

Numerical Methods in Scientific Computing

NGSSC

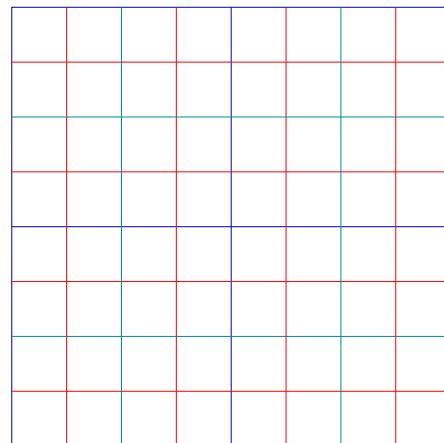
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Multigrid/Algebraic Multigrid methods

Run Jacobi demo...

Main idea:

reduce the error $e^{(k)} = \mathbf{x}_{exact} - \mathbf{x}^{(k)}$ on the given (fine) grid by successive residual corrections on a hierarchy of (nested) coarser grids.



1966, N.S. Bakhvalov and R. Fedorenko

Multilevel preconditioning methods: MG

Procedure MG : $\mathbf{u}^{(k)} \leftarrow MG \left(\mathbf{u}^{(k)}, \mathbf{f}^{(k)}, k, \{\nu_j^{(k)}\}_{j=1}^k \right)$;

if $k = 0$, **then** solve $A^{(0)} \mathbf{u}^{(0)} = \mathbf{f}^{(0)}$ exactly or by smoothing,

else

$\mathbf{u}^{(k)} \xleftarrow{s_1} \mathcal{S}_1^{(k)} (\mathbf{u}^{(k)}, \mathbf{f}^{(k)})$, perform s_1 pre-smoothing steps,

Correct the residual:

$\mathbf{r}^{(k)} = A^{(k)} \mathbf{u}^{(k)} - \mathbf{f}^{(k)}$; form the current residual,

$\mathbf{r}^{(k-1)} \leftarrow \mathcal{R} (\mathbf{r}^{(k)})$, restrict the residual on the next coarser grid,

$\mathbf{e}^{(k-1)} \leftarrow MG \left(\mathbf{0}, \mathbf{r}^{(k-1)}, k-1, \{\nu_j^{(k-1)}\}_{j=1}^{k-1} \right)$;

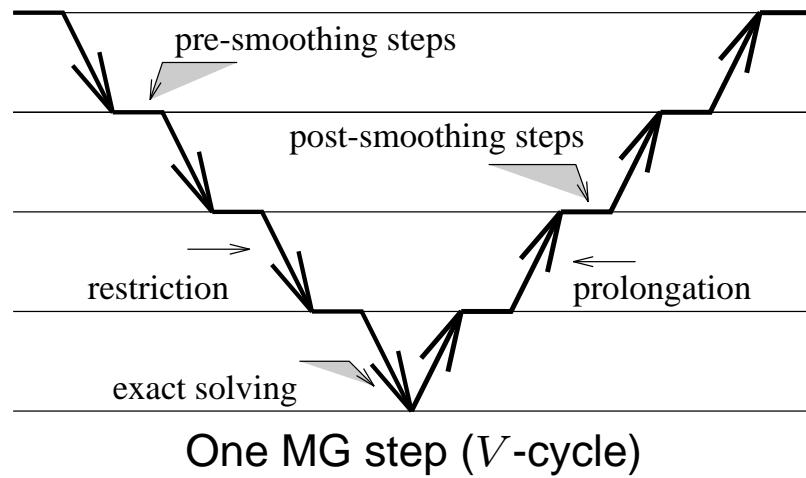
$\mathbf{e}^{(k)} \leftarrow \mathcal{P} (\mathbf{e}^{(k-1)})$; prolong the error from the next coarser to the current grid,

