - v_{i,j+1}}{h_y^2} + av_{ij} = f_{ij} \quad \text{with } 1 \le j \le n-1$$

$$v_{i0} = v_{in} = v_{0j} = v_{mj} = 0, \qquad 1 \le i, j \le m-1$$

Written in Matrix-Vector form



B looks almost like matrix from 1D - Problem - dimension is  $(n-1) \times (n-1)$ I is a  $(n-1) \times (n-1)$  identity matrix  $\Rightarrow$  Dimension of matrix is  $(m-1) \cdot (n-1) \times (m-1) \cdot (n-1)$ 

#### Model Problem 2D

#### Example: System for a=0, n=4, h=1



Again, written compactly: Av = f

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- To solve the PDE,  $\mathbf{u} = A^{-1}\mathbf{f}$  is too complicated
- Based on an estimated solution  $\mathbf{v}^{(0)} \rightarrow$  find better solution  $\mathbf{v}^{(1)}$  in next step
- Reduces norm of the error  $\mathbf{e} = \mathbf{u} \mathbf{v}$
- Use residual r = f − Av as a measure Relationship error / residual: Ae = r
   For exact solution v = u ⇒ r = 0
- For the following, split matrix A = D L U
   D: diagonal of A; L/U: lower/upper triangular part of A

- General approximation:  $\mathbf{v}^{(1)} = \mathbf{v}^{(0)} + B\mathbf{r}^0$ 
  - Try to find a B "close" to  $A^{-1}$ , as  $\mathbf{u} \mathbf{v} = A^{-1}\mathbf{r}$
- Jacobi scheme / Simultaneous displacement
  - j<sup>th</sup> component of v is calculated using the two neighbours from previous step

 $v_{j}^{(1)} = \frac{1}{2} (v_{j-1}^{(0)} + v_{j+1}^{(0)} + h^{2} f_{j}), \qquad 1 \le j \le n-1$ Jacobi iteration matrix :  $\mathbf{R}_{J} = D^{-1} (L+U)$  $v^{(1)} = R_{J} v^{(0)} + D^{-1} f$ 

 Solves the PDE locally (compare original problem: -u<sub>j-1</sub> + 2u<sub>j</sub> - u<sub>j+1</sub> = h<sup>2</sup>f<sub>j</sub>)

- Weighted or damped Jacobi method
  - Weighting factor  $0 < \omega < 1$

 $v_{j}^{(1)} = (1 - \omega)v_{j}^{(0)} + \omega v_{j}^{*}, \qquad 1 \le j \le n - 1$ with  $v_{j}^{*} = \frac{1}{2}(v_{j-1}^{(0)} + v_{j+1}^{(0)} + h^{2}f_{j})$  (like above) Weighted Jacobi iteration matrix :  $\mathbf{R}_{\omega} = (1 - \omega)I + \omega R_{J}$  $v^{(1)} = R_{\omega}v^{(0)} + \omega D^{-1}f$ 

- Gauss-Seidel
  - Like Jacobi, but components updated immediately
  - Reduces storage requirements

$$v_{j} \leftarrow \frac{1}{2}(v_{j-1} + v_{j+1} + h^{2}f_{j}), \leftarrow \text{meaning "overwrite"}$$
  
formally  $: v_{j}^{(1)} = \frac{1}{2}(v_{j-1}^{(1)} + v_{j+1}^{(0)} + h^{2}f_{j})$ 

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- Simplified problem: Au = 0
   ⇒ v should converge to 0, and e = v
- In what way does weighted Jacobi decrease the error?
   Analyse eigenvectors of iteration matrix
- Eigenvectors  $w_k$  of matrices A and  $R_{\omega}$  $w_{k,j} = \sin\left(\frac{jk\pi}{n}\right)$ , with  $1 \le k \le n-1$ ,  $0 \le j \le n$ 
  - Vector w<sub>k</sub> is also the k<sup>th</sup> Fourier mode
- Eigen values  $\lambda_k$  of matrix  $R_\omega$  (generally:  $R_\omega w_k = \lambda_k w_k$ )  $\lambda_k(R_\omega) = 1 - 2\omega \sin^2 \left(\frac{k\pi}{2n}\right)$ , with  $1 \le k \le n - 1$ 
  - For  $0 \le \omega < 1 \Rightarrow |\lambda_k| < 1$ , iteration converges

