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% % ELEC 4700 Assignment 1: Monte-Carlo Modelling of Electrons
% % Mike Crawford
% % 100952432
% % Question1
%%Would like the one on one evaluation please
% This code is designed to simulate a specified number of electrons movement
% in silicon using the Monte-Carlo method. The effective mass is known and
% each electron is given a random direction. Boundaries were put on the box
% to allow electrons to pass through the other side of the y plane and
% reflect off the x plane. The thermal velocity is calculated by the equation
%
%  $V_{th} = \sqrt{(C.k*T)/(m_n)}$ 
%
% The parameters are labeled below. The mean free path is can be found with
% the thermal velocity and the mean time between collisions, which is given
% as 0.2 ps. The simulation plots these electrons during the specified
% timeframe as well as the average semiconductor temperature over time.

%reset
clearvars
clearvars -GLOBAL
close all
format shortE
global C
global Ecount
global Vx Vy Vtotal x y
Ecount = 5000; %setting amount of electrons used
C.mo = 9.10938215e-31; %Electron mass
C.k = 1.3806504e-23; %Boltzmann constant

T = 300; %Temperature (K)
mn = 0.26*C.mo;
L = 200e-9; %Length of Border
W = 100e-9; %Width of Border
Vth = sqrt(2*(C.k*T)/mn); %Calculation of Thermal Velocity
dt = 10e-15; %timestep
Stop = 500*dt; %timeframe
x = zeros(Ecount, 2); %array of x positions
y = zeros(Ecount, 2); %array of y positions

Temperature = zeros(1,2); %array of temperatures
Time = 0;
VisibleEcount = 10; %number of visible electrons

for i = 1:Ecount %setting random locations for each electron
    x(i,1) = rand()*200e-9;
    y(i,1) = rand()*100e-9;
end

Vx(1:Ecount) = Vth * cos(2*pi*randn(1, Ecount)); %setting velocities

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Vy(1:Ecount) = Vth * sin(2*pi*randn(1, Ecount));
for i = 1:Ecount
    Vtotal(i) = sqrt(Vx(i)^2 + Vy(i)^2); %summing x and y velocities
end

figure(1) %plotting electrons
subplot(2,1,1);
axis([0 L 0 W]);
title('Monte-Carlo model of electron movement in Silicon');
xlabel('X');
ylabel('Y');
hold on;

subplot(2,1,2); %plotting avg temperature
axis([0 Stop 0 500]);
title('Semiconductor Temperature');
xlabel('Time (s)');
ylabel('Temperature (K)');
hold on;

for i = 1:Ecount %Sum of all initial temperatures
    Temperature(1,2) = Temperature(1,2) + (mn*Vtotal(i)^2)/(2*C.k);
end
%intitilal average temperature
AvgTemperature = Temperature(1,2)/Ecount;
TemperaturePlot = [300 AvgTemperature]; %setting temperature line to plot
TimePlot = [0 Time]; %setting time line to plot
plot(Time, AvgTemperature);

%reseting values before loop
VTotal = 0;
Temperature(1,2) = 0;
AvgTemperature = 0;

%looping through timesteps
while Time < Stop
    subplot(2,1,1)
    for j = 1:Ecount %looping through electrons
        x(j,2) = x(j,1); %updating previous positions
        y(j,2) = y(j,1);
        x(j,1) = x(j,1) + (dt * Vx(j));%updating new positions
        y(j,1) = y(j,1) + (dt * Vy(j));
        if x(j,1) > L %setting right side boundary condition
            x(j,2) = 0;
            x(j,1) = dt * Vx(j);
        end
        if x(j,1) < 0 %setting left side boundary condition
            x(j,2) = L;
            x(j,1) = x(j,2) + (dt * Vx(j));
        end
    end
    Time = Time + dt;
    Temperature(1,2) = Temperature(1,2) + (mn*Vtotal(j)^2)/(2*C.k);
end

