

Numerical Solutions to PDEs

Lecture Notes #23

Elliptic Equations — Steepest Descent and Conjugate Gradient

Peter Blomgren,
(blomgren.peter@gmail.com)

Department of Mathematics and Statistics
Dynamical Systems Group
Computational Sciences Research Center
San Diego State University
San Diego, CA 92182-7720
<http://terminus.sdsu.edu/>

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Last Time: Linear Iterative Schemes

We looked at the Jacobi, Gauss-Seidel, SOR, and SSOR iterations applied to linear systems $A\bar{\mathbf{x}} = \bar{\mathbf{b}}$, originating from the 5-point Laplacian.

We quantified under what circumstances we can guarantee convergence of these iterations (J&GS: irreducibly diagonally dominant matrices, (S)SOR: $\omega \in (0, 2)$), and discussed the convergence rates.

The discussion was extended to general linear systems, where A may be associated with the 9-point Laplacian, or something completely different. In this discussion we introduced **preconditioning**, where we find a matrix $M \approx A$, which is much easier to invert than A itself, and we leverage this in order to generate an efficient iterative solver.



Outline

- 1 Recap
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- 2 A Different Point of View
 - $A\bar{\mathbf{x}} = \bar{\mathbf{b}}$ as an Optimization Problem
 - Steepest Descent
 - Conjugate Gradient
- 3 Beyond Conjugate Gradient
 - Preconditioning, and Extensions
 - Beyond Finite Differences...



Another Point of View: Optimization

We consider a system of linear equations $A\bar{\mathbf{x}} = \bar{\mathbf{b}}$, where A is symmetric positive definite.

We define

$$F(\bar{\mathbf{y}}) = \frac{1}{2}(\bar{\mathbf{y}} - \bar{\mathbf{x}})^T A(\bar{\mathbf{y}} - \bar{\mathbf{x}}),$$

and note that since A is positive definite $F(\bar{\mathbf{y}}) \geq 0$, and $F(\bar{\mathbf{y}}) = 0 \Leftrightarrow \bar{\mathbf{y}} = \bar{\mathbf{x}}$. Further, we can define

$$E(\bar{\mathbf{y}}) = F(\bar{\mathbf{y}}) - F(\bar{\mathbf{0}}) = \frac{1}{2}\bar{\mathbf{y}}^T A\bar{\mathbf{y}} - \bar{\mathbf{y}}^T \bar{\mathbf{b}},$$

which has a unique minimum at the solution of $A\bar{\mathbf{x}} = \bar{\mathbf{b}}$.

Now the gradient of $E(\bar{\mathbf{y}})$ describes the direction of largest increase

$$G(\bar{\mathbf{y}}) = \nabla E(\bar{\mathbf{y}}) = A\bar{\mathbf{y}} - \bar{\mathbf{b}} = -\underbrace{\bar{\mathbf{r}}(\bar{\mathbf{y}})}_{\text{residual}}.$$



Optimization \rightsquigarrow Steepest Descent

Since the gradient points in the direction of steepest ascent, the residual points in the direction of steepest descent.

Given an approximation (guess) \bar{x}^k to the solution of $A\bar{x} = \bar{b}$, we find a better approximation by searching in the steepest descent direction

$$\bar{x}^{k+1} = \bar{x}^k + \alpha_k \bar{r}^k, \quad \text{where } \bar{r}^k = \bar{b} - A\bar{x}^k,$$

and we select α_k so that $E(\bar{x}^k + \alpha_k \bar{r}^k)$ is minimized:

$$\begin{aligned} E(\bar{x}^{k+1}) &= E(\bar{x}^k + \alpha_k \bar{r}^k) \\ &= \frac{1}{2} [\bar{x}^k]^T A \bar{x}^k + \alpha_k [\bar{r}^k]^T A \bar{x}^k + \frac{1}{2} \alpha_k^2 [\bar{r}^k]^T A \bar{r}^k - [\bar{x}^k]^T \bar{b} - \alpha_k [\bar{r}^k]^T \bar{b} \\ &= E(\bar{x}^k) - \alpha_k [\bar{r}^k]^T \bar{r}^k + \frac{1}{2} \alpha_k^2 [\bar{r}^k]^T A \bar{r}^k. \end{aligned}$$

