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# Assignment 3

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## Question 1

```
close all;
clear all;
Q1 = true;
Q2 = true;
Q3 = true;

xlimits = [0, 2e-7];
ylimits = [0, 1e-7];
PlotHowMany = 10;
Timestep = 5e-15;
NumParticles = floor((xlimits(2) - xlimits(1))*(ylimits(2) -
    ylimits(1))*10^19);
T = 300;
Kbolt = 1.38064852E-23;
tmn = 0.2e-12;
Colomb = -1.602e-19;
mo = 9.11e-31;
mn = mo * 0.26;
mycolors = hsv(PlotHowMany);
vx = (Kbolt*T/mn)^.5;
vy = (Kbolt*T/mn)^.5;
vth = (vx^2 + vy^2)^.5;
MeanFreePath = tmn*vth;
Pscat = 1-exp(-Timestep/tmn);

if Q1 == true
```

## A - Electric Field on Electrons

```
voltDiffx = 0.1;
voltDiffy = 0;
xLeng = 50;
yLeng = 50;

%Part 1 A)
% The Electric Field in the X Direction is
Efieldx = voltDiffx/(xlimits(2) - xlimits(1))
% The Electric Field in the Y Direction is
Efieldy = voltDiffy/(ylimits(2) - ylimits(1))

% In units of V/m.

Efieldx =

    5.0000e+05

Efieldy =

    0
```

## B - Force on the Electrons

```
%Part 1 B)
% The Force in the X Direction is
Forcex = Efieldx*Colomb
% The Force in the Y Direction is
Forcey = Efieldy*Colomb

% In units of VC/m.

Forcex =

   -8.0100e-14

Forcey =

    0
```

## C - Acceleration on the Electrons

```
%Part 1 C)
```

```

% The Acceleration in the X Direction is
Accelx = Forcex/mn
% The Acceleration in the Y Direction is
Accely = Forcey/mn

% In units of m/s.

xprev = zeros(1, NumParticles);
yprev = zeros(1, NumParticles);
x = zeros(3, NumParticles);
y = zeros(2, NumParticles);
temp = zeros(2, NumParticles);
scatTime = zeros(1, NumParticles);
endtime = Timestep*500;
graphfor = Timestep*500;

%Start the random distribution in x position
x(1,:) = xlimits(1) + rand(1, NumParticles).*(xlimits(2) -
    xlimits(1));
y(1,:) = ylimits(1) + rand(1, NumParticles).*(ylimits(2) -
    ylimits(1));

temp(1,:) = ((sqrt(Kbolt*T/mn)*randn(1, NumParticles)).^2 +
    (sqrt(Kbolt*T/mn)*randn(1, NumParticles)).^2).^0.5;
x(3,:) = rand(1, NumParticles)*2*pi;
x(2,:) = temp(1,:).*cos(x(3,:));
y(2,:) = temp(1,:).*sin(x(3,:));

figure
%Update the Position
ScatterTime = zeros(floor(endtime/Timestep)+ 1,1);
Current = zeros(floor(endtime/Timestep)+ 1,1);
Temperature = zeros(floor(endtime/Timestep)+ 1,1);
MFP = zeros(floor(endtime/Timestep)+ 1,1);
for i = 0:Timestep:endtime
    count = 0;
    xprev(1,:) = x(1,:);
    yprev(1,:) = y(1,:);
    for kt = 1:NumParticles
        if x(1,kt) + x(2,kt)*Timestep < xlimits(1)
            x(1,kt) = xlimits(2);
            xprev(1,kt) = xlimits(2);
            count = count - 1;
        elseif x(1,kt) + x(2,kt)*Timestep > xlimits(2)
            x(1,kt) = xlimits(1);
            xprev(1,kt) = xlimits(1);
            count = count + 1;
        end
        if y(1,kt) + y(2,kt)*Timestep < ylimits(1) || y(1,kt) +
            y(2,kt)*Timestep > ylimits(2)
            y(2,kt) = -y(2,kt);
        end
        x(1,kt) = x(1,kt) + x(2,kt).*Timestep;
    end
end