- **Eigen values**  $\lambda_k(R_\omega) = 1 2\omega \sin^2\left(\frac{k\pi}{2n}\right)$ , with  $1 \le k \le n-1$
- Smooth, low-frequency Fourier modes of  $e: 1 \le k \le \frac{1}{2}n$ 
  - $|\lambda_k|$  is close to 1  $\Rightarrow$  no satisfactory damping
- Oscillatory, high-frequency modes:  $\frac{1}{2}n \le k \le n-1$ 
  - For the right  $\omega$ ,  $|\lambda_k|$  is close to  $0 \Rightarrow$  good damping
  - Optimal damping for  $\omega = \frac{2}{3}$



Damping diagram for the weighted Jacobi method



- Oscillatory modes of the error are removed quite well
- Smooth modes are hardly damped.

#### **Error convergence**

- Example code in MATLAB
  - Grid n = 64
  - Initial error modes 2 and 16
  - Solves –u''(x) = 0

```
n = 64;
% components of A
D = 2 * diag(ones(n-1,1),0); U = d
% iteration matrices
w=2/3; RJ = inv(D) * (L+U); RW =
% init f=0, v with modes 2 and 16
f = zeros(n-1,1);
v = transpose(sin((1:n-1) * 2 * pi
plot(v); hold on
% do 10 iterations
for i = 1:10
v = RW*v + 0; end
plot(v);
```



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- Fundamental idea of multigrid
  - Make smooth modes look oscillatory!
  - Smooth mode on Ω<sup>h</sup> looks oscillatory on grid Ω<sup>nh</sup>
  - A "hierarchy of discretizations" is used to solve the problem of small damping for smooth modes



#### **Multiple Grids**

- Intergrid Transfer coarse  $\rightarrow$  fine: Interpolation
  - $\Omega^{2h} \rightarrow \Omega^{h}$ , "Upsampling"
  - Linear interpolation is effective

$$\begin{aligned} v_{2j}^{h} &= v_{j}^{2h} \\ v_{2j+1}^{h} &= \frac{1}{2} (v_{j}^{2h} + v_{j+1}^{2h}), \quad 0 \leq j \leq \frac{n}{2} - 1 \end{aligned}$$

$$I_{2h}^{h} \mathbf{v}^{2h} = \frac{1}{2} \begin{bmatrix} 1 & & \\ 2 & & \\ 1 & 1 & \\ & 2 & \\ & 1 & 1 \\ & & 2 \\ & & & 1 \end{bmatrix} \cdot \begin{bmatrix} v_1 \\ v_2 \\ v_3 \end{bmatrix}_{2h} = \begin{bmatrix} v_1 \\ v_2 \\ v_3 \\ v_4 \\ v_5 \\ v_6 \\ v_7 \end{bmatrix}_{h} = \mathbf{v}^{h}$$

#### **Multiple Grids**

- Intergrid Transfer fine  $\rightarrow$  coarse: **Restriction** 
  - $\Omega^h \rightarrow \Omega^{2h}$ , "Downsampling"
  - Simplest method: Injection
  - Better: Full weighting

$$v_j^{2h} = \frac{1}{4}(v_{2j-1}^h + 2v_{2j}^h + v_{2j+1}^h), \quad 1 \le j \le \frac{n}{2} - 1$$

• Restriction operator:

$$I_{h}^{2h} = \frac{1}{4} \begin{bmatrix} 1 & 2 & 1 & & \\ & 1 & 2 & 1 & \\ & & 1 & 2 & 1 \end{bmatrix}$$
$$\mathbf{v}^{2h} = I_{h}^{2h} \mathbf{v}^{h}$$

