

MATH7003-00: Assignment #9

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Problem.

#1. Prove the inequality (8.5.13) on slide 5 of week11 [1].

$$\parallel x - x^{(m+1)} \parallel_{\infty} \le \frac{c}{1-c} \parallel x^{(m+1)} - x^{(m)} \parallel_{\infty}.$$

#2. Consider the example on slide 18 and write down m-files for Gauss-Seidel and SOR. Generate Table 8.9. Discuss about the result. [1].

Example. Solve

$$\frac{\partial^2 u}{\partial x^2} + \frac{\partial^2 u}{\partial y^2} = 0 \qquad 0 \le x, y \le 1$$

$$u(0,y) = \cos(\pi y), \quad u(1,y) = e^{\pi}\cos(\pi y), \quad u(x,0) = e^{\pi x}, \quad u(x,1) = -e^{\pi x}.$$

The true solution is

$$u(x,y) = e^{\pi x} \cos(\pi y).$$

Example. Recall the previous example (8.8.11). This was solved with both the Gauss-Seidel method and the SOR method. The initial guess for the iteration was taken to be the "bilinear" interpolation formula for the boundary data f:

$$u_h^{(0)}(x,y) = (1-x)f(0,y) + xf(1,y) + (1-y)f(x,0) + yf(x,1)$$
$$- [(1-y)(1-x)f(0,0) + (1-y)xf(1,0) + y(1-x)f(0,1) + xyf(1,1)]$$

at all interior grid points. The error test to stop the iteration was

$$\underset{1 \le j,k, \le N-1}{\operatorname{Max}} \left| u_h(x_j, y_k) - u_h^{(m)}(x_j, y_k) \right| \le \epsilon$$

with $\epsilon > 0$ given and the right-hand size of (8.7.5) used to predict the error in the iterate. The numerical results for the necessary number of iterates are given in **Table 8.9**.

Solution. (problem #1)

For the initial guess $x^{(0)}$, we know the fact that $x^{(1)} - x^{(0)} = Cr^{(0)}$ where $r^{(0)} = b - Ax^{(0)}$, C is the approximation of A^{-1} . Then, in general, we can define this form recursively, and we obtain a recursion formula for the error as

$$\begin{aligned} \mathsf{C}x - x^{(m+1)} &= x - x^{(m)} - C r^{(m)} = x - x^{(m)} - C \big[b - A x^{(m)} \big] \\ &= x - x^{(m)} - C \big[A x - A x^{(m)} \big] \end{aligned}$$

Thus, $x-x^{(m+1)}=(I-CA)(x-x^{(m)})$. By induction, we also obtain $x-x^{(m)}=(I-CA)^m(x-x^{(0)})$ for $m\geq 0$. In this, if we assume $\|I-CA\|<1$ for some matrix norm, then

$$||x-x^{(m)}|| = ||I-CA||^m ||x-x^{(0)}||$$

and, this converges th zero as $m \rightarrow \infty$ for any choice of initial guess $x^{(0)}$. This implies that

$$(I-CA)^m \to 0$$
 as $m \to \infty$.

Now, note that the theorem and the definition.

Theorem 7.9 Let A be a square matrix of order n. Then A^m converges to the zero matrix as $m \to \infty$ if and only if $r_{\sigma}(A) < 1$.

Definition Let A be an arbitrary matrix. The **spectrum** of A is the set of all eigenvalues of A, and it is denoted by $\sigma(A)$. The **spectral radius** is the maximum size of these eigenvalues, and it is denoted by

$$r_{\sigma}(A) = \max_{\lambda \in \sigma(A)} |\lambda|$$

Then, by Theorem 7.9, this is equivalent to $r_{\sigma}(I-CA)<1$. For the case of I-AC, we obtain

$$I - AC = A(I - CA)A^{-1}$$

and thus I-AC and I-CA are similar matrices and have the same eigenvalues, by the diagonalization theorem. (see Chapter 5 in [4]) So if $\|I-AC\|<1$, then $r_{\sigma}(I-CA)<1$ is true, even if $\|I-CA\|<1$ is not true, and convergence will still occur.