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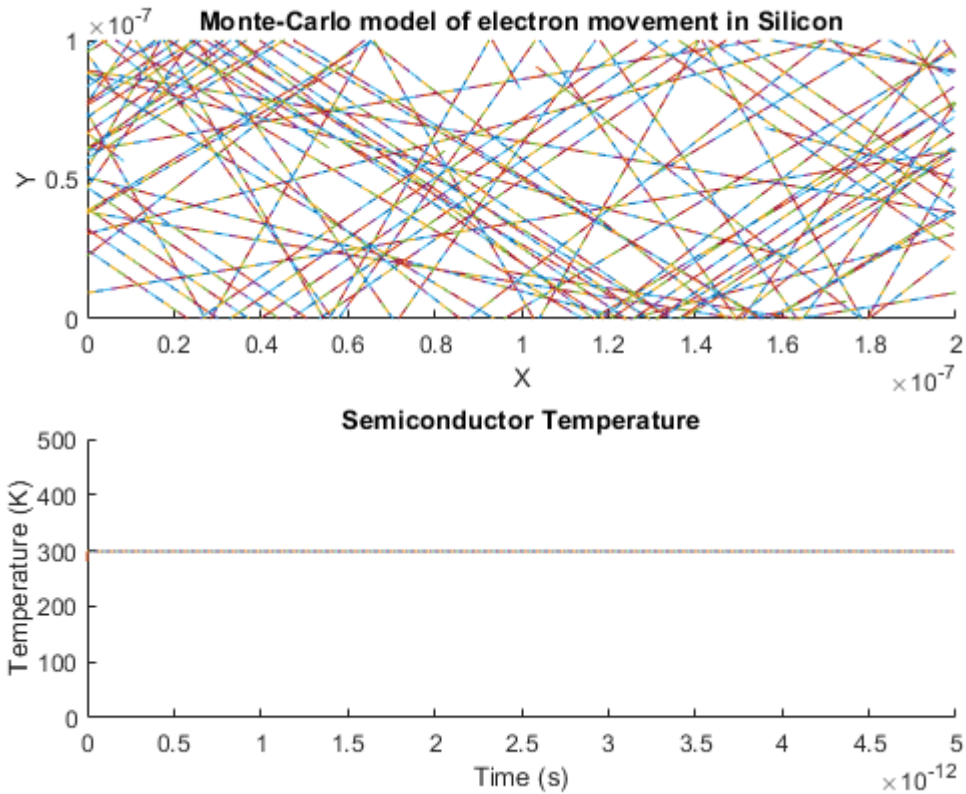
end
%setting roof/floor boundary condition
if y(j,1) > W || y(j,1) < 0
    Vy(j) = -Vy(j);
end
%setting up position line vectors to plot
XPlot = [x(j,2) x(j,1)];
YPlot = [y(j,2) y(j,1)];
if j < VisibleEcount
    %plot visible position line vectors
    plot(XPlot,YPlot);
end

%sum of x and y velocities
VTotal = sqrt(Vx(j)^2 + Vy(j)^2);

%sum of all temperatures
Temperature(1,2) = Temperature(1,2) + (mn*Vtotal(j)^2)/(2*C.k);

end
%averaging temperature sum and plotting against each timestep
AvgTemperature = Temperature(1,2)/Ecount;
TemperaturePlot = [Temperature(1,1) AvgTemperature];
TimePlot = [(Time - dt) Time];
subplot(2,1,2);
plot(TimePlot, TemperaturePlot);
Temperature(1,1) = AvgTemperature;
AvgTemperature = 0;
Temperature(1,2) = 0;
pause(1e-19)
Time = Time + dt;
end

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% % ELEC 4700 Assignment 1: Monte-Carlo Modelling of Electrons
% % Mike Crawford
% % 100952432
% % Question2
% This next part is similar to the previous, with the same model of electrons
% and average temperatures. The difference is it adds a collision factor
% with a scattering probability as well as a random velocity with each
% electron. A histogram showing the distribution of different velocities is
% also given. This part is using the same code skeleton as the first part
% so much of the code will look the same

%reset
clearvars
clearvars -GLOBAL
close all
format shortE
global C
global Ecount
global Vx Vy Vtotal x y
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Ecount = 1000; %setting amount of electrons used
C.mo = 9.10938215e-31;%Electron mass
C.k = 1.3806504e-23; %Boltzmann constant

T = 300; %Temperature (K)
mn = 0.26*C.mo;
L = 200e-9; %Length of Border
W = 100e-9; %Width of Border
Vth = sqrt((2*C.k*T)/mn); %Calculation of Thermal Velocity
dt = 10e-15; %timestep
Stop = 100*dt; %timeframe
x = zeros(Ecount, 2); %array of x positions
y = zeros(Ecount, 2); %array of y positions

Temperature = zeros(1,2); %array of temperatures
Time = 0;
VisibleEcount = 50; %visibbble electrons
tmn = 0.2e-12; %mean time between collisions
PScat = 1 - exp(-dt/tmn); %scattering probability equation
VTotalHistogram = zeros(Ecount, 1);%array of thermal velocities
for i = 1:Ecount
    x(i,1) = rand()*200e-9;%setting random electron positions
    y(i,1) = rand()*100e-9;
end
%setting random velocities
Vx(1:Ecount) = Vth*rand * cos(2*pi*randn(1, Ecount));
Vy(1:Ecount) = Vth*rand * sin(2*pi*randn(1, Ecount));
%sum of x and y velocities
for i = 1:Ecount
    Vtotal(i) = sqrt(Vx(i)^2 + Vy(i)^2);
    VTotalHistogram(i) = sqrt(Vx(i)^2 + Vy(i)^2);

end

%plot of electrons
figure(1)
subplot(2,1,1);
axis([0 L 0 W]);
title('Monte-Carlo model of electron movement in Silicon');
xlabel('X');
ylabel('Y');
hold on;
%plot of average temperatures
subplot(2,1,2);
axis([0 Stop 0 500]);
title('Semiconductor Temperature');
xlabel('Time (s)');
ylabel('Temperature (K)');
hold on;
%initial temperature sum