```

```

y(1,kt) = y(1,kt) + y(2,kt).*Timestep;
if kt <= PlotHowMany && i < graphfor
    plot([xprev(1,kt), x(1,kt)], [yprev(1,kt), y(1,kt)], 'color',
mycolors(kt,:) );
    hold on;
end
%Scattering Check
if Pscat>rand()
    temp = (normrnd(0, sqrt(Kbolt*T/mn)).^2 + normrnd(0,
sqrt(Kbolt*T/mn)).^2).^5;
    x(3,kt) = rand*2*pi;
    x(2,kt) = temp*cos(x(3,kt));
    y(2,kt) = temp*sin(x(3,kt));
    scatTime(1,kt) = 0;
end
end
hold on;
AvgScat = mean(scatTime(1,:));
scatTime(1,:) = scatTime(1,:) + Timestep;
VelSquared = mean((x(2,:).^2 + y(2,:).^2));
CalcTemp = VelSquared*mn/2/Kbolt;
%title(['Average Temperature: ' num2str(CalcTemp) ' Average
Scatter Time: ' num2str(AvgScat)]);
xlim([xlimits(1), xlimits(2)]);
ylim([ylimits(1), ylimits(2)]);
ScatterTime(round(i/Timestep) + 1,1) = AvgScat;
MFP(round(i/Timestep) + 1,1) = VelSquared^.5*AvgScat;
Temperature(round(i/Timestep) + 1,1) = CalcTemp;
x(2,:) = x(2,:) + Accelx*Timestep;
y(2,:) = y(2,:) + Accely*Timestep;
Current(round(i/Timestep) + 1,1) = Colomb*count/Timestep;
end
title('Question 1 - Electron Position Plot');
xlabel('X Position (m)');
ylabel('Y Position (m)');
hold off;

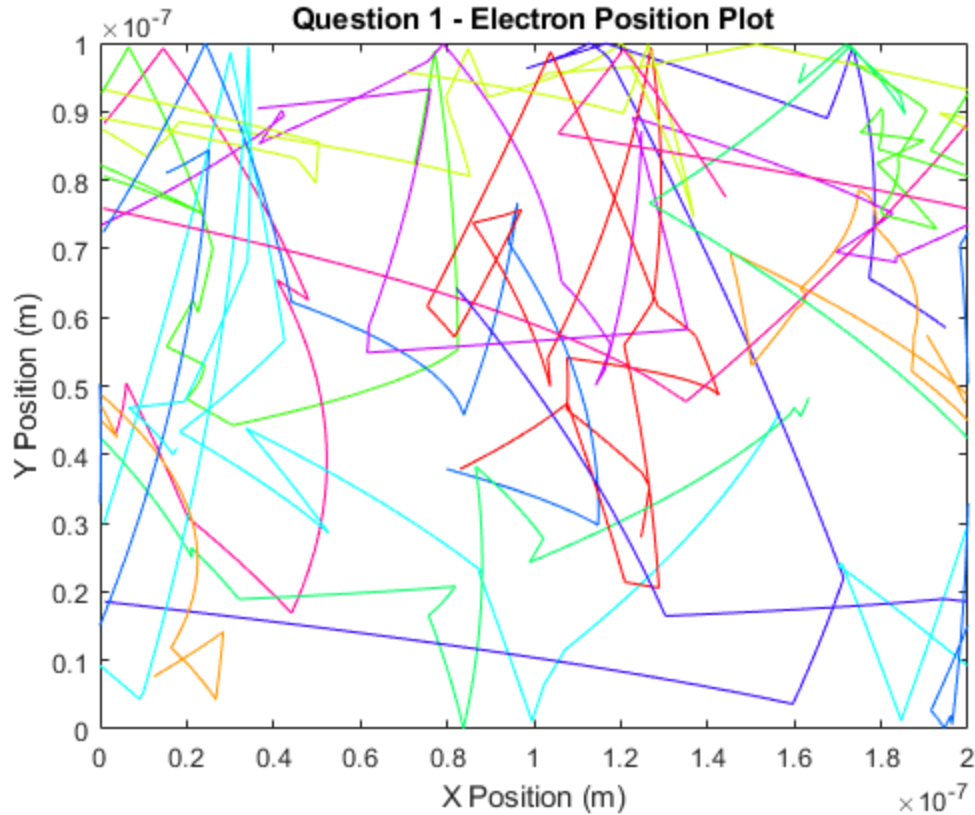
```

Accelx =

-3.3817e+17

Accely =

0



## D - Relationship between Electron Drift and Average Carrier Velocity

The electron drift current density is given by:

$$J = q\rho\mu E$$

where  $q$  is the electron charge,  $\rho$  is the resistivity,  $\mu$  is the electron mobility and  $E$  is the electric field applied. This equation represents the drift current and then the average velocity of the particle is given by:

$$v_d = \mu E$$

where  $v_d$  is the average velocity, and  $\mu$  is electron mobility and  $E$  is electric field again. By combining these two we can create an equation of current density in terms of average velocity:

