Assignment 6 Solutions

Aaron Cahn University of Wisconsin-Madison cahn@cs.wisc.edu

May 8, 2015

1 Solutions

1.1 Question 1

Listing 1: Matlab Commands

```
% Qustion 1
% generate the A matrix with intervals of eigenvalues
A=generate interval matrix (100, [-3 \ 4 \ 1], 1);
b = rand(100,1);
fig=figure;
for n=1:10-1
    % run my_gmres to obtain Q and S
    [Q,S]=my\_gmres(A,b,n);
    % compare spectrums
    specA = eig(A);
    specS = eig(S);
    % plot the eigenvalues of A and Sn
    subplot(3,3,n)
    plot (real (specA), zeros (size (specA, 1)), 'bo', ...
          real(specS), zeros(size(specS,1)), 'rx')
    title (sprintf ('n=%d',n))
end
```

The code above was run on various incarnations of the matrix A. For each expirement the matrix A is 100×100 in dimension. Each expirement varied the number of intervals the eigenvalues of A were centered around and how large this interval was. The number of intervals ranged from 1 to 5.

The conclusion of the expirements was very clear. The speed of convergence of $\sigma(S_n)$ towards $\sigma(A)$ was proportional to the number of intervals the eigenvalues were centered around. For example, consider the plots below. This expirement was run with three intervals centered at -3, 4, and 1. From Figure 1 we see that after three iterations the $\sigma(S_n)$ is starts to align well with the $\sigma(A)$ (i.e., the three eigenvalues of S_n align with the three intervals). This is true in general. Specifically the convergence happens at a rate of $\frac{1}{|I|}$ where |I| is the number of intervals the eigenvalues of A are centered around.

Listing 2: Matlab Commands

```
% runs the gmres algorithm for n iterations
function [Q, S, xtil] = my\_gmres(A, b, n)
    % track Q, S, and X's
    Q = []; S = []; X={};
    % pre-process steps
    b1 = norm(b);
    Q(:,1) = b/norm(b);
    % run n iterations of gmres
    for i = 1:n
       v=A*Q(:, i);
       S(:, i)=Q'*v;
       v=v-Q(:,1:i)*S(:,i);
       Q(:, i+1)=v/norm(v);
       S(i+1,i)=norm(v);
       b1 = [b1; 0];
       X{i}=Q(:,1:i)*(S\b1);
    end
    % return Sn and final x
    x til = X\{n\};
    S=S(1:end-1,:);
end
```

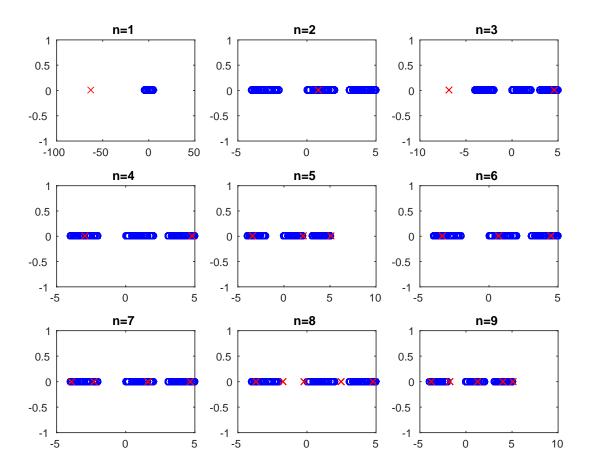


Figure 1: Plots of $\sigma(S_n)$ (red) and $\sigma(A)$ (blue). Plots correspond to stopping GMRES after n iterations.

1.2 Question 2

Listing 3: Matlab Commands % Part (ii) n: 5 error: 0.104672 n: 10 error: 0.000154 n: 20 error: 0.000000 n: 90 error: 0.000000

Listing 4: Matlab Commands

```
% Question 2
% Generate A from generate_interval_matrix
m = 100; n = [4 \ 5 \ 10]; e = 1;
A = generate_interval_matrix (m, n, e);
% Generate a random b
b = rand(100,1);
% Find the actual x (partt(i))
x = A \setminus b
% Use function from Q1 (part(ii))
n = [5 \ 10 \ 20 \ 90];
errs = []; Q = {}; S = {};
for i = 1:4
    % gather data
    [Q\{i\},S\{i\},xtil]=my\_gmres(A,b,n(i));
    errs = [errs norm(A*xtil-b)];
    disp(sprintf('n: \%d error: \%f', n(i), norm(A*xtil-b)));
end
plot(n, errs, 'b—o')
```