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for i = 1:Ecount
    Temperature(1,2) = Temperature(1,2) + (mn*Vtotal(i)^2)/(2*C.k);
end
%initial average temperature
AvgTemperature = Temperature(1,2)/Ecount;
TemperaturePlot = [300 AvgTemperature];
TimePlot = [0 Time];
plot(Time, AvgTemperature);
%resetting before loop
VTotal = 0;
Temperature(1,2) = 0;
AvgTemperature = 0;
%looping timestep
while Time < Stop
    subplot(2,1,1)
    %looping electrons
    for j = 1:Ecount
        %Scattering probability conditioned with random value to simulate
        %random scattering at set rate
        if PScat > rand
            Vx(j) = Vth * randn;
            Vy(j) = Vth * randn;

        end
        %update previous/new x and y positions
        x(j,2) = x(j,1);
        y(j,2) = y(j,1);
        x(j,1) = x(j,1) + (dt * Vx(j));
        y(j,1) = y(j,1) + (dt * Vy(j));
        %check right wall border
        if x(j,1) > L
            x(j,2) = 0;
            x(j,1) = dt * Vx(j);
        end
        %check left wall border
        if x(j,1) < 0
            x(j,2) = L;
            x(j,1) = x(j,2) + (dt * Vx(j));
        end
        %check roof/floor border
        if y(j,1) > W || y(j,1) < 0
            Vy(j) = -Vy(j);
        end
        %set line vectores for x and y positions
        XPlot = [x(j,2) x(j,1)];
        YPlot = [y(j,2) y(j,1)];
        %plot visible x and y line vectors
        if j < VisibleEcount
            plot(XPlot,YPlot);
        end
    end
end