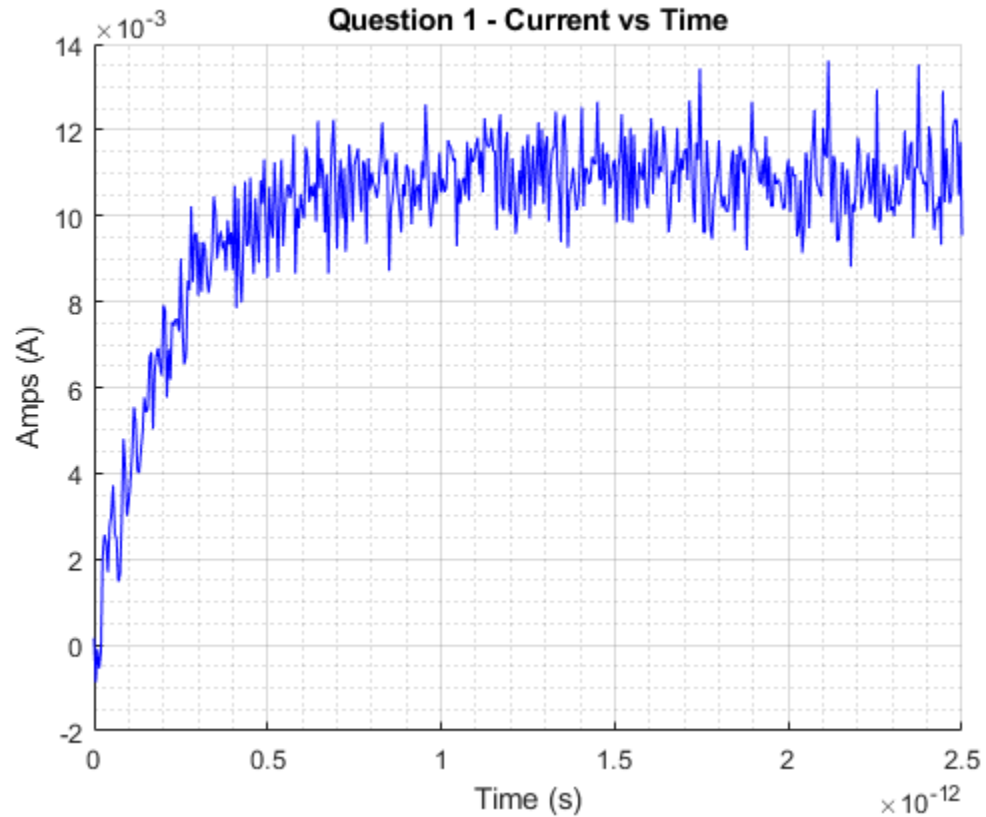
$$J = q\rho v_d$$

The calculation of current can be done by multiplying the current density by the cross sectional area.

$$I = JA = Aq\rho v_d$$

figure;

```
title('Question 1 - Current vs Time');
xlabel('Time (s)');
ylabel('Amps (A)');
hold on;
plot(0:Timestep:endtime, Current(:, 1), 'b');
grid on;
grid minor;
hold off;
```



The current starts at zero or very small, this is due to the fact that only the electrons that are moving randomly out of the box through the x axis limits are counting towards current flow. Then as more acceleration is added to the electrons with more time steps; more begin to accelerate over the boundary in the direction by the electric field. This approaches a limit and is somewhat stabilizes with some fluctuation due to the inherent randomness with diffusion and scattering.

## E - Electron Density Map and Temperature Map

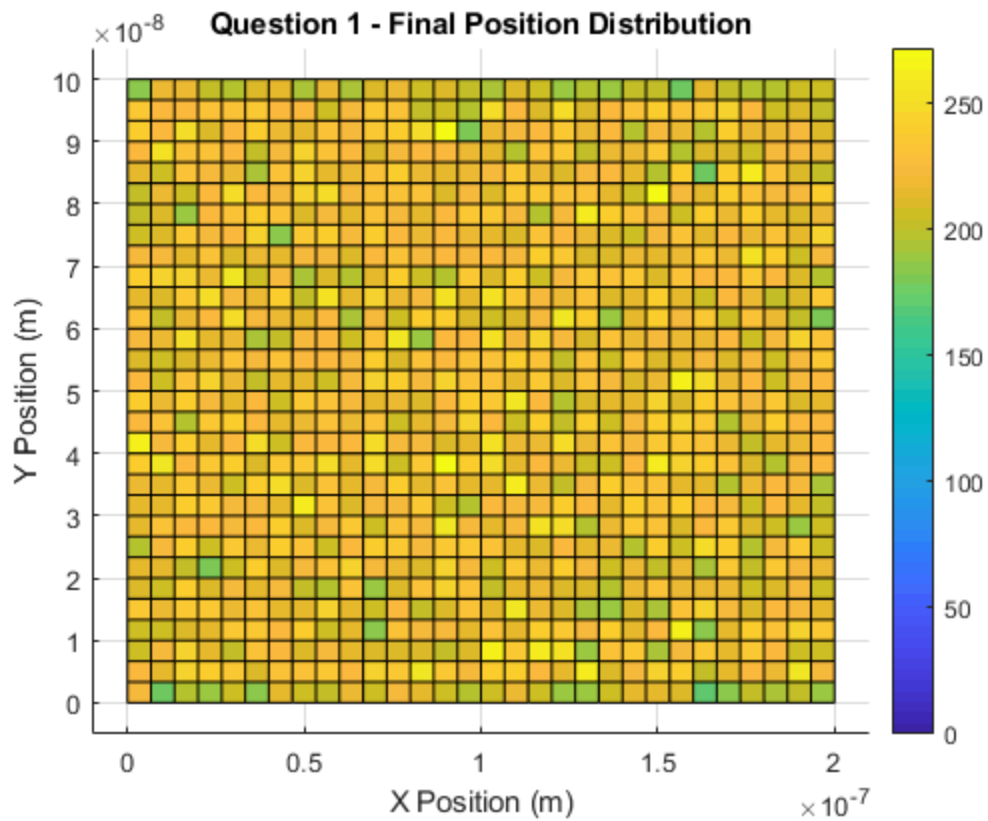
```
figure;
Z = [transpose(x(1,:)), transpose(y(1,:))];
hist3(Z, 'Nbins',[30 30], 'CdataMode', 'auto');
hold on;
colorbar;
view(2);
grid on;
title('Question 1 - Final Position Distribution');
xlabel('X Position (m)');
```