Therefore, $x - x^{(m+1)} = (I - CA)(x - x^{(m)})$ shows that the rate of convergence of $x^{(m)}$ to x is linear:

$$\|x-x^{(m+1)}\| \le c \|x-x^{(m)}\|$$
 for $m \ge 0$, with $c < 1$ unknown.

And, c is estimated computationally with

$$c = \text{Max} \frac{\| x^{(m+2)} - x^{(m+1)} \|}{\| x^{(m+1)} - x^{(m)} \|}.$$

Then, we now easily can produce an error bound that

$$\| x^{(m+1)} - x^{(m)} \| \ge \| [x - x^{(m)}] - [x - x^{(m+1)}] \| \cdots (I)$$

$$\ge \| x - x^{(m)} \| - \| x - x^{(m+1)} \| \cdots (II)$$

$$\ge \| x - x^{(m)} \| - c \| x - x^{(m)} \| \cdots (III)$$

$$\text{(I)} \quad \text{is} \quad x^{(m+1)} - x^{(m)} = x^{(m+1)} + (x-x) - x^{(m)} = \left[x^{(m+1)} - x \right] + \left[x - x^{(m)} \right] = \left[x - x^{(m)} \right] - \left[x - x^{(m+1)} \right],$$

(II) is by triangle inequality, and (III) is $\parallel x-x^{(m+1)}\parallel \ \leq c \parallel x-x^{(m)}\parallel$.

Thus, we obtain

$$\parallel x^{(m+1)} - x^{(m)} \parallel \ge \parallel x - x^{(m)} \parallel - c \parallel x - x^{(m)} \parallel$$

$$= (1 - c) \parallel x - x^{(m)} \parallel$$

and, finally we have

$$\parallel x-x^{(m)}\parallel \ \leq \ \frac{1}{1-c}\parallel x^{(m+1)}-x^{(m)}\parallel \ \ \Rightarrow \ \ \parallel x-x^{(m+1)}\parallel \ \leq \ \frac{c}{1-c}\parallel x^{(m+1)}-x^{(m)}\parallel$$

$$\text{since } \parallel x-x^{(m+1)}\parallel \leq c\parallel x-x^{(m)}\parallel \ \Rightarrow \ c\parallel x-x^{(m)}\parallel \ \leq c\left(\frac{1}{1-c}\parallel x^{(m+1)}-x^{(m)}\parallel \ \right).$$

Solution. (problem #2)

For j = 1, 2, ..., N-1 and k = 1, 2, ..., N-1, the 'matrix-free' form of the Gauss-Seidel (G-S) method of the given partial differential equation is as follows.

$$u_h^{(m+1)}(x_j,y_k) = \frac{1}{4} \left[u_h^{(m)}(x_{j+1},y_k) + u_h^{(m)}(x_j,y_{k+1}) + u_h^{(m+1)}(x_{j-1},y_k) + u_h^{(m+1)}(x_j,y_{k-1}) \right].$$

For the boundary points, use

$$u_h^{(m)}(x_i, y_k) = f(x_i, y_k) \text{ for all } m \ge 0.$$

Since the equation of this example is the homogeneous equation, $g(x_j, y_k) = 0$. (refer to (8.8.12) in [1]) Also, the successive overrelaxation (SOR) method is as follows.

$$\boldsymbol{v}_{h}^{(m+1)}(x_{j},y_{k}) = \frac{1}{4} \left[u_{h}^{(m)}(x_{j+1},y_{k}) + u_{h}^{(m)}(x_{j},y_{k+1}) + u_{h}^{(m+1)}(x_{j-1},y_{k}) + u_{h}^{(m+1)}(x_{j},y_{k-1}) \right],$$

$$u_h^{(m+1)}(x_j,y_k) = \omega \, \boldsymbol{v} \, _h^{\,(m+1)}(x_j,y_k) + (1-\omega) u_h^{(m)}(x_j,y_k) \ \, \text{for} \ \, j=1,2,\dots,N-1, \ \, k=1,2,\dots,N-1.$$