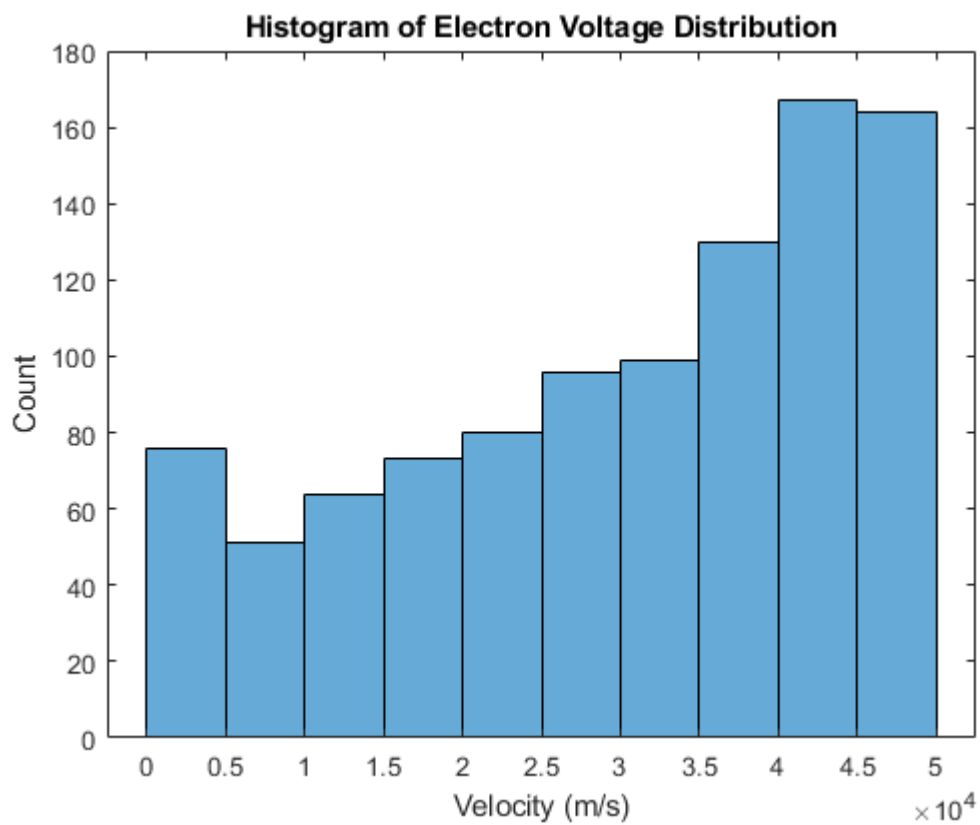
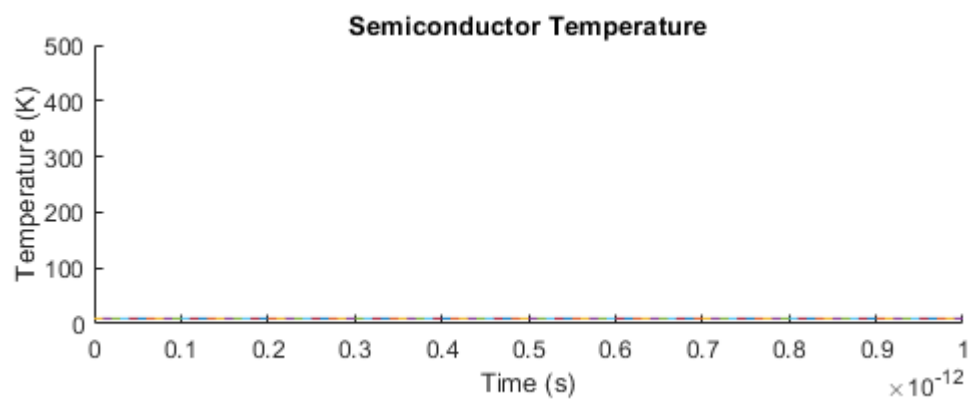
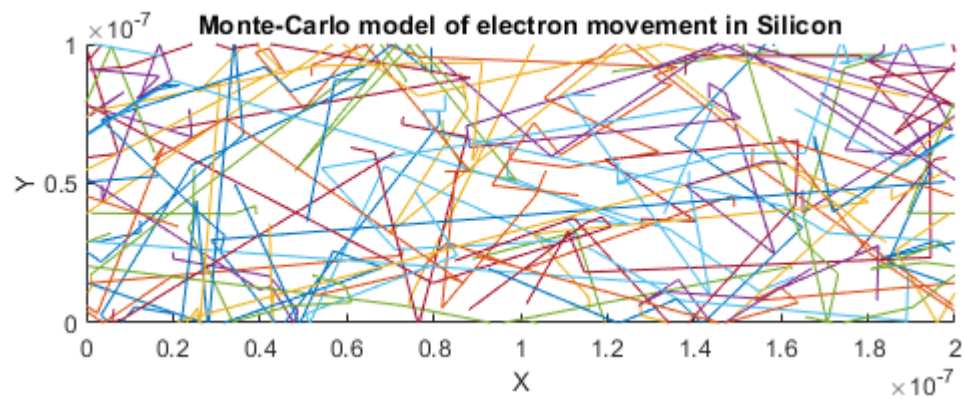
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    %sum of themal velocities and temperatures
    VTtotal = sqrt(Vx(j)^2 + Vy(j)^2);
    Temperature(1,2) = Temperature(1,2) + (mn*VTtotal(j)^2)/(2*C.k);