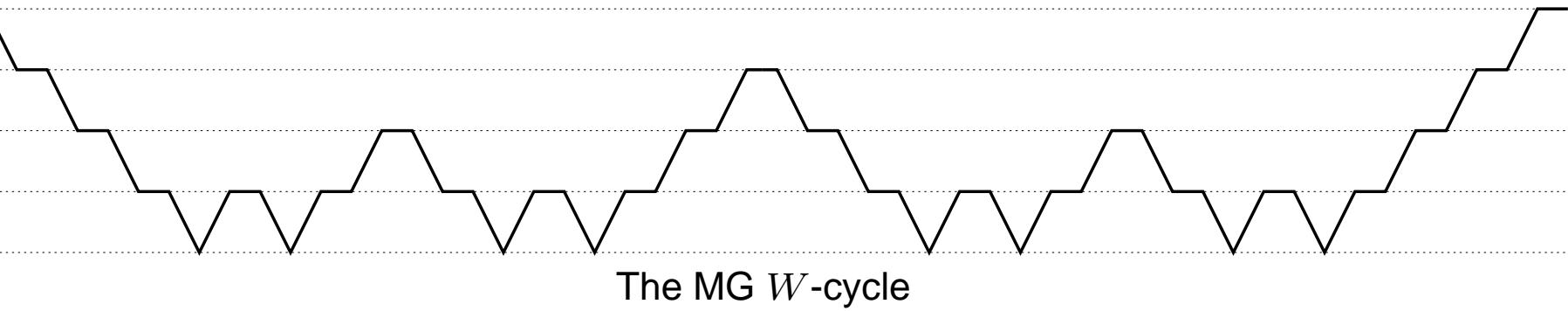
$\mathbf{u}^{(k)} = \mathbf{u}^{(k)} - \mathbf{e}^{(k)}$; update the solution,

$\mathbf{u}^{(k)} \xleftarrow{s_2} \mathcal{S}_2^{(k)} (\mathbf{u}^{(k)}, \mathbf{f}^{(k)})$, perform s_2 post-smoothing steps.

