

THE INCOMPLETE LU-DECOMPOSITION AS A
RELAXATION METHOD IN MULTI-GRID
ALGORITHMS

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ABSTRACT

We consider relaxation methods for use with Multi-Grid (MG-) algorithms to solve a sparse linear system

$$A_\ell x_\ell = f_\ell, \quad \ell > 0,$$

which is supposed to be the discretization of a continuous boundary value problem $Ax = f$. With the same equation coarser discretizations are related:

$$A_k x_k = f_k, \quad k = 0, 1, \dots, \ell-1.$$

We first give a brief exposition of the framework of MG-methods. Next, we describe the incomplete LU-decomposition as a relaxation method and in the 3rd section we compare different variants of it. The conclusion is in favour of the ILU variant.

THE MULTI-GRID FRAMEWORK

Considering MG methods in which each iteration step consists of γ coarse grid correction steps, preceded by p and followed by q relaxation sweeps, we see that in each MG iteration step the residual is multiplied by the operator (cf. [2])

$$M_k^{-MLA} = M_k^{-TLA} + (M_k^{-REL})^q A_k P_{k-1}^k A_{k-1}^{-1} (M_{k-1}^{-TLA})^\gamma R_k^{k-1} (M_k^{-REL})^p,$$

where

$$M_0^{-TLA} = 0$$

$$M_k^{-TLA} = (M_k^{-REL})^q (I - A_k P_{k-1}^k A_{k-1}^{-1} R_k^{k-1}) (M_k^{-REL})^p, \quad k = 1, 2, \dots, \ell.$$

P_{k-1}^k denotes the prolongation from level $k-1$ to level k ,

R_k^{k-1} denotes the restriction from level k to level $k-1$,

A_{k-1} denotes the $k-1$ level discretisation of the operator A , e.g.

$$A_{k-1} = R_k^{k-1} A_k P_{k-1}^k.$$

M_k^{-REL} is the operator on level k by which the residual is multiplied in one relaxation sweep (see next section).

Communication between finer and coarser grids takes place via the prolongations and restrictions. By the spectral decomposition of a gridfunction

$$u_h(jh) = \left(\frac{1}{\sqrt{2\pi}}\right)^n \int_{\omega \in [-\pi/h, \pi/h]^n} e^{+ijh\omega} \hat{u}_h(\omega) d\omega$$

where

$$\hat{u}_h(\omega) = \left(\frac{h}{\sqrt{2\pi}}\right)^n \sum_{j \in \mathbb{Z}^n} e^{-ijh\omega} u_h(jh)$$

denotes the spectrum of the gridfunction u_h , defined on $[-\pi/h, \pi/h]^n$, we see that only low frequency components of a gridfunction can be represented on coarse grids. The relation between the spectra of a gridfunction, its prolongation and its restriction are given by ($H = qh$)

$$\widehat{Ru}_h(\omega) = \left(\frac{\sqrt{2\pi}}{h}\right)^n \sum_{0 \leq p < q} \hat{a}_h(\omega + 2\pi p/H) \hat{u}_h(\omega + 2\pi p/H),$$

$$\widehat{Pu}_h(\omega) = \left(\frac{\sqrt{2\pi}}{H}\right)^n \hat{b}_h(\omega) \hat{u}_H(\omega);$$

where a_h and b_h are grid functions that characterize the particular prolongations and restrictions (cf. [3]).

With the choice $A_{k-1} = RA_kP$, we find the relation

$$(I - \tilde{P}R)(I - A_k P A_{k-1}^{-1} R) = (I - A_k P A_{k-1}^{-1} R)$$

for any prolongation \tilde{P} . Hence, taking e.g. Shannon's interpolation for \tilde{P} , we see that - in the residual - coarse grid corrections annihilate all frequencies that can be represented on the coarser grid. Similar results hold for other reasonable choices of A_{k-1} .

From this it is clear that an efficient relaxation method in an MG-algorithm should damp those high frequencies that cannot be represented on the coarser grid (in the 2-D case ($n=2$) and with mesh doubling ($H=qh=2h$) this is the shadowed portion of fig. 1). This behaviour is analyzed by considering $\hat{M}(\omega)$, the spectrum of the operator \tilde{M}_k^{REL} . Following BRANDT [1], we define the smoothing rate

$$\mu = \sup_{\substack{\omega \in [-\pi/h, \pi/h]^n \\ \omega \notin [-\pi/2h, \pi/2h]^n}} |\hat{M}(\omega)|.$$

For any linear difference operator A_k with constant coefficients and each relaxation method this μ can easily be determined numerically. E.g. it is well known that for the usual 5-point discretization of Poisson's equation ($n=2, q=2$) and GS-relaxation we have $\mu = 0.5$.