end
%plotting average temperature and resetting values for next timestep
AvgTemperature = Temperature(1,2)/Ecount;
TemperaturePlot = [Temperature(1,1) AvgTemperature];
TimePlot = [(Time - dt) Time];
subplot(2,1,2);
plot(TimePlot, TemperaturePlot);
Temperature(1,1) = AvgTemperature;
AvgTemperature = 0;
Temperature(1,2) = 0;
pause(1e-19)
Time = Time + dt;
end
%histogram of thermal velocities
figure(2)
histogram(VTtotalHistogram)
title('Histogram of Electron Voltage Distribution')
xlabel('Velocity (m/s)')
ylabel('Count')

```




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% % ELEC 4700 Assignment 1: Monte-Carlo Modelling of Electrons
% % Mike Crawford
% % 100952432
% % Question3
% This final part add onto the previous by adding two boxes that from a
% bottleneck inside the borders of the simulation. These boxes should reflect
% any electrons coliding into it and stop elctrons from beign positioned
% inside of it. The boxes should be able to ither be specular or diffusive.
% Meaning after an electron bounces, its either given the reflected velocity
% a random velocity respectively. After the simulation is completed, a map
% of both the electron density and the temperature will be displayed. \

%reset
clearvars
clearvars -GLOBAL
close all
format shortE
global C
global Ecount
global Vx Vy Vtotal x y
Ecount =1000; %setting amount of electrons used
C.mo = 9.10938215e-31; %Electron mass
C.k = 1.3806504e-23; %Boltzmann constant

T =300; %Temperature (K)
mn = 0.26*C.mo;
L = 200e-9; %Length of Border
W = 100e-9; %Width of Border
Vth = sqrt((2*C.k*T)/mn); %Calculation of Thermal Velocity
dt = 10e-15; %timestep
Stop = 100*dt; %timeframe
x = zeros(Ecount, 2); %array of x positions
y = zeros(Ecount, 2); %array of y positions
Temperature = zeros(1,2); %array of temperatures
Time = 0;
VisibleEcount = 50; %visibbble electrons
tmn = 0.2e-12; %mean time between collisions
PScat = 1 - exp(-dt/tmn); %scattering probability equation
VTotalHistogram = zeros(Ecount, 1); %array of thermal velocities
DrawBoxX = [80e-9 80e-9 120e-9 120e-9 80e-9]; %X border limits of boxes
%(Same X values for both)
DrawBoxY1 = [100e-9 60e-9 60e-9 100e-9 100e-9]; %Y border limits of top box
DrawBoxY2 = [40e-9 0 0 40e-9 40e-9]; %Y border limits of bottom box
Specular = true; %boolean indicating whether specular or diffusive
InsideBox = true; %boolean showing if electron spawned inside boxes
MappingStep = 10e-9; %size of each suare used on mapping grid
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%setting arrays of density and temperature based off border demesions and
%mapping step
DensityMap = zeros(W/MappingStep, L/MappingStep);
TemperatureMap = zeros(W/MappingStep, L/MappingStep);
for i = 1:Ecount
    x(i,1) = rand()*200e-9; %setting random electron positions
    y(i,1) = rand()*100e-9;
    InsideBox = true;
    %checks if set electron is inside box, if so reinitialize its position
    while InsideBox == true
        if (x(i) >= 40e-9 && x(i) <= 120e-9) && (y(i) >= 60e-9 ||...
            y(i) <= 40e-9)
            x(i,1) = rand * 200e-9;
            y(i,1) = rand * 100e-9;
        else
            InsideBox = false;
        end
    end
end

end

%set random velocities
Vx(1:Ecount) = Vth*rand * cos(2*pi*randn(1, Ecount));
Vy(1:Ecount) = Vth*rand * sin(2*pi*randn(1, Ecount));
%sum of x and y velocities
for i = 1:Ecount
    Vtotal(i) = sqrt(Vx(i)^2 + Vy(i)^2);
    VTotalHistogram(i) = sqrt(Vx(i)^2 + Vy(i)^2);

end

%plot of electrons with boxes
figure(1)
subplot(2,1,1);
plot(DrawBoxX, DrawBoxY1, DrawBoxX, DrawBoxY2)
axis([0 L 0 W]);
title('Monte-Carlo model of electron movement in Silicon');
xlabel('X');
ylabel('Y');
hold on;
%set up plot for temperatures
subplot(2,1,2);
axis([0 Stop 0 500]);
title('Semiconductor Temperature');
xlabel('Time (s)');
ylabel('Temperature (K)');
hold on;
%summing temperatures
for i = 1:Ecount
    Temperature(1,2) = Temperature(1,2) + (mn*Vtotal(i)^2)/(2*C.k);
end

```

```

%averaging temperatures and plotting over timestep
AvgTemperature = Temperature(1,2)/Ecount;
TemperaturePlot = [300 AvgTemperature];
TimePlot = [0 Time];
plot(Time, AvgTemperature);
%reset variables before loop
VTotal = 0;
Temperature(1,2) = 0;
AvgTemperature = 0;
%looping timestep
while Time < Stop
    subplot(2,1,1)
    %looping electrons
    for j = 1:Ecount
        %set to assume leaking to start condition for diffusive bounces
        leaking = true;
        %scattering probability condition
        if PScat > rand
            Vx(j) = Vth * randn;
            Vy(j) = Vth * randn;

        end
        %updating old/new x and y positions
        x(j,2) = x(j,1);
        y(j,2) = y(j,1);
        x(j,1) = x(j,1) + (dt * Vx(j));
        y(j,1) = y(j,1) + (dt * Vy(j));
        %checking if collision with top box
        if (x(j,1) >= 80e-9 && x(j,1) <= 120e-9) && y(j,1) >= 60e-9
            %checking which side of box collision came from
            %if left side, reflect X velocity and reset positions outside
            %of box
            if y(j,2) < 60e-9
                Vy(j) = -Vy(j);
                y(j,1) = 60e-9;
                y(j,2) = 60e-9;
                %if right side, reflect X velocity and reset position
                %outside of box
            elseif x(j,2) < 80e-9
                Vx(j) = -Vx(j);
                x(j,1) = 80e-9;
                x(j,2) = 80e-9;
                %if bottom side, reflect Y velocity and reset position
                %outside of box
            elseif x(j,2) > 120e-9
                Vx(j) = -Vx(j);
                x(j,1) = 120e-9;
                x(j,2) = 120e-9;
            end
            %check if specular or diffusive collisions
            if Specular == true