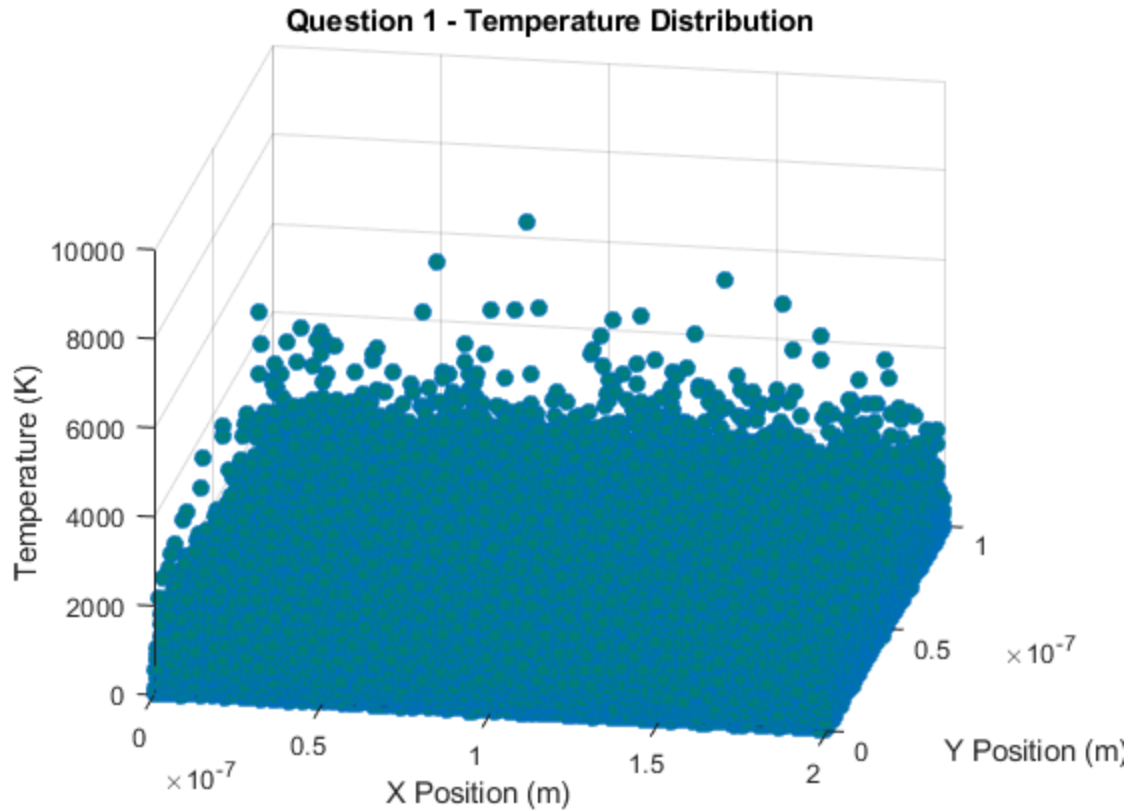
```

ylabel('Y Position (m)');
hold off;

figure
VelSquared = (x(2,:).^2 + y(2,:).^2);
CalcTemp = VelSquared.*mn./2./Kbolt;
h = scatter3(x(1,:), y(1,:), CalcTemp(1,:));
h.MarkerFaceColor = [0 0.5 0.5];
grid on;
xlabel('X Position (m)');
ylabel('Y Position (m)');
xlim([xlims(1), xlims(2)]);
ylim([ylims(1), ylims(2)]);
zlabel('Temperature (K)');
title('Question 1 - Temperature Distribution');
view(10,25);
hold off;