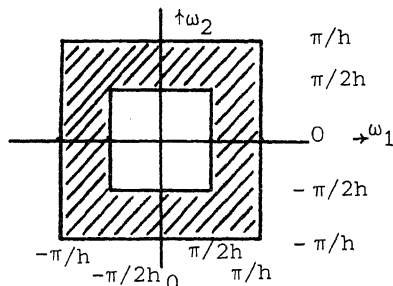


Fig. 1. Frequency region

We consider similar relaxation methods in order to find methods that take less work per iteration sweep and smaller values of μ .

INCOMPLETE LU-DECOMPOSITION AS A RELAXATION METHOD

By B_k we denote the approximate inverse in a stationary relaxation method for the solution of $A_k x_k = f_k$. Then $M_k^{REL} = I - A_k B_k$ is the operator by which the residual is multiplied in each relaxation sweep. E.g. with Jacobi-iteration $B_k = (\text{diag}(A_k))^{-1}$ and with forward - or backward Gauss-Seidel iteration

$$B_k = (\text{lower triag } (A_k))^{-1} \quad \text{resp.} \quad B_k = (\text{upper triag } (A_k))^{-1}.$$

We are interested in incomplete LU-decomposition, where for some D , L and U (which are diagonal, lower and upper triangular respectively) we take

$$B_k = (LD^{-1}U)^{-1}$$

and we may write

$$A_k = LD^{-1}U + R.$$

Although the treatment of incomplete LU-decomposition can be given for the general case, here we confine ourselves to infinite Toeplitz-matrices A_k, L, D and U of the form: $D = wI$,

$$(*) \quad A_k = \begin{pmatrix} C & G & & E \\ & B & & \\ & & & \\ H & & & \end{pmatrix}, \quad L = \begin{pmatrix} a & & & \\ b & & & \\ & c & & \\ & & & \end{pmatrix}, \quad U = \begin{pmatrix} \alpha & \beta & & \\ & & \gamma & \\ & & & \\ & & & \end{pmatrix}.$$

It can be shown that the analysis for this case is representative for the local behaviour in the interior of the domain of the PDE. Multiplication of $LD^{-1}U$ and identification of the sub- and superdiagonals yields

$$c = H, \quad b = B, \quad \beta = G, \quad \gamma = E, \quad a = \alpha = w^{-1},$$

and R has the form

$$R = \begin{pmatrix} rm & 0 & & ru & 0 & \\ 0 & & & & & \\ & & & & & \\ & & & & & \\ rl & & & & & \\ 0 & & & & & \end{pmatrix}, \quad \begin{aligned} ru &= BE/a, \\ rl &= HG/a, \\ rm &= a - C + (GD+HE)/a. \end{aligned}$$

Here one free parameter a is left. If no corrections are admitted to the main diagonal: $rm = 0$ and a is prescribed by

$$(ILU) \quad a = \frac{1}{2}(C \pm \sqrt{C^2 - 4(GD+HE)}).$$

To get minimal ru and rl the sign which yields maximal absolute value is used. Other choices for a are possible:

$$(SGS) \quad a=C,$$

or the 1-st order approximation

$$(MILU) \quad a=C - (GD+HE)/C.$$

Remark.

Notice that the choice $a = C$ yields a method which is equivalent with symmetric Gauss-Seidel relaxation. In this case namely $A_k = L + U - D$ and we see

$$\bar{M}^{\text{REL}} = I - A_k(LD^{-1}U)^{-1} = (I - A_k U^{-1})(I - A_k L^{-1}),$$

i.e. one GS-forward step followed by one GS-backward step is identical with this particular incomplete LU-iteration step. *Symmetric GS-iteration (SGS) is one (simple) form of incomplete LU-iteration.*

THE EFFICIENCY OF INCOMPLETE LU-ITERATION

To study the efficiency of the incomplete LU-relaxation, first we compute the smoothing rate $\mu(a)$ for the usual 5-point discretization of Poisson's equation, depending on the parameter a . The maximum of

$$\hat{M}(\omega) = \frac{(a^2 - 4a + 2) + 2 \cos(\omega_1 - \omega_2)}{(a^2 + 2) - 2a(\cos \omega_1 - \cos \omega_2) + 2 \cos(\omega_1 - \omega_2)}$$

on the shaded area of fig. 1 can be computed analytically and is ($a > 2$)

$$\mu(a) = \frac{8 \pm A\sqrt{A^2 + 12}}{8 - 2Aa \pm (A+4a)\sqrt{A^2+12}}, \quad \text{where } A = a^2 - 4a + 2,$$

and is attained at $\omega = (\arccos(\frac{8 + A\sqrt{A^2 + 12}}{A^2 + 16}), \frac{\pi}{2})$.