```

```

        %if specular, simply update positions
        x(j,1) = x(j,2) + Vx(j)*dt;
        y(j,1) = y(j,2) + Vy(j)*dt;
    else
        %if diffusive, update velocities with random value
        Vx(j) = Vth * randn;
        Vy(j) = Vth * randn;
        %assume random velocity will direct it inside the box
        while leaking == true
            %check if new velocity is directed towards the box by
            %examining position point and new velocity
            if(x(j,2) < 80e-9 && Vx(j) >= 0) || ...
                (x(j,2) > 120e-9 && Vx(j) <= 0) || ...
                (y(j,2) < 60e-9 && Vy(j) >= 0)
                %if new velocity conflicts with boxes, reinitialize
                %the velocity
                Vx(j) = Vth * randn;
                Vy(j) = Vth * randn;
            else
                leaking = false;
            end
        end
        %update positions with random velocity
        x(j,1) = x(j,2) + Vx(j)*dt;
        y(j,1) = y(j,2) + Vy(j)*dt;
    end
end

%checking if collision with bottom box
if (x(j,1) >= 80e-9 && x(j,1) <= 120e-9) && y(j,1) <= 40e-9
    %checking which side of box collision came from
    %if top side, reflect Y velocity and reset positions outside
    %of box
    if y(j,2) > 40e-9
        Vy(j) = -Vy(j);
        y(j,1) = 40e-9;
        y(j,2) = 40e-9;
        %if left side, reflect X velocity and reset positions outside
        %of box
        elseif x(j,2) < 80e-9
            Vx(j) = -Vx(j);
            x(j,1) = 80e-9;
            x(j,2) = 80e-9;
            %if right side, reflect X velocity and reset position
            %outside of box
            elseif x(j,2) > 120e-9
                Vx(j) = -Vx(j);
                x(j,1) = 120e-9;
                x(j,2) = 120e-9;
            end
            %check if specular or diffusive collisions

```

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    if Specular == true
        %if specular, simply update positions
        x(j,1) = x(j,2) + Vx(j)*dt;
        y(j,1) = y(j,2) + Vy(j)*dt;
    else
        %if diffusive, update velocities with random value
        Vx(j) = Vth * randn;
        Vy(j) = Vth * randn;
        %assume random velocity will direct it inside the box
        while leaking == true
            %check if new velocity is directed towards the box by
            %examining position point and new velocity
            if (x(j,2) < 80e-9 && Vx(j) >= 0) || ...
                (x(j,2) > 120e-9 && Vx(j) <= 0) || ...
                (y(j,2) > 40e-9 && Vy(j) <= 0)
                %if new velocity conflicts with boxes, reinitialize
                %the velocity
                Vx(j) = Vth * randn;
                Vy(j) = Vth * randn;
            else
                leaking = false;
            end
        end
        %update positions with random velocity
        x(j,1) = x(j,2) + Vx(j)*dt;
        y(j,1) = y(j,2) + Vy(j)*dt;
    end
end

%check right wall border
if x(j,1) > L
    x(j,2) = 0;
    x(j,1) = dt * Vx(j);
end

%check left wall border
if x(j,1) < 0
    x(j,2) = L;
    x(j,1) = x(j,2) + (dt * Vx(j));
end

%check roof/floor border
if y(j,1) > W || y(j,1) < 0
    Vy(j) = -Vy(j);
end

%set line vectores for x and y positions
XPlot = [x(j,2) x(j,1)];
YPlot = [y(j,2) y(j,1)];
%plot visible x and y line vectors
if j < VisibleEcount
    plot(XPlot,YPlot);
end

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        %sum of themal velocities and temperatures
        VTotal = sqrt(Vx(j)^2 + Vy(j)^2);
        Temperature(1,2) = Temperature(1,2) + (mn*Vtotal(j)^2)/(2*C.k);