```





end

## Question 2

```

if Q2 == true

v0 = 1;
L = xlimits(2)- xlimits(1);
W = ylimits(2)- ylimits(1);
nx =75;
ny= 50;
wb = W/3;
lb = L/3;
map = @(j,i) j + (i-1)*ny;
sigma = ones(ny,nx);
for i = 1:nx
if i > (nx/L)*(L-lb)/2 && i < (nx/L)*(L+lb)/2
    for j = 1:ny
        if j < (ny/W)*wb || j> ceil((ny/W)*(W-wb))
            sigma(j,i) = 0.01;
        end
    end
end
end
end
end

```



```
for i = 1:nx
for j = 1:ny
    n = map(j,i);
    if i==1
        B(n) = v0;
        G(n,:) = 0;
        G(n,n) = 1;
    elseif i == nx
        B(n) = 0;
        G(n,:) = 0;
        G(n,n) = 1;
    elseif j ==1
        up = (sigma(j,i) + sigma(j+1,i))/2;
        left = (sigma(j,i) + sigma(j,i-1))/2;
        right = (sigma(j,i) + sigma(j,i+1))/2;

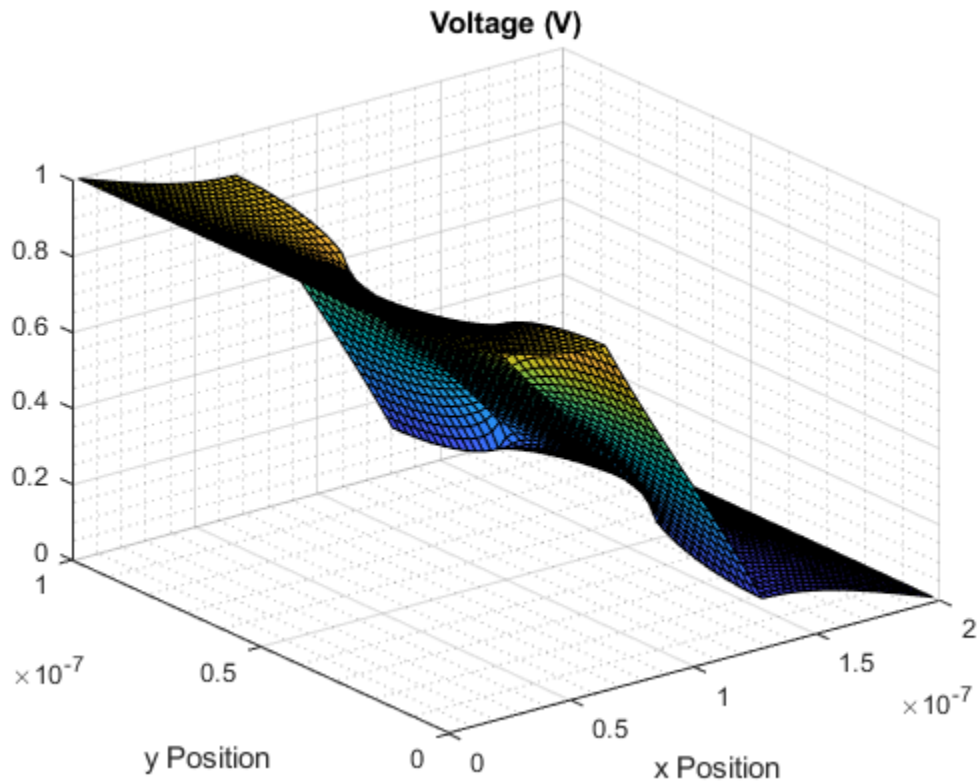
        G(n,n) = -(up+left+right);
        G(n,map(j,i-1)) = left;
        G(n,map(j,i+1)) = right;
        G(n,map(j+1,i)) = up;
    elseif j == ny
        left = (sigma(j,i) + sigma(j,i-1))/2;
        right = (sigma(j,i) + sigma(j,i+1))/2;
        down = (sigma(j,i) + sigma(j-1,i))/2;

        G(n,n) = -(down+left+right);
        G(n,map(j,i-1)) = left;
        G(n,map(j,i+1)) = right;
        G(n,map(j-1,i)) = down;
    else
        up = (sigma(j,i) + sigma(j+1,i))/2;
        left = (sigma(j,i) + sigma(j,i-1))/2;
        right = (sigma(j,i) + sigma(j,i+1))/2;
        down = (sigma(j,i) + sigma(j-1,i))/2;

        G(n,n) = -(up+left+right+down);
        G(n,map(j,i-1)) = left;
        G(n,map(j,i+1)) = right;
        G(n,map(j+1,i)) = up;
        G(n,map(j-1,i)) = down;
    end
end
end
E = G\B';
d = zeros(ny,nx);
for i = 1:nx
    for j = 1:ny
        n = map(j,i);
        d(j,i) = E(n);
    end
end
end
```