A graph of this function is given in fig. 2. For our three particular choices of a we find:

$$\begin{aligned} \text{SGS} : a = 4 & , \quad \mu(a) = 0.2500, \quad \omega = (\arccos(4/5), \pi/2); \\ \text{MILU} : a = 3.5 & , \quad \mu(a) = 0.1649, \quad \omega = (0.555\pi, \pi/2); \\ \text{ILU} : a = 2 + \sqrt{2} & , \quad \mu(a) = 0.2035, \quad \omega = (\pi/3, \pi/2). \end{aligned}$$

The minimal value of $\mu(a)$ is 0.1607 and is attained for $a = 3.510$ which is remarkably close to our choice MILU.

The amount of work in each iteration step is $5N$ for the solution and $3N$ for the computation of the residual ($2N$ in the case ILU, where $rm = 0$). Summarizing, the efficiency of the different methods is

$$\begin{aligned} \text{SGS} : \mu = 0.25, \text{ work } 8N, \text{ efficiency } \frac{\sqrt[3]{0.25}}{\sqrt[3]{0.25}} = 0.84; \\ \text{MILU} : \mu = 0.16, \text{ work } 8N, \text{ efficiency } \frac{\sqrt[3]{0.16}}{\sqrt[3]{0.16}} = 0.80; \\ \text{ILU} : \mu = 0.20, \text{ work } 7N, \text{ efficiency } \frac{\sqrt[3]{0.20}}{\sqrt[3]{0.20}} = 0.79. \end{aligned}$$

For the Poisson equation we conclude that, although MILU has optimal $\mu(a)$, ILU has slightly better efficiency because of the cheaper computation of the residual.

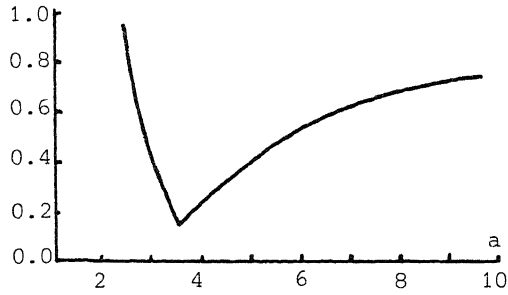


Fig. 2 The smoothing rate $\mu(a)$.

Beside Poisson's equation we also considered other model equations viz. the convection-diffusion equation

$$\varepsilon \Delta \phi + \cos(\alpha) \phi_x + \sin(\alpha) \phi_y = f,$$

discretized with Il'in's method and the anisotrope Poisson equation

$$\varepsilon \phi_{xx} + \phi_{yy} = f.$$

For the convection diffusion we see from (*) that, asymptotically for $\varepsilon \rightarrow 0$, either H or E and either B or G vanish. Hence, asymptotically the methods SGS, MILU and ILU coincide. Moreover, for $0 < \alpha < \pi/2$, $\text{mod}(\pi)$, asymptotically $\mu \rightarrow 0$, i.e. the system is solved exactly by only one iteration sweep. However, for a convection direction α with $-\pi/2 \leq \alpha \leq 0$, $\text{mod}(\pi)$, μ takes a positive value. For different values of ε and α , these μ are given in table 1. Taking into account the number of operations, we see again that, although $\mu_{\text{MILU}} \leq \mu_{\text{ILU}} \leq \mu_{\text{SGS}}$, ILU is in most cases more efficient than MILU.

$\varepsilon = 1.0$	α :				
	$-\pi/4$	$-\pi/8$ $-3\pi/8$	$\alpha = 0$ $\pi/2$	$\pi/8$ $3\pi/8$	$\alpha = \pi/4$
SGS	0.261	0.258	0.243	0.225	0.215
MILU	0.200	0.195	0.174	0.153	0.144
ILU	0.229	0.226	0.205	0.182	0.171
$\varepsilon = 0.1$					
SGS	0.446	0.655	0.405	0.128	0.679
MILU } ILU }	0.447	0.660	0.427	0.131	0.488
$\varepsilon = 0.001$					
SGS } MILU } ILU }	0.447	0.679	0.499	0.000	0.000

Table 1. The smoothing rate μ for the convection diffusion equation with Il'in's discretization.

	$\epsilon = 1$	$\epsilon = 0.5$	$\epsilon = 0.1$	$\epsilon = 0.01$	$\epsilon = 0.001$
SGS	0.250	0.321	0.697	0.961	0.996
MILU	0.165	0.190	0.387	0.889	0.988
ILU	0.204	0.250	0.477	0.768	0.916

Table 2. μ for the anisotrope Poisson equation.

For the anisotrope Poisson equation (table 2) we see that in the three cases $\mu \rightarrow 1$ as $\epsilon \rightarrow 0$. However, MILU and ILU have smaller smoothing rates and for small ϵ we even have $\mu_{ILU} < \mu_{MILU}$.

We conclude that also in this case the ILU-decomposition is the best choice in our class of incomplete LU-decomposition relaxation methods and it is significantly more efficient than symmetric Gauss-Seidel relaxation.

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