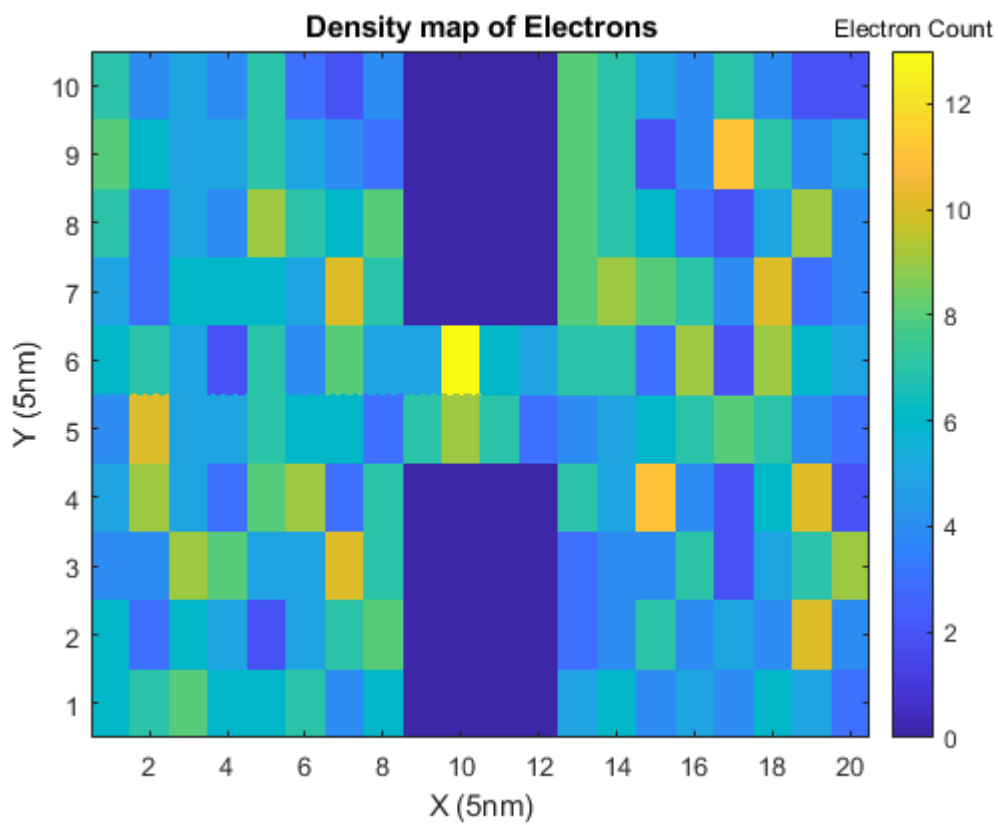
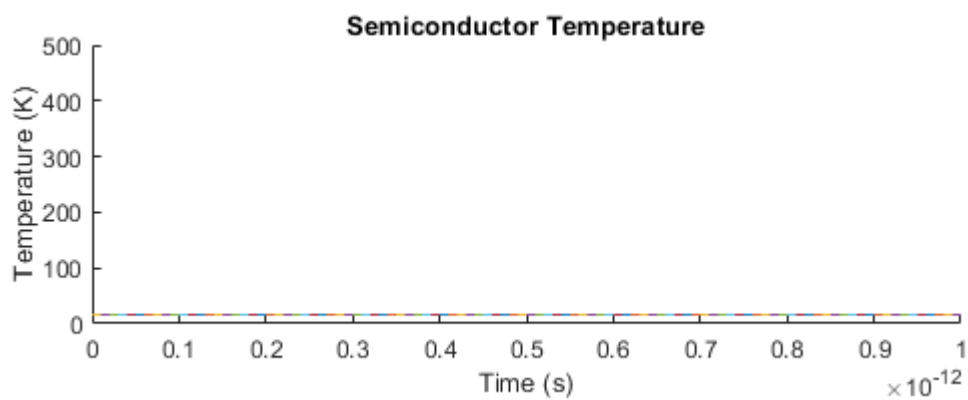
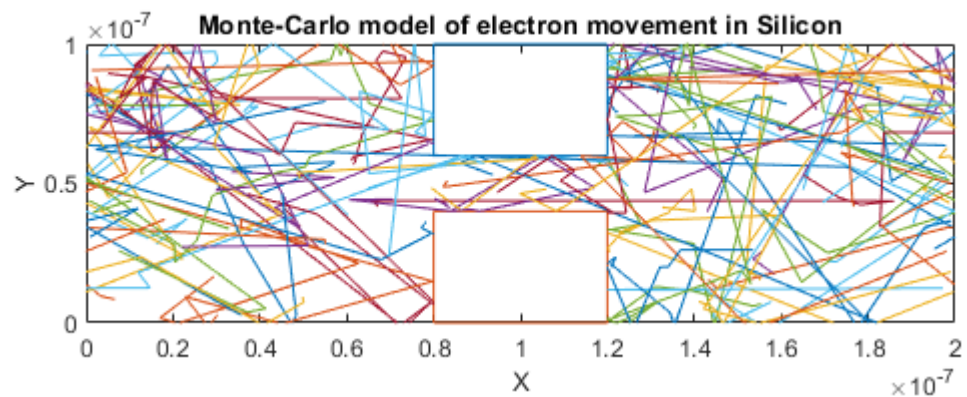
    end
    %plotting average temperature and ressetting values for next timestep
    AvgTemperature = Temperature(1,2)/Ecount;
    TemperaturePlot = [Temperature(1,1) AvgTemperature];
    TimePlot = [(Time - dt) Time];
    subplot(2,1,2);
    plot(TimePlot, TemperaturePlot);
    Temperature(1,1) = AvgTemperature;
    AvgTemperature = 0;
    Temperature(1,2) = 0;
    pause(1e-19)
    Time = Time + dt;
end
%looping through X indexes of the Mappingstep grid
for i = 1:(L/MappingStep)
    %looping through Y indexes of the Mapping grid
    for j = 1:(W/MappingStep)
        %looping through electrons
        for m = 1:Ecount
            %Checks to see if the electons position is with the map grip
            %being indexed. does this by checking all 4 sides of the
            %square grid index
            if(x(m,1) > MappingStep*(i -1)) && ...
                (x(m,1) < MappingStep*(i)) && ...
                (y(m,1) > MappingStep*(j - 1)) && ...
                (y(m,1) < MappingStep*(j))
                %sum of x and y velocities
                Vtotal(m) = sqrt(Vx(m)^2 + Vy(m)^2);
                %counts increased on the indexed location of both Maps
                DensityMap(j, i) = DensityMap(j, i) + 1;
                TemperatureMap(j, i) = TemperatureMap(j,i) + ...
                    (mn*Vtotal(m)^2)/(2*C.k);
            end
            %Averages the temperature values given in each grid index
            TemperatureMap(j,i) = TemperatureMap(j,i)/DensityMap(j,i);
        end
    end
end