Setting $\partial E(\bar{x}^k + \alpha_k \bar{r}^k) / \partial \alpha_k = 0$ gives us

$$\alpha_k = \frac{[\bar{r}^k]^T \bar{r}^k}{[\bar{r}^k]^T A \bar{r}^k} = \frac{\|\bar{r}^k\|_2^2}{[\bar{r}^k]^T A \bar{r}^k}.$$



Steepest Descent

The steepest descent algorithm is given by $\bar{x}^0 = \bar{0}$, $\bar{r}^0 = \bar{b}$:

$$\begin{aligned} \alpha_k &= \frac{\|\bar{r}^k\|^2}{[\bar{r}^k]^T A \bar{r}^k} \\ \bar{x}^{k+1} &= \bar{x}^k + \alpha_k \bar{r}^k \\ \bar{r}^{k+1} &= \bar{r}^k - \alpha_k A \bar{r}^k. \end{aligned}$$

Where the update formula for the residual comes from

$$\begin{aligned} \bar{x}^{k+1} &= \bar{x}^k + \alpha_k \bar{r}^k \\ A\bar{x}^{k+1} &= A\bar{x}^k + \alpha_k A \bar{r}^k \\ \bar{b} - A\bar{x}^{k+1} &= \bar{b} - A\bar{x}^k - \alpha_k A \bar{r}^k \\ \bar{r}^{k+1} &= \bar{r}^k - \alpha_k A \bar{r}^k. \end{aligned}$$



Steepest Descent

We note that the steepest descent algorithm only requires one matrix-vector product $A\bar{r}^k$ and two vector-vector inner products ($\|\bar{r}^k\|^2$, $[\bar{r}^k]^T A \bar{r}^k$) per iteration.

When A is sparse the matrix-vector product can be implemented in $\mathcal{O}(N)$ operations.

Theorem

If A is a positive definite matrix for which $A^T A^{-1}$ is also positive definite, then the steepest descent algorithm converges to the unique solution $\bar{x}^ = A^{-1} \bar{b}$ for any initial \bar{x}^0 .*

Theorem

If A is SPD, then the steepest descent algorithm converges to the unique solution $\bar{x}^ = A^{-1} \bar{b}$ for any initial \bar{x}^0 .*



Steepest Descent

It turns out, maybe somewhat counter-intuitively, that the steepest descent algorithm converges very slowly unless A is a (near-)multiple of the identity matrix.

The residuals tend to oscillate so that \bar{r}^{k+2} points in the same direction as \bar{r}^k , and very little progress is made.

Next we quantify this convergence rate, and discuss the **conjugate gradient method** which is an “accelerated version of steepest descent.”



Convergence Rate for Steepest Descent

Theorem (Convergence Rate for Steepest Descent)

If A is a symmetric positive definite matrix whose eigenvalues lie in the interval $[a, b]$, then the error vector \bar{e}^k for the steepest descent method satisfies

$$[\bar{e}^k]^T A \bar{e}^k \leq \left[\frac{b-a}{b+a} \right]^{2k} [\bar{e}^0]^T A \bar{e}^0 \equiv \left[\frac{\kappa-1}{\kappa+1} \right]^{2k} [\bar{e}^0]^T A \bar{e}^0$$

The larger the interval $[a, b]$, i.e. the more ill-conditioned A is, the slower the convergence rate we get.

The **condition number** κ of a matrix is defined as

$$\kappa = \frac{b}{a} = \frac{|\lambda|_{\max}}{|\lambda|_{\min}}$$

it is an intrinsic measure of difficult the matrix is to invert.



The Steepest Descent Method: Zig-Zagging

The “zig-zagging” ($\bar{p}^{k+2} \approx \bar{p}^k$) is what causes the steepest descent method to slow down. The amount of zig-zagging is directly proportional to the ratio $|\lambda|_{\max}/|\lambda|_{\min}$, or more generally for a non-square matrix A , $\sigma_{\max}/\sigma_{\min}$, where σ_{ν} are the singular values of A .