• Transfer Operations  $\Omega^h \leftrightarrow \Omega^{2h}$  sufficient

- Aliasing: Oscillatory modes on Ω<sup>h</sup> will be represented as smooth modes on Ω<sup>2h</sup>
- A basic two-grid correction scheme
  - On grid  $\Omega^h$ , relax  $\upsilon_1$  times on  $A^h \mathbf{v}^h = \mathbf{0}$  with initial guess  $\mathbf{v}^{(0)h}$ 
    - Restrict fine-grid residual **r**<sup>h</sup> to the coarse grid
    - On grid  $\Omega^{2h}$ , relax  $\upsilon_2$  times on  $A^{2h}e^{2h} = r^{2h}$  with initial guess  $e^{(0)h} = 0$
    - Interpolate the coarse-grid error
  - Correct the fine-grid approximation:  $\mathbf{v}^{h} \leftarrow \mathbf{v}^{h} + \mathbf{e}^{h}$
  - On grid  $\Omega^h$ , relax  $\upsilon_1$  times on  $A^h \mathbf{v}^h = \mathbf{0}$  with initial guess  $\mathbf{v}^h$

- Multigrid strategies
  - Nested iteration: Use coarse grids to generate improved initial guesses
  - Coarse grid correction: Approximate the error by relaxing on the residual equation on a course grid



- The V-Cycle Scheme (Coarse Grid Correction)
  - V-Cycle(v<sup>h</sup>, f<sup>h</sup>)
    - Relax  $v_1$  times on  $A^h v^h = 0$  with initial guess  $v^h$
    - If (current grid = coarsest grid) goto last point
      - Else:  $\mathbf{f}^{2h} = \text{Restrict}(\mathbf{f}^h A^h \mathbf{v}^h)$
      - $v^{2h} = 0$
      - Call v<sup>2h</sup> = V-Cycle(v<sup>2h</sup>, f<sup>2h</sup>)
    - Correct v<sup>h</sup> += Interpolate(v<sup>2h</sup>)
    - Relax  $v_2$  times on  $A^h v^h = 0$  with initial guess  $v^h$
  - Recursive algorithm



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Performance

- Storage requirements
  - Vectors v and f for n = 16 with boundary values

V	Ω <sup>h</sup>	$\Omega^{2h}$	$\Omega^{4h}$	$\Omega^{8h}$
	17	9	5	3
f	Ω <sup>h</sup>	$\Omega^{2h}$	$\Omega^{4h}$	$\Omega^{8h}$
	17	9	5	3

StorageSpace  $< \frac{2n^d}{1-2^{-d}}$ , with d: Dimension of problem

 For d = 1, memory requirement is less then twice that of the fine-grid problem alone

- Computational costs
  - 1 work unit (WU): one relaxation sweep on Ω<sup>h</sup>
  - O(WU) = O(N), with N: Total number of grid points
  - Intergrid transfer is neglected
  - One relaxation sweep per level ( $v_i = 1$ )

$$Cost_{\text{V-Cycle}} < \frac{2}{1 - 2^{-d}} WU$$
$$Cost_{\text{FMG Cycle}} < \frac{2}{(1 - 2^{-d})^2} WU$$

 1D problem: Single V-Cycle costs ~4WU, Complete FMG cycle ~8WU

- Diagnostic Tools
  - Help to debug your implementation
  - Methodical Plan for testing modules
  - Starting Simply with small, simple problems
  - *Exposing Trouble* difficulties might be hidden
  - Fixed Point Property relaxation may not change exact solution
  - Homogenous Problem: norm of error and residual should decrease to zero

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- The Impact of Intergrid Transfer and the iterative method may be expressed and proven in a formal way
- Two-grid correction TG consists of matrices for Interpolation, Restriction and Relaxation
- Spectral picture of multigrid
  - Relaxation damps oscillator modes
  - Interpolation & Restriction damp smooth modes
- Algebraic picture of multigrid
  - Decompose Space of the error:  $\Omega^h = R \oplus N$
  - $R = Range(I_{2h}^{h}) = Ker(TG)$   $N = N(I_{h}^{2h}A^{h}) = Ker(I_{h}^{2h}A^{h})$
  - L similar to R, H similar to N

#### Theory

- Operations of multigrid, visualized
  - Plane represents Ω<sup>h</sup>
  - Error e<sup>h</sup> is successively projected on one of the axes
    - Relaxations on the fine grid (1)
    - Two-grid correction (2)
    - Again, relaxation on the fine grid (3)



# Thanks for your attentionAny questions?