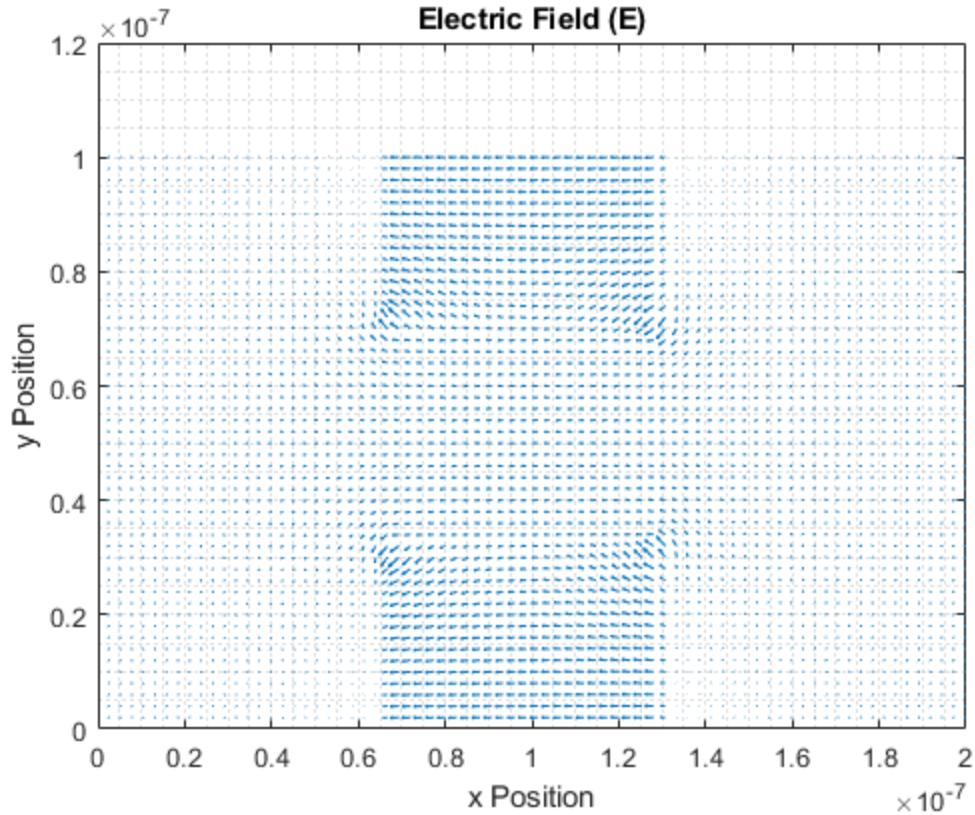
## A - Surface Plot of the bottleneck Voltage

```
figure
[X, Y] = meshgrid(L/nx:L/nx:L, W/ny:W/ny:W);
surf(X, Y, d);
grid minor;
title('Voltage (V)');
xlabel('x Position');
ylabel('y Position');
```



## B - Electric Field Vector Plot

```
[Ex, Ey] = gradient(d, L/nx);
figure
quiver(X,Y,Ex, Ey);
title('Electric Field (E)');
grid minor;
xlabel('x Position');
ylabel('y Position');
```



end

## Question 3

```
if Q3 == true

NumParticles = 100000;

xboxLim = [.4*(xlims(2)-xlims(1)), .6*(xlims(2)-xlims(1))];
yboxLim1 = [ylims(1), .4*(ylims(2)-ylims(1))];
yboxLim2 = [.6*(ylims(2)-ylims(1)), ylims(2)];
xbox = xboxLim([1 1 2 2 1]);
ybox1 = yboxLim1([1 2 2 1 1]);
ybox2 = yboxLim2([1 2 2 1 1]);
endtime = Timestep*1000;
xprev = zeros(1, NumParticles);
yprev = zeros(1, NumParticles);
x = zeros(3, NumParticles);
y = zeros(2, NumParticles);
temp = zeros(2, NumParticles);
scatTime = zeros(1, NumParticles);

%Start the random distribution in x position
x(1,:) = rand(1, NumParticles);
y(1,:) = rand(1, NumParticles);
x(1,:) = xlims(1) + x(1,:).*(xlims(2) - xlims(1));
```

```

y(1,:) = ylimits(1) + y(1,:).*(ylimits(2) - ylimits(1));

%Get the Particle Indexes that are out of bounds
IDX = uint32(1:NumParticles);
ind = IDX((y(1,:) >= yboxLim2(1) | y(1,:) <= yboxLim1(2)) & x(1,:) >=
    xboxLim(1) & x(1,:) <= xboxLim(2));

%Reassign Positions
counter = 1;
%size(ind, 2)
while counter <= size(ind,2)
    x(1,ind(counter)) = xlimits(1) + rand*(xlimits(2) - xlimits(1));
    y(1,ind(counter)) = ylimits(1) + rand*(ylimits(2) - ylimits(1));
    if ~(y(1,ind(counter)) >= yboxLim2(1) || y(1,ind(counter)) <=
        yboxLim1(2)) && x(1,ind(counter)) >= xboxLim(1) && x(1,ind(counter))
        <= xboxLim(2))
        counter = counter + 1;
    end
end

%Assign the random velocity and random angle
temp(1,:) = ((sqrt(Kbolt*T/mn)*randn(1, NumParticles)).^2 +
    (sqrt(Kbolt*T/mn)*randn(1, NumParticles)).^2).^5;
x(3,:) = rand(1, NumParticles)*2*pi;
x(2,:) = temp(1,:).*cos(x(3,:));
y(2,:) = temp(1,:).*sin(x(3,:));

DiffusionBarrierProbability = 0.2;