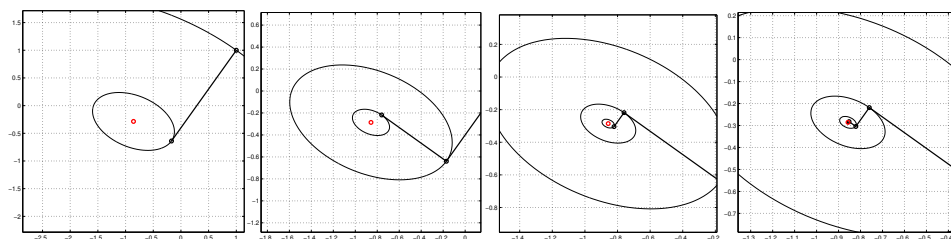


Figure: Illustration of the “zig-zagging” of the search directions in the steepest descent algorithm. If $\kappa = 1$, then all the level curves of $\|A\bar{x} - \bar{b}\| = c$ are circles (hyper-spheres in \mathbb{R}^n) and the steepest descent direction points straight in toward the central point. The more elongated the ellipse becomes, the more zig-zagging we get...



The Conjugate Gradient Method

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The Conjugate Gradient method can be viewed as an acceleration of the steepest descent method, in which we by adding a little bit of “memory” to the algorithm can avoid the zig-zagging.

We consider

$$\bar{x}^{k+1} = \bar{x}^k + \alpha_k \underbrace{\left[\bar{r}^k + \gamma_k \underbrace{(\bar{x}^k - \bar{x}^{k-1})}_{\alpha_{k-1} \bar{p}^{k-1}} \right]}_{\bar{p}^k}$$

clearly, if $\gamma_k \equiv 0$, we can recover the steepest descent algorithm.

We form the new search direction \bar{p}^k as a linear combination of the steepest descent direction \bar{r}^k and the previous search direction \bar{p}^{k-1} , i.e.

$$\bar{p}^k = \bar{r}^k + \beta_{k-1} \bar{p}^{k-1}.$$



The Conjugate Gradient Method

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The conjugate gradient iteration involves updates for the approximate solution \bar{x} , the residual \bar{r} , and the search direction \bar{p} :

$$\begin{aligned} \bar{x}^{k+1} &= \bar{x}^k + \alpha_k \bar{p}^k, \\ \bar{r}^{k+1} &= \bar{r}^k - \alpha_k A \bar{p}^k, \\ \bar{p}^{k+1} &= \bar{r}^{k+1} + \beta_k \bar{p}^k. \end{aligned}$$

Where we want to select α_k and β_k in an optimal way. A minimization of the error $E(\bar{x}^{k+1})$ with respect to α (just as in the steepest descent case), and a similar analysis of $E(\bar{x}^{k+1})$ with respect to β gives

$$\alpha_k = \frac{\|\bar{r}^k\|_2^2}{[\bar{p}^k]^T A \bar{p}^k}, \quad \beta_k = -\frac{[\bar{r}^{k+1}]^T A \bar{p}^k}{[\bar{p}^k]^T A \bar{p}^k} \equiv \frac{\|\bar{r}^{k+1}\|_2^2}{\|\bar{r}^k\|_2^2}.$$



The Conjugate Gradient Method

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Algorithm: The Conjugate Gradient Method

$$\bar{p}^0 = \bar{r}^0 = \bar{b} - A\bar{x}^0, \quad k = 0$$

while ($\|\bar{r}^k\| > \epsilon_{\text{tol}}\|\bar{r}^0\|$)

$$\alpha_k = \frac{\|\bar{r}^k\|_2^2}{[\bar{p}^k]^T A \bar{p}^k}$$

$$\bar{x}^{k+1} = \bar{x}^k + \alpha_k \bar{p}^k$$

$$\bar{r}^{k+1} = \bar{r}^k - \alpha_k A \bar{p}^k$$

$$\beta_k = \frac{\|\bar{r}^{k+1}\|_2^2}{\|\bar{r}^k\|_2^2}$$

$$\bar{p}^{k+1} = \bar{r}^{k+1} + \beta_k \bar{p}^k$$

endwhile ($k := k + 1$)



The Conjugate Gradient Method

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The CG method only requires one matrix-vector product $A\bar{p}^k$, and two vector-vector inner products $[\bar{p}^k]^T A \bar{p}^k$ and $\|\bar{r}^k\|_2^2$ per iteration, hence if A has $\mathcal{O}(N)$ non-zero entries, the work/iteration is $\mathcal{O}(N)$.

