

ENGR 8102: COMPUTATIONAL ENGINEERING

Solution Set 1 (due in class on Monday, 10/26)

Questions:

1. (7 pts.) Modify the code `laplace.m` given in class, to feature a 10x10 grid (including boundaries), instead of 5x5. Replace the zero boundary with the last two digits of your uga-ID (810-??-??XX), keep the other boundaries the same. Plot the temperature distribution and isothermal lines. Include a copy of your code, the temperature distribution matrix, and the graphs along with your solution. Submit a soft copy of your code to `caner@uga.edu`.

SOLUTION. Assuming $XX = 25$:

```
function laplace

% create operator matrix m
m = diag( -4*ones(64,1) );
band1 = ones(1,63);
band1(8:8:60)=0;
m = m + diag( band1, -1 ) + diag( band1, 1 );
m = m + diag( ones(56,1), -8 ) + diag( ones(56,1), 8 );

% construct vector b using boundary conditions

bndry = [-175 -100 -100 -100 -100 -100 -100 -150 -75  0  0  0  0 ...
 0  0 -50 -75  0  0  0  0  0  0 -50 -75  0 ...
 0  0  0  0  0 -50 -75  0  0  0  0  0  0 ...
 -50 -75  0  0  0  0  0  0 -50 -75  0  0  0 ...
 0  0  0 -50 -100 -25 -25 -25 -25 -25 -25 -75]';
res = inv(m) * bndry;

% convert solution from vector to matrix format
mres = [ res(1:8)'; res(9:16)'; res(17:24)'; res(25:32)'; ... \
 res(33:40)'; res(41:48)'; res(49:56)'; res(57:64)' ];

% attach boundary

t_dist(:,1) = 75*ones(1,10);
t_dist(:,10) = 50*ones(1,10);
t_dist(1,:) = 100*ones(1,10);
t_dist(10,:) = 25*ones(1,10);
t_dist(2:9,2:9) = mres

figure; % plot temperature distribution
surf( t_dist );
view(-220,20)
title('Temperature distribution');

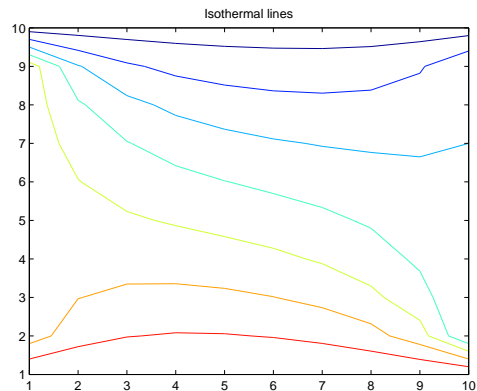
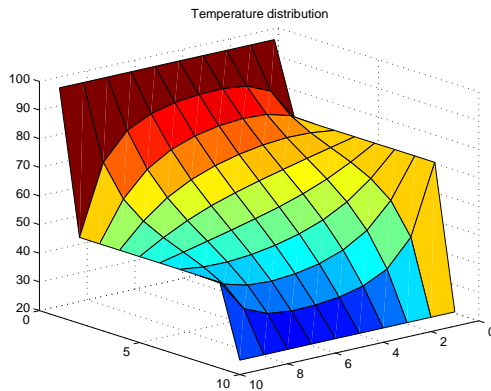
figure; % plot isothermal lines
contour( t_dist )
title('Isothermal lines')
```

Here is the output of the code:

```

>> hw21_laplace
t_dist =
100.0000 100.0000 100.0000 100.0000 100.0000 100.0000 100.0000 100.0000 100.0000 100.0000
 75.0000  86.1291  89.7163  90.6861  90.5187  89.5683  87.5922  83.5455  74.3145  50.0000
 75.0000  79.8000  82.0502  82.5092  81.8204  80.1623  77.2549  72.2751  63.7127  50.0000
 75.0000  76.0205  76.1754  75.4801  74.0913  72.0058  68.9900  64.5874  58.2611  50.0000
 75.0000  73.1067  71.1508  69.1444  67.0589  64.7795  62.1121  58.8233  54.7444  50.0000
 75.0000  70.2556  66.1767  62.8879  60.2205  57.9411  55.8556  53.8492  51.8933  50.0000
 75.0000  66.7389  60.4126  56.0100  52.9942  50.9087  49.5199  48.8246  48.9795  50.0000
 75.0000  61.2873  52.7249  47.7451  44.8377  43.1796  42.4908  42.9498  45.2000  50.0000
 75.0000  50.6855  41.4545  37.4078  35.4317  34.4813  34.3139  35.2837  38.8709  50.0000
 25.0000  25.0000  25.0000  25.0000  25.0000  25.0000  25.0000  25.0000  25.0000  25.0000
>>