endif

end Procedure MG



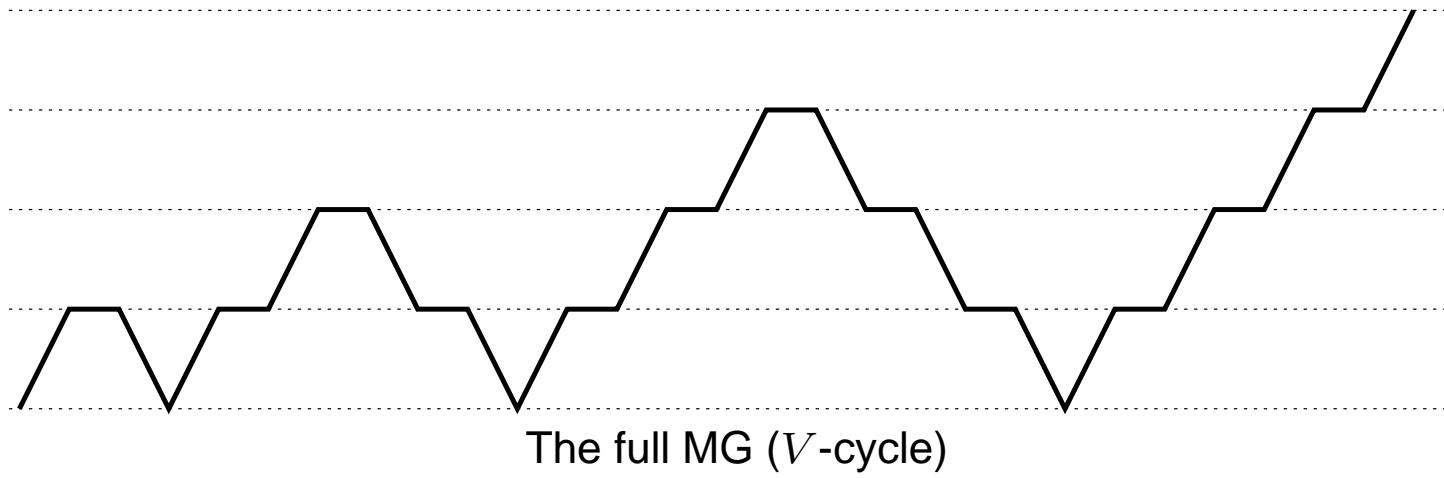


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Procedure NI:  $\mathbf{u}^{(\ell)} \leftarrow NI \left( \mathbf{u}^{(0)}, \{\mathbf{f}^{(k)}\}_{k=1}^{(\ell)}, \ell, \{\nu^{(k)}\}_{k=1}^{\ell} \right);$ 
 $\mathbf{u}^{(0)} = A^{(0)}^{-1} \mathbf{f}^{(0)},$ 
for  $k=1$  to  $\ell$  do
     $\mathbf{u}^{(k)} = \mathcal{P} (\mathbf{u}^{(k-1)});$ 
     $\mathbf{u}^{(k)} \leftarrow MG \left( \mathbf{u}^{(k)}, \mathbf{f}^{(k)}, k, \{\nu_j^{(k)}\}_{j=1}^k \right);$ 
endfor
end Procedure NI

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The so-called *full MG* corresponds to **Procedure** $NI(\cdot, \cdot, \ell, \{1, 1, \dots, 1\})$



A compact formula presenting the MG procedure in terms of a recursively defined iteration matrix:

- (i) Let $M^{(0)} = 0$,
- (ii) For $k = 1$ to ℓ , define

$$M^{(k)} = S^{(k)^{s_2}} \left(A^{(k)^{-1}} - P_{k-1}^k \left(I - M^{(k-1)^{\nu}} \right) A^{(k-1)^{-1}} R_k^{k-1} \right) A^{(k)} S^{(k)^{s_1}},$$

where $S^{(k)}$ is a smoothing iteration matrix (assuming S_1 and S_2 are the same), R_k^{k-1} and P_{k-1}^k are matrices which transfer data between two consecutive grids and correspond to the restriction and prolongation operators \mathcal{R} and \mathcal{P} , respectively, and $\nu = 1$ and $\nu = 2$ correspond to the V - and W -cycles.

It turns out that in many cases the spectral radius of $M^{(\ell)}$, $\rho(M^{(\ell)})$, is independent of ℓ , thus the rate of convergence of the NI method is optimal. Also, a mechanism to make the spectral radius of $M^{(\ell)}$ smaller is to choose s_1 and s_2 larger. The price for the latter is, clearly, a higher computational cost.

Rate of convergence

Multigrid demo ...

Procedure $AMLI$: $\mathbf{u}^{(k)} \leftarrow AMLI \left(\mathbf{f}^{(k)}, k, \nu_k, \{a_j^{(k)}\}_{j=0}^{\nu_k} \right);$

$[\mathbf{f}_1^{(k)}, \mathbf{f}_2^{(k)}] \leftarrow \mathbf{f}^{(k)},$

$\mathbf{w}_1^{(k)} = B_{11}^{(k)} \mathbf{f}_1^{(k)},$

$\mathbf{w}_2^{(k)} = \mathbf{f}_2^{(k)} - A_{21}^{(k)} \mathbf{w}_1^{(k)},$

$k = k - 1,$

if $k = 0$ **then** $\mathbf{u}_2^{(0)} = A^{(0)} \mathbf{w}_2^{(1)}$, solve on the coarsest level exactly;

else

$\mathbf{u}_2^{(k)} \leftarrow AMLI \left(a_{\nu_k}^{(k)} \mathbf{w}_2^{(k)}, k, \nu_k, \{a_j^{(k)}\}_{j=0}^{\nu_k} \right);$

for $j = 1$ **to** $\nu_k - 1$:

$\mathbf{u}_2^{(k)} \leftarrow AMLI \left(A^{(k)} \mathbf{u}_2^{(k)} + a_{\nu_k-j}^{(k)} \mathbf{w}_2^{(k)}, k, \nu_k, \{a_j^{(k)}\}_{j=0}^{\nu_k} \right);$

endfor

endif

$k = k + 1,$

$\mathbf{u}_1^{(k)} = \mathbf{w}_1^{(k)} - B_{11}^{(k)} A_{12}^{(k)} \mathbf{u}_2^{(k)},$

$\mathbf{u}^{(k)} \leftarrow [\mathbf{u}_1^{(k)}, \mathbf{u}_2^{(k)}]$

end Procedure $AMLI$