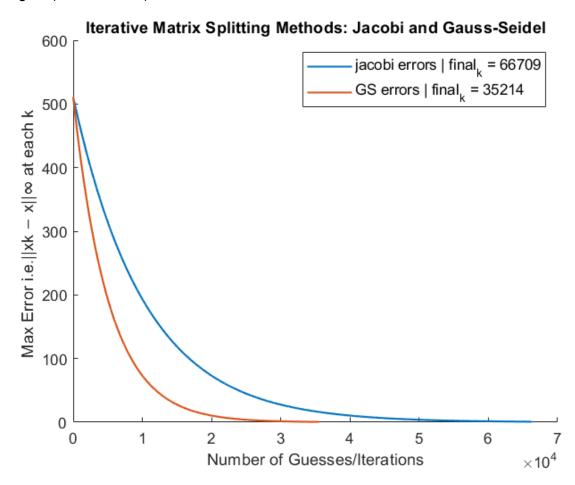
```
Andrew Garwood
Written HW 4
Problem 1
a)
% find ||Ax - b||
max_error = max(abs(A * x - b)); \% = 6.7540e-13
% considering that the max error is extremely small, I believe it is
% reasonable to state that this error is zero
b, c, d)
% (b) Solve the system of equations Ax = b using the Jacobi method with a
% tolerance of 10-4 and an initial guess of all ones. At each guess xk,
% find the error between the guess and the true solution.
% In other words, find ||xk - x|| = x at each k.
% Solve for phi using Jacobi method.
% Jacobi Partition
P = diag(diag(A));
T = A - P;
% Get M Matrix
M = -P \T;
lambda = eig(M);
max_lambda = max(abs(lambda))
% construct initial guess
phi0 = zeros(20, 1);
phi0(:, 1) = 1;
% max lambda is less than 1, so no infinite while loop
% designate template for previous guess and new guess
new phi = zeros(20, 1);
% Set previous guess to phi0
prev_phi = phi0;
tolerance = 1e-4; % initialize tolerance level
error = tolerance + 1; % ensure loop runs at least once
```

k = 1; % indexing variable

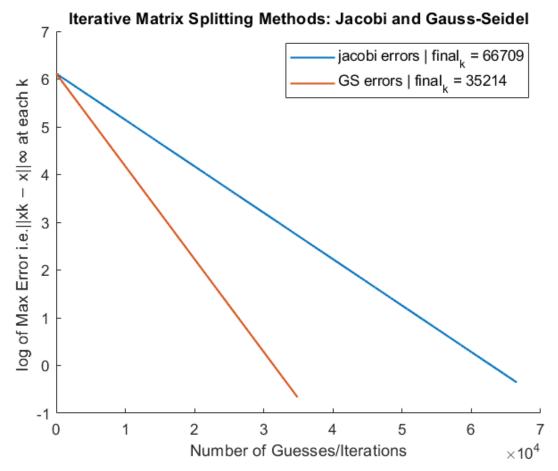
```
% At each guess Xk find error between guess and solution
true solution = x;
% arbitrarily set size of guess error vector
jacobi errors = zeros(1, 101);
% add first guess error measurement to vector
jacobi_errors(k) = max(abs(phi0 - true_solution));
if max lambda < 1
  while error >= tolerance
    new_phi = P(-T * prev_phi + b);
     error = max(abs(new_phi - prev_phi));
    k = k + 1; % increment/set current guess number
    jacobi_errors(k) = max(abs(new_phi - true_solution));
     prev_phi = new_phi; % set previous guess to new one to progress!!
  end
end
% (c) Repeat part (b) using the Gauss-Seidel method.
% Solve for phi using GS method.
% GS Partition
P = tril(A);
T = A - P;
% Get M Matrix
M = -P \T;
lambda = eig(M);
max_lambda = max(abs(lambda));
% construct initial guess
phi0 = zeros(20, 1);
phi0(:, 1) = 1;
% max_lambda is less than 1, so no infinite while loop
% designate template for previous guess and new guess
```

```
new phi = zeros(20, 1);
% Set previous guess to phi0
prev phi = phi0;
tolerance = 1e-4; % initialize tolerance level
error = tolerance + 1; % ensure loop runs at least once
k = 1; % indexing variable
% At each guess Xk find error between guess and solution
true solution = x;
% arbitrarily set size of guess error vector
GS errors = zeros(1, 101);
% add first guess error measurement to vector
GS_errors(k) = max(abs(phi0 - true_solution));
if max_lambda < 1
  while error >= tolerance
     new phi = P(-T * prev phi + b);
     error = max(abs(new_phi - prev_phi));
    k = k + 1; % increment/set current guess number
     GS errors(k) = max(abs(new phi - true solution));
     prev phi = new phi; % set previous guess to new one to progress!!
  end
end
% Plot that ish
% x jacobi = 0:length(jacobi errors) - 1;
% x GS = 0:length(GS errors) - 1;
% figure()
% hold on
% semilogy(x jacobi, log(jacobi errors), 'DisplayName', 'jacobi errors | final k = 66709',
'LineWidth', 1.25)
% semilogy(x_GS, log(GS_errors), 'DisplayName', 'GS errors | final_k = 35214', 'LineWidth',
1.25)
%
% hold off
% xlabel('Number of Guesses/Iterations')
```