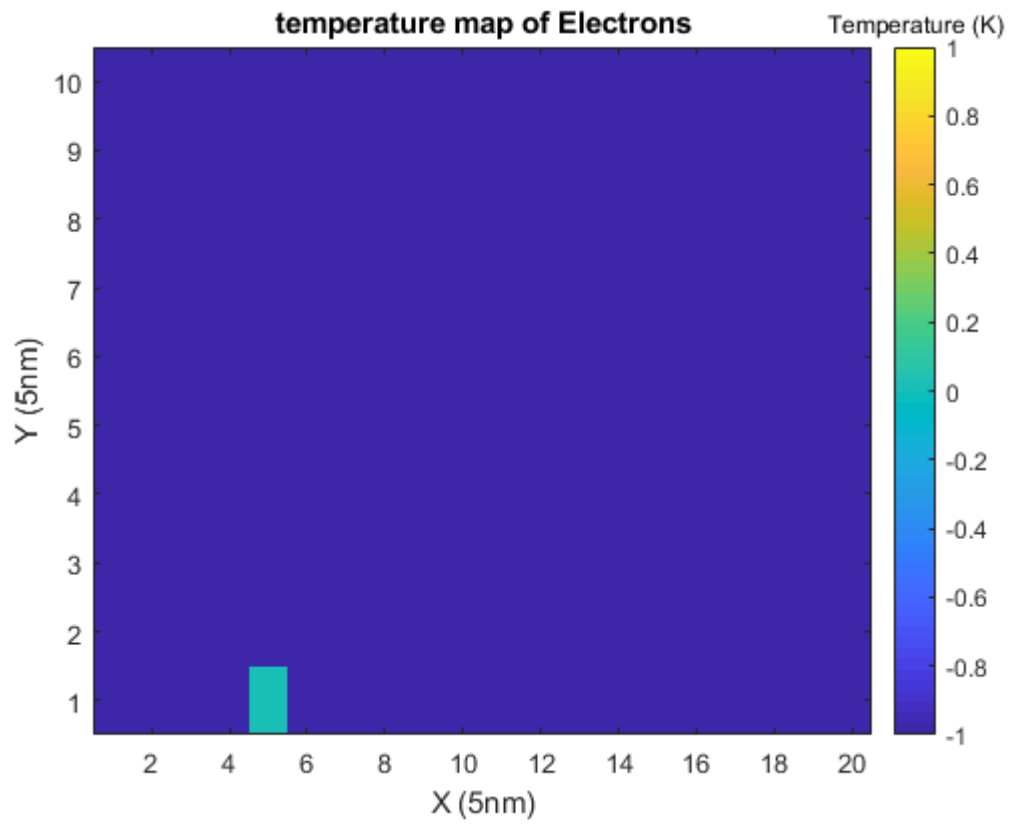
figure(2)
imagesc(DensityMap)
title('Density map of Electrons')
xlabel('X (5nm)');
ylabel('Y (5nm)');
set(gca, 'Ydir', 'Normal')

```

```
c = colorbar;
title(c, 'Electron Count')

figure(3)
imagesc(TemperatureMap)
title('temperature map of Electrons');
xlabel('X (5nm)')
ylabel('Y (5nm)')
set(gca, 'Ydir', 'Normal')
c = colorbar;
title(c, 'Temperature (K)')
```





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