```

## A - Plot of Particle Trajectories

```

figure;
title('Question 3 - Electron Position');
xlabel('X Position (m)');
ylabel('Y Position (m)');
xlim([xlimits(1), xlimits(2)]);
ylim([ylimits(1), ylimits(2)]);
hold on;
rectangle('Position', [xboxLim(1), yboxLim1(1), xboxLim(2)-xboxLim(1),
    yboxLim1(2)]);
rectangle('Position', [xboxLim(1), yboxLim2(1), xboxLim(2)-xboxLim(1),
    yboxLim2(2) - yboxLim2(1)]);

counter = 1;
for i = 0:Timestep:endtime
    pause(.1);
    xprev(1,:) = x(1,:);
    yprev(1,:) = y(1,:);
    for kt = 1:NumParticles

        %PARTICLE'S HITTING TOP and
        if x(1,kt) + x(2,kt)*Timestep < xlimits(1)
            x(1,kt) = xlimits(2);

```

```

        xprev(1,kt) = xlimits(2);
elseif x(1,kt) + x(2,kt)*Timestep > xlimits(2)
    x(1,kt) = xlimits(1);
    xprev(1,kt) = xlimits(1);
end
if y(1,kt) + y(2,kt)*Timestep < ylimits(1) || y(1,kt) +
y(2,kt)*Timestep > ylimits(2)
    if rand > DiffusionBarrierProbability
        y(2,kt) = -y(2,kt);
    else
        if y(1,kt) + y(2,kt)*Timestep < ylimits(1)
            x(3,kt) = rand*pi;
        elseif y(1,kt) + y(2,kt)*Timestep > ylimits(2)
            x(3,kt) = -rand*pi;
        end
        temp = (normrnd(0, sqrt(Kbolt*T/mn))^2 + normrnd(0,
sqrt(Kbolt*T/mn))^2)^.5;
        x(2,kt) = temp*cos(x(3,kt));
        y(2,kt) = temp*sin(x(3,kt));
    end
end

%PARTICLE'S HITTING BOX
if (y(1,kt) + y(2,kt)*Timestep >= yboxLim2(1) && x(1,kt)
+ x(2,kt)*Timestep >= xboxLim(1) && x(1,kt) + x(2,kt)*Timestep <=
xboxLim(2))
    [xinter1, xinter2, yinter1, yinter2] =
BoxIntercept( x(1,kt), y(1,kt), x(1,kt) + x(2,kt)*Timestep, y(1,kt)
+ y(2,kt)*Timestep, xboxLim(1), xboxLim(2), yboxLim2(1), ylimits(2));
    if rand > DiffusionBarrierProbability
        if xinter2 || xinter1
            x(2,kt) = - x(2,kt);
        end
        if yinter2 || yinter1
            y(2,kt) = - y(2,kt);
        end
    else
        if xinter1
            x(3,kt) = pi/2 + rand*pi;
        elseif xinter2
            x(3,kt) = pi/2 - rand*pi;
        elseif yinter2
            x(3,kt) = rand*pi;
        elseif yinter1
            x(3,kt) = -rand*pi;
        end
        temp = (normrnd(0, sqrt(Kbolt*T/mn))^2 +
normrnd(0, sqrt(Kbolt*T/mn))^2)^.5;
        x(2,kt) = temp*cos(x(3,kt));
        y(2,kt) = temp*sin(x(3,kt));
    end
elseif y(1,kt) + y(2,kt)*Timestep <= yboxLim1(2) && x(1,kt)
+ x(2,kt)*Timestep >= xboxLim(1) && x(1,kt) + x(2,kt)*Timestep <=
xboxLim(2)

