Optimization methods with Applications - Spring 2020

Home assignment 2

Due on 26/04/2020.

April 3, 2020

- 1. The efficiency of different iterative methods for solving a linear system.
 - (a) Implement programs for the four methods: Jacobi, Gauss-Seidel, Steepest (or Gradient) Descent (SD) and Conjugate Gradient (CG). See section 4.3.4 and Algorithm 6 for CG.
 - (b) Define the random symmetric positive definite matrix A below (which we define as sparse, see section 4.3.6). In Julia (after adding the SparseArrays package): using SparseArrays;

```
n = 256;
A = sprandn(n,n,5/n);
A = A'*spdiagm(0=>rand(n))*A + 0.1*spdiagm(0=>ones(n));
In python:
import numpy as np
from scipy.sparse import random
import scipy.sparse as sparse
n = 256
A = random(n, n, 5 / n, dtype=float)
v = np.random.rand(n)
v = sparse.spdiags(v, 0, v.shape[0], v.shape[0], 'csr')
A = A.transpose() * v * A + 0.1*sparse.eye(n)
```

For $\mathbf{x}^{(0)} = 0$, and a random \mathbf{b} , ($\mathbf{b} = \text{rand}(\mathbf{n})$ in Julia or $\mathbf{b} = \text{rand}(\mathbf{n}, 1)$ in Matlab) solve the system $A\mathbf{x} = \mathbf{b}$ with each of the methods, with at most 100 iterations. For Jacobi, use the weights $\omega = 1.0$ (standard Jacobi) and if it doesn't converge, search a reasonable $0 < \omega < 1$ that does lead for convergence most of

the times (remember, it is a random experiment).

Use the semilogy() plotting function to plot a convergence graph of $||A\mathbf{x}^{(k)} - \mathbf{b}||_2$. Also plot the convergence factor $\frac{||A\mathbf{x}^{(k)} - \mathbf{b}||_2}{||A\mathbf{x}^{(k-1)} - \mathbf{b}||_2}$.

2. Convergence properties

(a) Show that for any symmetric positive definite matrix $A \in \mathbb{R}^{n \times n}$, the Richardson method

$$\mathbf{x}^{(k+1)} = \mathbf{x}^{(k)} + \frac{1}{\|A\|} \left(\mathbf{b} - A\mathbf{x}^{(k)} \right),$$

converges to the solution of $A\mathbf{x} = \mathbf{b}$, where $\|\cdot\|$ is any induced matrix norm.

- (b) Show that if A in the previous section is indefinite (has both positive and negative eigenvalues), then the Richardson method diverges.
- (c) We will now prove the convergence of Steepest Descent with optimal choice of $\alpha^{(k)} = \alpha_{opt}$ as shown in Section 4.3.2 in the NLA notes.
 - i. Show that (recall: $A \succ 0$)

$$f(\mathbf{x}^{(k+1)}) = f(\mathbf{x}^{(k)}) - \frac{1}{2} \frac{\langle \mathbf{r}^{(k)}, A\mathbf{e}^{(k)} \rangle^2}{\langle \mathbf{r}^{(k)}, A\mathbf{r}^{(k)} \rangle} < f(\mathbf{x}^{(k)}),$$

where $f(\mathbf{x}) = \frac{1}{2} ||\mathbf{x}^* - \mathbf{x}||_A^2$, defined in Eq. 17 in the notes. The strict inequality holds as long as that $\mathbf{r}^{(k)} \neq 0$.

ii. Using the previous section, find a scalar factor $C^{(k)}$ such that

$$f(\mathbf{x}^{(k+1)}) = C^{(k)} \cdot f(\mathbf{x}^{(k)})$$

where $C^{(k)} < 1$ (this upper bound is achieved from the inequality above).

iii. Now, using the fact that for every symmetric matrix $A \in \mathbb{R}^{n \times n}$:

$$\forall \mathbf{v} \in \mathbb{R}^n : \lambda_{min} \leq \frac{\mathbf{v}^{\top} A \mathbf{v}}{\mathbf{v}^{\top} \mathbf{v}} \leq \lambda_{max},$$

(this is called a Rayleigh quotient) show that $C^{(k)} \leq 1 - \frac{\lambda_{min}}{\lambda_{max}} < 1$.

iv. Conclude that SD converges: $\lim_{k\to\infty} f(\mathbf{x}^k) = 0$, and hence $\lim_{k\to\infty} \mathbf{x}^k = \mathbf{x}^*$

3. In this question we will develop a method called GMRES(1). Assume that $A \in \mathbb{R}^{n \times n}$ is full rank, positive definite, but non-symmetric. We want to solve a linear system $A\mathbf{x} = \mathbf{b}$ using a method that is similar to Steepest Descent (SD). We define

$$\mathbf{x}^{(k+1)} = \mathbf{x}^{(k)} + \alpha^{(k)}(\mathbf{b} - A\mathbf{x}^{(k)}).$$

and we choose $\alpha^{(k)}$ to minimize the residual norm $\|\mathbf{r}^{(k+1)}\|_2 = \|\mathbf{b} - A\mathbf{x}^{(k+1)}\|_2$ with respect to $\alpha^{(k)}$ (in SD we minimized the A-norm of the error $\mathbf{e}^{(k+1)}$ instead of the residual norm).