Note that the optimal acceleration parameter is

$$\omega^* = \frac{2}{1 + \sqrt{1 - \xi^2}} \quad \text{where } \xi = 1 - 2\sin^2\left(\frac{\pi}{2N}\right).$$

If $\omega^*=1$, then the formula of SOR is directly equals to the formular of Gauss-Seidel i.e., we can easily know that the SOR method is the generalized version of Gauss-Seidel method.

We use the stop criterion in problem #1 as

$$\| x - x^{(m+1)} \|_{\infty} \le \frac{c}{1-c} \| x^{(m+1)} - x^{(m)} \|_{\infty}, \quad \text{where} \quad c \doteq \frac{\| x^{(m+1)} - x^{(m)} \|_{\infty}}{\| x^{(m)} - x^{(m-1)} \|_{\infty}},$$

in order for our implementation. Now, the results of problem #2 is attached below. (see Figure 1.)

Table 8.9 Number of iterates necessary to solve (8.8.5)

N	ě	Gauss-Seidel	SOR
8	.01	25	12
8	.001	40	16
16	.001	142	32
32	.001	495	65
8	.0001	54	18
16	.0001	201	35
32	.0001	733	71

Table 8.9 Number of iterates necessary to solve (8.8.5) Gauss-Seidel 25 12 0.01 0.001 16 0.001 142 31 32 495 0.001 54 16 0.0001 201 31 32 0.0001 733 65 fx >>

Figure 1. The results of the problem 2 using MATLAB. The left-side is in the textbook.

Discussion. (problem #2)

Our problem has the equivalent step size h = 1/N for x and y. Hence the finite mesh grid depends on the step parameter N. This implies that if N is larger, then the iteration number is basically larger.

Moreover, if the stop criterion $\epsilon > 0$ is smaller, we need to iterate more than previous, in order to meet more rigorous criterion ϵ .

In Table 8.9., we suggest the Gauss-Seidel and SOR results for N=8,16,32 and $\epsilon=0.01,0.001,0.0001$, and we can see that if N and ϵ is larger, the iteration of both methods is larger.

Especially, we already learned the fact that the SOR method is the acceleration of Gauss-Seidel method, so we intuitively can see that the iteration number of SOR results are better than Gauss-Seidel results in Table 8.9. So we conclude that the use of SOR greatly reduces the resulting work, although it still is large when N is large.

With either method, note that doubling N will increase the number of equations to be solved by a factor of 4, and thus the work per iteration will increase by the same amount.