% ylabel('log of Max Error i.e.||xk - x|| ∞ at each k ') % title('Iterative Matrix Splitting Methods: Jacobi and Gauss-Seidel') % % legend('FontSize', 10, 'Position', [0.5 0.5 0.1 0.2]) % legend('FontSize', 10)



e, f)



% (f) What do the slopes of the lines from part (e) represent? Can you relate them % to a quantity from lecture?

% slope = (log(method_error_k1) - log(method_error_k2)) / (k1 - k2)

% Jacobi log error slope = -0.00009721799004

% Jacobi max abs lambda = 0.999902786604301

% Jacobi max abs lambda + abs(jacobi_log_error) = 1.000000004594341

% GS log error slope = -0.0001944366197

% GS max abs Lambda = 0.999805582659051

% GS max abs lambda + abs(gs log error) = 1.000000019278751

% I grow tired, my apologies.

```
%SOR method
% The coding homework asked you to use \omega = 1.5 to solve the system A114\phi = \rho.
% You should have found that SOR (with \omega = 1.5) was slightly faster than
% GaussSeidel. In this problem, we will try to find an optimal value of \omega
% and see how that impacts the speed of our splitting method.
% (a) For every value of \omega between \omega = 1 and \omega = 1.999 in increments of 0.001,
% calculate M = -P^-1 * T using the corresponding P and T from equation (1)
% and then find the absolute value of the largest eigenvalue \lambda 1 of this matrix M.
% Make a plot of the absolute value of \lambda 1 versus \omega. (That is, \omega should be on the
% x-axis and the absolute value of the largest eigenvalue of M should be on the
% y-axis.)
% A = diagonal + upperTri + lowerTri
% Then P = (1 / omega) * D + L, T = ((omega - 1) / omega) * D + U
% omega between 1 and 2. 1 => GS
D = diag(diag(A));
L = tril(A) - D;
U = triu(A) - D;
max lambdas = zeros(1, length(1:.001:1.999));
i = 1:
for omega = 1:.001:1.999
  P = (1 / omega) * D + L;
  T = ((omega - 1) / omega) * D + U;
  M = -P \T:
  lambda = eig(M);
  max lambdas(i) = max(abs(lambda));
  i = i + 1;
end
x = 1:.001:1.999:
figure()
hold on
plot(x, max_lambdas, 'DisplayName', 'Max_eigenvals', 'LineWidth', 1.25)
hold off
xlabel('omega 1 to 1.999, intervals of 0.001')
```

```
ylabel('abs max(eigenvalue) of M') title('SOR method, finding optimal omega') legend('FontSize', 10, 'Position', [0.5 0.5 0.1 0.2]) % legend('FontSize', 10)
```

%с

% optimal_omega approx 1.973 % min(max lambdas) = 0.973

% d

% SOR with optimal omega took 461 guesses/iterations

% maximum error was 8.142494574814307e-06!!!

- % This is a lot faster than the Jacobi And Gauss-Seidel Methods % Jacobi took 10348, Gauss_seidel took 6104
- % (e) In lecture, we mentioned that calculating eigenvalues was a fairly slow process.
- % Do you think that the speed increase over Jacobi or Gauss-Seidel was worth
- % the time spent finding an optimal ω ? Would your answer change if we had
- % to solve many systems of equations with the same matrix A? Explain your
- % reasoning
- % For a small matrix, it would be worth it, however, because the
- % eig(Matrix) function is O(n^3) i do not think it would be worthwhile to
- % find an optimal omega in applications with huge amounts of data. . i.e.
- % large matrices
- % For the same matrix A? If I am reading the question right, then you mean
- % the A that we read in at the beginning of this question. Then yes I think
- % it would be worth it because A is relatively small so we would be
- % spending marginally more time to find a more precise answer; however, as
- % I stated above, if A is arbitrarily large, then I do not think it would
- % be worthwhile.