(a) Show that

$$\alpha^{(k)} = \frac{(\mathbf{r}^{(k)})^{\top} A \mathbf{r}^{(k)}}{(\mathbf{r}^{(k)})^{\top} A^{\top} A \mathbf{r}^{(k)}}.$$

- (b) (non-mandatory) Show that as in SD, at each iteration we have to compute only one matrix vector multiplication $A\mathbf{r}^{(k)}$.
- (c) Demonstrate the convergence of the GMRES method for the following matrix:

$$A = \begin{bmatrix} 5 & 4 & 4 & -1 & 0 \\ 3 & 12 & 4 & -5 & -5 \\ -4 & 2 & 6 & 0 & 3 \\ 4 & 5 & -7 & 10 & 2 \\ 1 & 2 & 5 & 3 & 10 \end{bmatrix}.$$

Choose $\mathbf{b} = [1, 1, 1, 1, 1]^{\mathsf{T}}$, and $\mathbf{x}^{(0)} = [0, 0, 0, 0, 0]^{\mathsf{T}}$, and apply 50 iterations. Plot a graph of the residual norm vs. the iterations using the semilogy() plotting function.

- (d) The graph that you get in the previous subsection is monotone. Explain why?
- (e) Explicitly define the method GMRES(2):

$$\mathbf{x}^{(k+1)} = \mathbf{x}^{(k)} + \alpha_1^{(k)} \mathbf{r}^{(k)} + \alpha_2^{(k)} \mathbf{r}^{(k-1)}.$$

where the vector $[\alpha_1^{(k)}, \alpha_2^{(k)}]^{\top}$ is chosen to minimize $\|\mathbf{b} - A\mathbf{x}^{(k+1)}\|_2$. Guidance: Define a vector $\vec{\alpha}^{(k)} = [\alpha_1^{(k)}, \alpha_2^{(k)}]$, and a $n \times 2$ matrix $R^{(k)} = [\mathbf{r}^{(k)}, \mathbf{r}^{(k-1)}]$, and write in matrix form $\mathbf{x}^{(k+1)} = \mathbf{x}^{(k)} + R^{(k)}\vec{\alpha}^{(k)}$. Find a closed form for $\vec{\alpha}$.

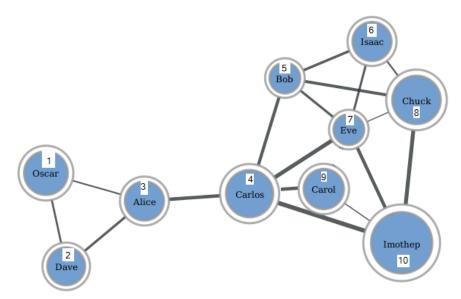
Remark for general knowledge: the method GMRES(1) that approximates only one parameter α stagnates for indefinite systems, choosing $\alpha^{(k)} = 0$ from some iteration on. The general GMRES(m) method approximates m such α 's (for m previous search directions) similarly to Conjugate Gradient and is able to solve indefinite systems as long as there aren't more than m negative eigenvalues. The GMRES(m) method also includes orthogonalizations in the minimization process for the α 's to avoid numerical errors.

4. Graph-Laplacians

Graph Laplacians are a powerful tool is computer science theory and have many applications. You may have a look at the following links if you're interested - there is a large amount of research devoted at these:

- https://en.wikipedia.org/wiki/Laplacian_matrix
- https://en.wikipedia.org/wiki/Algebraic_connectivity
- https://sites.google.com/a/yale.edu/laplacian/
- http://www.cs.yale.edu/homes/spielman/eigs/
- https://www.cs.yale.edu/homes/spielman/PAPERS/icm10post.pdf
- https://theory.epfl.ch/vishnoi/Lxb-Web.pdf

In many cases listed above, we are needed to solve linear systems or compute eigenvalues/eigenvectors with these graph Laplacians. Below we see a graph of a small social network:



And it's corresponding graph Laplacian matrix is:

Note that this matrix is symmetric and positive semi-definite. It has a null-space of a single vector: the constant [1, 1, 1, 1, ..., 1]. Let's define a vector $\mathbf{b} = [1, -1, 1, -1, ..., 1]$. That is a vector with a sum of zero, so there is a solution $L\mathbf{x} = \mathbf{b}$.

- (a) Solve the system using the Jacobi method, up to five digits of accuracy. Show the convergence graph of the residual norm. How many iterations are needed?
- (b) We will now accelerate Jacobi with a very powerful block preconditioner, instead of the simple diagonal one in Jacobi (the matrix D). We will partition the unknowns of L into two subgroups: $\{1,2,3\}$, and $\{4,5,6,7,8,9,10\}$. According to this partitioning we will define the submatrices $M_1 = L_{1..3,1..3} \in \mathbb{R}^{3\times 3}$ and $M_2 = L_{4..10,4..10} \in \mathbb{R}^{7\times 7}$. Now we will use the block preconditioner M which can be inverted by inverting the two smaller submatrices (which is cheaper):

$$M = \begin{bmatrix} M_1 & 0 \\ 0 & M_2 \end{bmatrix} \quad M^{-1} = \begin{bmatrix} M_1^{-1} & 0 \\ 0 & M_2^{-1} \end{bmatrix}$$

Repeat section 4a with the iteration $\mathbf{x}^{(k+1)} = \mathbf{x}^{(k)} + M^{-1}(\mathbf{b} - L\mathbf{x})$. Does the method converge in fewer iterations than standard Jacobi? Note that the matrices M_1 and M_2 are not singular like L, so this is well-defined.

(c) A student argued that M_2 is 7×7 and is still too large compared to the 10×10 in L, and that he wants to add another group. Using what we've learned in class, divide the unknowns $\{1, ..., 10\}$ into groups of balanced sizes 4,3,3 so that the iterations using a block diagonal M with the three blocks M_1, M_2, M_3 will be most efficient. Repeat section 4a with the new partitioning. Do you see a connection to the structure of the graph? Explain.