```



2. (7 pts.) Modify the code `gauss_seidel.m` to feature a 10x10 grid (including boundaries) instead of 5x5. Replace the zero boundary with the last two digits of your uga-ID (810-??-??XX), keep the other boundaries the same. Run enough iterations until the absolute value of the difference between componentwise temperature distribution difference between (1) and (2) is bounded above by 1×10^{-2} . Plot the temperature distribution and isothermal lines. Include a copy of your code, the temperature distribution matrix, difference of temperature distributions (between Laplace and Gauss-Seidel), and the graphs along with your solution. Submit a soft copy of your code to `caner@uga.edu`.

SOLUTION. Assuming $XX = 25$:

```

function gs

% create grid with boundary
T = zeros(10,10);
T(:,1) = 75*ones(1,10);
T(:,10) = 50*ones(1,10);
T(10,:) = 25*ones(1,10);
T(1,:) = 100*ones(1,10);
T_new = zeros(10,10); % create new iteration matrix

% main iteration
error = 1;
while ( error > 0.0005 )
    for i=2:9
        for j=2:9
            T_new(i,j) = ( T(i+1,j) + T(i-1,j) + T(i,j+1) + T(i,j-1) )/4;
        end
    end
end

```

```

end
error = max(max( abs(T(2:9,2:9) - T_new(2:9,2:9)) ));
T(2:9,2:9) = T_new(2:9,2:9);    % update T
end
T

figure;          % plot temperature distribution
surf( T );
view(-220,20)
title('Temperature distribution');

figure;          % plot isothermal lines
contour(T)
title('Isothermal lines')

```

Here is the output of the code. Note that we had to set the approximate error a lot less to achieve the exact error of 0.01.

```

>> hw21_gs
T =
100.0000 100.0000 100.0000 100.0000 100.0000 100.0000 100.0000 100.0000 100.0000 100.0000
 75.0000  86.1282  89.7146  90.6837  90.5161  89.5657  87.5899  83.5437  74.3136  50.0000
 75.0000  79.7982  82.0470  82.5049  81.8155  80.1574  77.2506  72.2719  63.7110  50.0000
 75.0000  76.0182  76.1711  75.4743  74.0847  71.9992  68.9842  64.5831  58.2588  50.0000
 75.0000  73.1041  71.1459  69.1378  67.0514  64.7719  62.1055  58.8183  54.7418  50.0000
 75.0000  70.2530  66.1718  62.8813  60.2130  57.9335  55.8489  53.8442  51.8906  50.0000
 75.0000  66.7366  60.4083  56.0041  52.9876  50.9021  49.5141  48.8203  48.9772  50.0000
 75.0000  61.2856  52.7217  47.7408  44.8327  43.1747  42.4865  42.9466  45.1983  50.0000
 75.0000  50.6846  41.4528  37.4055  35.4291  34.4787  34.3116  35.2820  38.8700  50.0000
 25.0000  25.0000  25.0000  25.0000  25.0000  25.0000  25.0000  25.0000  25.0000  25.0000

>> abs( T - t_dist )
ans =
    0         0         0         0         0         0         0         0         0         0
    0    0.0009    0.0017    0.0024    0.0026    0.0026    0.0023    0.0018    0.0009    0
    0    0.0018    0.0032    0.0043    0.0049    0.0049    0.0043    0.0032    0.0017    0
    0    0.0023    0.0043    0.0058    0.0066    0.0066    0.0058    0.0043    0.0023    0
    0    0.0026    0.0049    0.0066    0.0075    0.0076    0.0066    0.0050    0.0026    0
    0    0.0026    0.0049    0.0066    0.0075    0.0076    0.0067    0.0050    0.0027    0
    0    0.0023    0.0043    0.0059    0.0066    0.0066    0.0058    0.0043    0.0023    0
    0    0.0017    0.0032    0.0043    0.0050    0.0049    0.0043    0.0032    0.0017    0
    0    0.0009    0.0017    0.0023    0.0026    0.0026    0.0023    0.0017    0.0009    0
    0         0         0         0         0         0         0         0         0         0
>>

```