The CG gets its name (somewhat incorrectly, it should be “**the A-conjugate search-direction method**”) from the fact that the generated residuals are orthogonal, and the search directions are A -conjugate, *i.e.*

$$[\bar{r}^k]^T \bar{r}^j = [\bar{p}^k]^T A \bar{p}^j = 0, \quad \text{for } k \neq j.$$

A direct corollary of these (easily checked) facts, is

Corollary

If A is an $N \times N$ symmetric positive definite matrix, then the CG algorithm converges in at most N steps.



The Conjugate Gradient Method

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The N -step termination theorem tells us that for the 5-point Laplacian on an $N \times N$ grid we need at most

$$W_{\text{CG}} = \underbrace{5(N \times N)}_{\text{Matrix Entries}} \cdot \underbrace{N \times N}_{\text{iterations}} = \mathcal{O}(N^4),$$

operations to compute the exact solution to $A\bar{x} = \bar{b}$. This may not seem so impressive, since optimal SOR does a better job

$$W_{\text{SOR}}^* \approx \frac{N^3}{\pi^2} \log(\epsilon^{-1}) = \mathcal{O}(N^3).$$

However, in practice the iterates \bar{x}^k generated by the CG-iteration converge to \bar{x} very rapidly, and the iteration can be stopped for $k \ll N \times N$ iterations. Applied to the 5-point Laplacian, the CG iteration and optimal SOR both require $\sim N \log(\epsilon^{-1})$ iterations to reach a specified tolerance. CG has the advantage over SOR in that (i) *there is no parameter (ω) which must be optimally chosen*; further (ii) *the CG-iteration can be accelerated further by preconditioning PCG(M).*



Convergence Rate for the Conjugate Gradient Method

Theorem (Convergence Rate for Conjugate Gradient)

If A is a symmetric positive definite matrix whose eigenvalues lie in the interval $[a, b]$, then the error vector \bar{e}^k for the steepest descent method satisfies

$$[\bar{e}^k]^T A \bar{e}^k \leq \left[\frac{\sqrt{b} - \sqrt{a}}{\sqrt{b} + \sqrt{a}} \right]^{2k} [\bar{e}^0]^T A \bar{e}^0 \equiv \left[\frac{\sqrt{\kappa} - 1}{\sqrt{\kappa} + 1} \right]^{2k} [\bar{e}^0]^T A \bar{e}^0.$$



Conjugate Gradient vs. Steepest Descent

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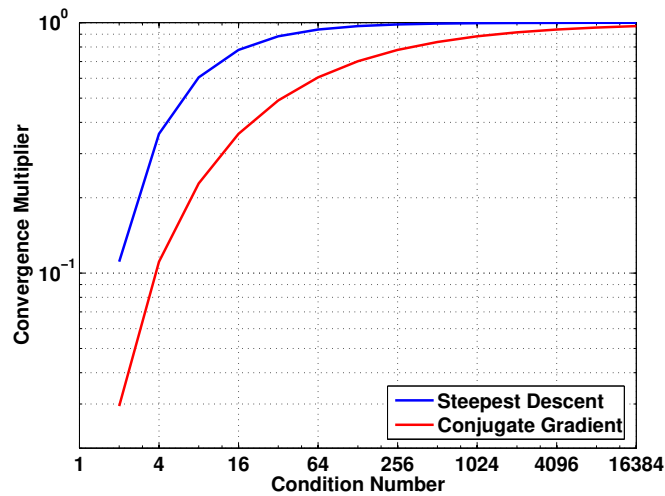


Figure: The convergence multipliers $m_{SD} = \left[\frac{\kappa-1}{\kappa+1}\right]^2$, and $m_{CG} = \left[\frac{\sqrt{\kappa}-\sqrt{1}}{\sqrt{\kappa}+\sqrt{1}}\right]^2$.