Table_8_9_HW9.m

```
%% MATH7003-00: Assignment #9, 2019310290 Sangman Jung.
N = [8 16 32]; % step N
epsilon = [0.01 0.001 0.0001]; % error criterion
f = 0(x,y) \exp(pi*x).*\cos(pi*y); % exact solution of our problem
ksi = 1-2.*((sin(pi./(2*N))).^2);
omega = 2./(1+sgrt(1-(ksi).^2)); % optimal acceleration parameter
uval\{length(N), length(epsilon), 2\} = []; % pre-allocation of the numerical solution
\texttt{iteration} = \texttt{zeros}(\texttt{length}\,(\texttt{N})\,, \texttt{length}\,(\texttt{epsilon})\,)\,;\,\,\,\$\,\,\,\texttt{pre-allocation}\,\,\,\texttt{of}\,\,\,\texttt{G-S}\,\,\,\texttt{iteration}
iteration sor = zeros(length(N),length(epsilon)); % pre-allocation of SOR iteration
for w iter = 1:2 % w iter = 1 is G-S, w iter = 2 is SOR
    for N_iter = 1:length(N) % iterate per step N
       h = 1/N(N_iter); % step size
       x = 0:h:1; y = x; % spatial variables (grid)
       nx = length(x); ny = length(y); % the size of the grid
       u ini = zeros(nx,ny); % initial guess
       u_exact = zeros(nx,ny); % pre-allocation of exact solution
        u_new = zeros(nx,ny); % pre-allocation of numerical solution
       for eps iter = 1:length(epsilon) % iteration of the criterion
           m_iter = 1; % initialize the iteration number
           for k = 1:ny
              for j = 1:nx
                  u_{ini}(j,k) = (1-x(j))*f(0,y(k))+x(j)*f(1,y(k))+(1-y(k))*f(x(j),0)+...
                     y(k) *f(x(j),1) - ((1-y(k)) *(1-x(j)) *f(0,0) + (1-y(k)) *x(j) *f(1,0) + ...
                      y(k)*(1-x(j))*f(0,1)+x(j)*y(k)*f(1,1)); % compute the initial guess
                  u_{exact(j,k)=f(x(j),y(k))}; % allocate the exact solution
              end
           u old = u ini; % update the m-iteration
           % boundary values of the problem
           u_new(1,:) = u_exact(1,:); % left vertical
           u_new(:,1) = u_exact(:,1); % bottom
           u_new(end,:) = u_exact(end,:); % right vertical
           u new(:,end) = u exact(:,end); % top
           error = 1: % initialize the error
           while 1 % Main Loop
for k = 2:ny-1
                 for j = 2:nx-1
                      \verb"u_new(j,k) = (\verb"u_old(j+1,k) + \verb"u_old(j,k+1) + \verb"u_new(j-1,k) + \verb"u_new(j,k-1)) / 4;
                      if w iter > 1
                          % Successive OverRelaxation (SOR) method
                         \label{eq:u_new} \verb"u"_new"(j,k) = \verb"omega"(N_iter)" * \verb"u"_new"(j,k) + (1-omega"(N_iter)") * \verb"u"_old"(j,k);
              if m_iter > 1 % compute the error (8.7.5) in Chapter 8.7
                 c = max(max(abs(u_new-u_old)))/max(max(abs(u_old-u_2old)));
error = c/(1-c)*(max(max(abs(u_new-u_old))));
              if error <= epsilon(eps_iter) % error criterion
              m iter = m iter + 1; % update the error after passing the criterion
              u_2old = u_old; % uptate from m-1 to m
              u_old = u_new; % update from m to m+1
           if w_iter == 1 % G-S iteration
              iteration(N_iter,eps_iter) = m_iter;
           else % SOR iteration
              iteration_sor(N_iter,eps_iter) = m_iter;
           uval(eps_iter,N_iter,w_iter) = (u_new); % save the numerical solution
% summarize our results in order to obtain the table
N_table = [N(1) N N]';
eps_table = [epsilon(1) repmat(epsilon(2),1,3) repmat(epsilon(3),1,3)]';
iteration = reshape(iteration,[9 1]);
iteration = iteration([1 4:9]);
iteration_sor = reshape(iteration_sor,[9 1]);
iteration_sor = iteration_sor([1 4:9]);
fprintf('\nTable 8.9 \nmber of iterates necessary to solve (8.8.5) \n')
fprintf(' N
fprintf('----
for t = 1:length(N_table)
                                         %3.0f
                                                           %3.0f\n',...
                           %6.4a
      [N_table(t) eps_table(t) iteration(t) iteration_sor(t)]);
fprintf('-
```

References.

- [1] Atkinson, K. E. (2008). An introduction to numerical analysis. John wiley & sons.
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- [3] Atkinson, K. E., & Han, W. (1985). Elementary numerical analysis (p. 17). New York et al.: Wiley.
- [4] Lay, D. C. (2012). Linear algebra and its applications. Addison-Wesley.