As with the expirement from Question 1 the number of intervals was kept constant at 3 and the interval ranges were varied from 0.5 to 2. The results obtained reflect the sentiments expressed above. The convergence (relative error decreased) of $\sigma(S_n)$ to $\sigma(A)$ occurred at a rate of $\frac{1}{|I|}$. Therefore after ten iterations the error was extremely small (0.000154) and became almost negligible after that. After ninety iterations the $\sigma(S_n)$ was essentially the same as the $\sigma(A)$ and after one hundred there was complete convergence. However, it makes little sense to iterate this long as the cost trade off of using an indirect method is lost (*i.e.*, the complexity is $O(m^3)$). Figure 2 demonstrates the reduction in error at the different values of n.

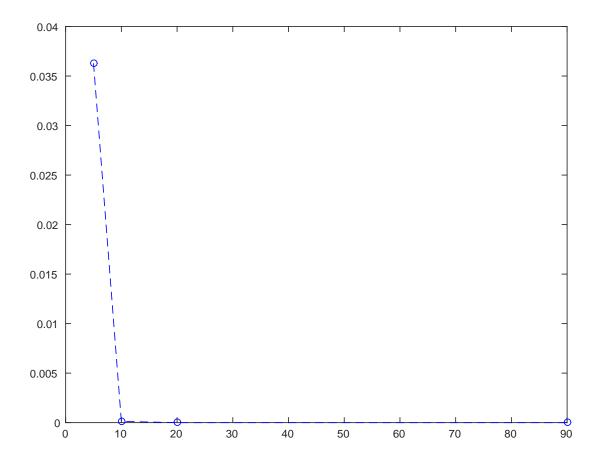


Figure 2: Plot showing the relative error $||A\tilde{x} - b||_2$ for different values of n.

1.3 Question 3

```
Listing 5: Matlab Commands

% Generate B

B=rand(30,10);

D=rand(10,10);

for k=100:100:999
% Generate C
C=D;
C(1:2,:)=C(1:2,:)*k;
% Generate A
A=B*C;
[U,S,V] = svd(A);
end
```

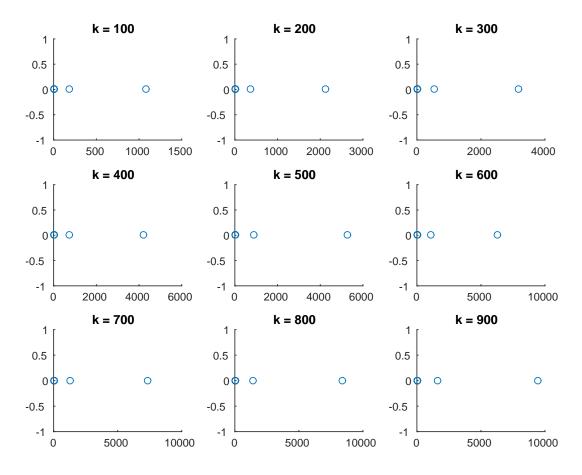


Figure 3: Plots of singular values of A at various k. Note, the first two singular values are much larger than the rest (k*10 and k times larger respectively).

The experiment setup was similar to the two expirements above, except the value of k was the variable in this case. Figure 3 shows one of these expirements. It is clear from Figure 3 that the first two singular values of A are much larger than the other singular values of A. Specifically, the first two singular values are roughly k*10 times and k times larger than the other singular values. This makes intuitive sense as A is built out of an underlying matrix where the first two rows are k times larger than the other rows. Therefore, the first two singular values and left singular vectors of A capture the structure of A quite well.

Listing 6: Matlab Commands

```
function [A] = generate_interval_matrix(m,n,e)
    I = [];
    for i = 1 : size(n, 2)
        % generate random values in the interval
        I = [I (((n(i)+e)-(n(i)-e)).* ...
            rand (round (m/size (n,2)), 1)+(n(i)-e))'];
    end
    % m/n has remainder (add those values)
    if size(I,2) < m
        I = [I (((n(i)+e)-(n(i)-e)).* ...
            rand (m-size(I,2),1)+(n(i)-e));
    end
    % generate A with eigenvalues from I
    D = diag(sort(I, 'descend'));
    P = rand(m);
    A = P*D*inv(P);
end
```