Conjugate Gradient vs. Steepest Descent

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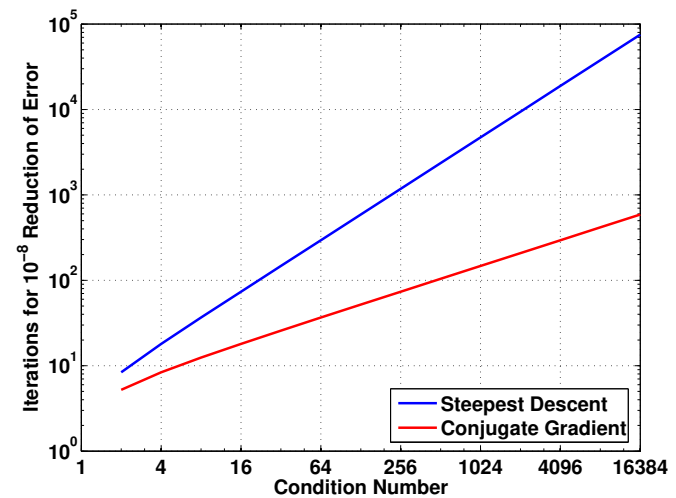


Figure: The number of iterations necessary to reduce the initial error by a factor of 10^{-8} .



GS vs. SOR vs. CG

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n	n^2	$\kappa(A)$	GS	SOR*	CG
8	64	47	252	65	11
16	256	169	837	121	27
32	1,024	641	2,870	223	52
64	4,096	2,489	9,983	414	98
128	16,384	9,807	34,706	777	192
256	65,536	38,926	—	1,473	370
512	262,144	155,103	—	2,813	715

Table: Number of iterations needed to achieve 10^{-8} relative update. 5-point Laplacian ∇_{5pt}^2 in 2D discretized on an $n \times n$ grid $\rightsquigarrow n^2 \times n^2$ matrix, with $\sim 5n^2$ non-zero elements.



GS vs. SOR vs. CG

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n	n^2	$\kappa(A)$	GS	SOR*	CG
8	64	47	—	71	10
16	256	169	—	136	28
32	1,024	641	—	261	59
64	4,096	2,489	—	504	119
128	16,384	9,807	—	984	239
256	65,536	38,926	—	1,938	470
512	262,144	155,103	—	3,844	941

Table: Number of iterations needed to achieve 10^{-8} residual reduction. 5-point Laplacian ∇_{5pt}^2 in 2D discretized on an $n \times n$ grid $\rightsquigarrow n^2 \times n^2$ matrix, with $\sim 5n^2$ non-zero elements.

Bottom Line: Even in the “homework case” where the optimal SOR parameter is known, the Conjugate Gradient approach is better.



Speeding Up Conjugate Gradient — PCG(M)

The conjugate gradient algorithm is not the end of the story (it is just barely the end of the beginning). By combining the CG-algorithm with the idea of preconditioning ($M \approx A$, and M easily invertible) the Preconditioned CG algorithm can be derived.

Further, the CG-method can be extended to work for non-symmetric matrices as well:

Symmetry	Linear System $A\bar{x} = \bar{b}$	Eigenvalue Problem $A\bar{x} = \lambda\bar{x}$
$A = A^*$	CG	Lanczos
$A \neq A^*$	GMRES CGNE / CGNR BiCG, etc...	Arnoldi



Finite Differences vs. Finite Elements

This ends our overview of finite difference schemes for hyperbolic, parabolic, and elliptic problems. We have seen quite a few tools useful for both analysis and implementation of these schemes...

More Topics...

- Spectral Methods
- Mimetic Methods (a different view of the Finite Difference problem)
- Finite Element Methods — a different approach to approximation.
 - The FEM formulation is better suited for complex domains, and includes local error estimates which help us locally improve the solution exactly where these errors are large.
 - The biggest disadvantage, from a pedagogical point of view, is that whereas FD methods are quite straight-forward to implement, setting up a meaningful FEM-solver requires more “technology.” There are some nice (\$\$\$) commercial packages available (e.g. Comsol Multiphysics: <http://www.comsol.com/>).