```

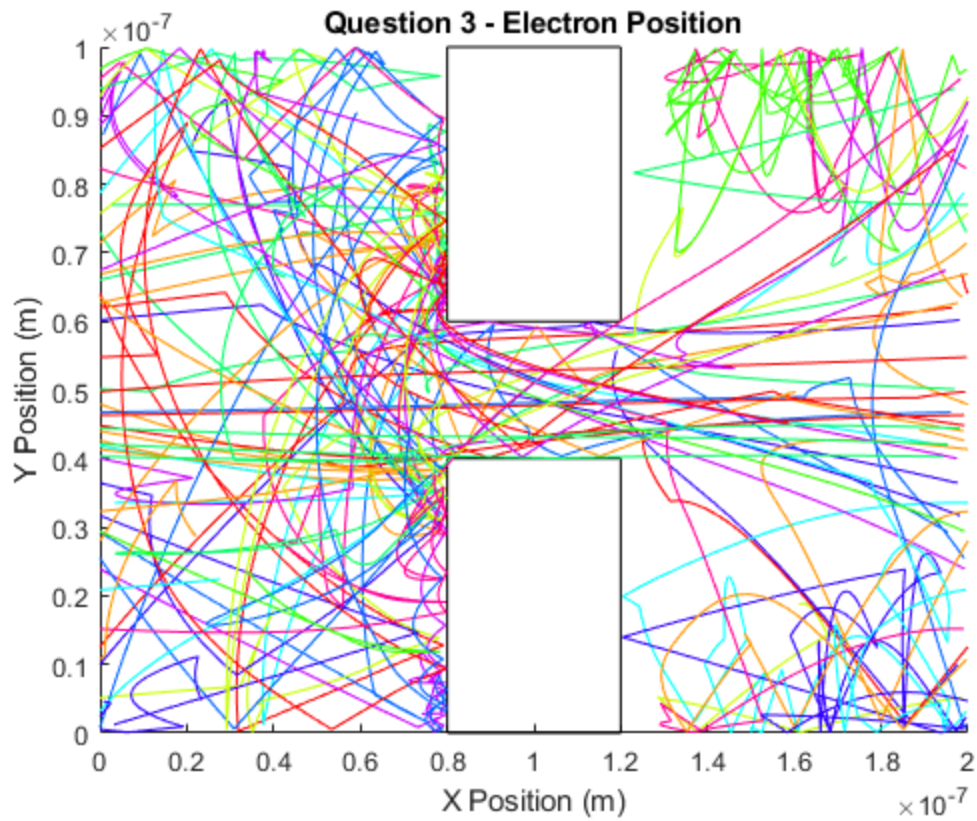
```
[xinter1, xinter2, yinter1, yinter2] =
BoxIntercept( x(1,kt), y(1,kt), x(1,kt) + x(2,kt)*Timestep,
y(1,kt) + y(2,kt)*Timestep, xboxLim(1), xboxLim(2), ylimits(1),
yboxLim(2));
if rand > DiffusionBarrierProbability
if xinter1 || xinter2
x(2,kt) = - x(2,kt);
end
if yinter1 || yinter2
y(2,kt) = - y(2,kt);
end
else
if xinter1
x(3,kt) = pi/2 + rand*pi;
elseif xinter2
x(3,kt) = pi/2 - rand*pi;
elseif yinter2
x(3,kt) = rand*pi;
elseif yinter1
x(3,kt) = -rand*pi;
end
temp = (normrnd(0, sqrt(Kbolt*T/mn))^2 + normrnd(0,
sqrt(Kbolt*T/mn))^2)^.5;
x(2,kt) = temp*cos(x(3,kt));
y(2,kt) = temp*sin(x(3,kt));
end
end
%UPDATE POSITION
x(1,kt) = x(1,kt) + x(2,kt).*Timestep;
y(1,kt) = y(1,kt) + y(2,kt).*Timestep;
%PLOT
if kt <= PlotHowMany
plot([xprev(1,kt), x(1,kt)], [yprev(1,kt), y(1,kt)], 'color',
mycolors(kt,:) );
hold on;
end
%Scattering Check
if Pscat>rand()
x(3,kt) = rand*2*pi;
temp = (normrnd(0, sqrt(Kbolt*T/mn))^2 + normrnd(0,
sqrt(Kbolt*T/mn))^2)^.5;
x(2,kt) = temp*cos(x(3,kt));
y(2,kt) = temp*sin(x(3,kt));
scatTime(1,kt) = 0;
end
xpo = round(nx*x(1,kt)/(xlimits(2)-xlimits(1)));
ypo = round(ny*y(1,kt)/(ylimits(2)-ylimits(1)));
if xpo <= 0
xpo = 1;
end
if ypo <= 0
ypo = 1;
end
if xpo >= 51
```

```

        xpo = 50;
    end
    if ypo >= 51
        ypo = 50;
    end
    x(2,kt) = x(2,kt) + Ex(ypo,xpo)*Colomb/mn*Timestep;
    y(2,kt) = y(2,kt) + Ey(ypo,xpo)*Colomb/mn*Timestep;
    end
    hold on;

end

```

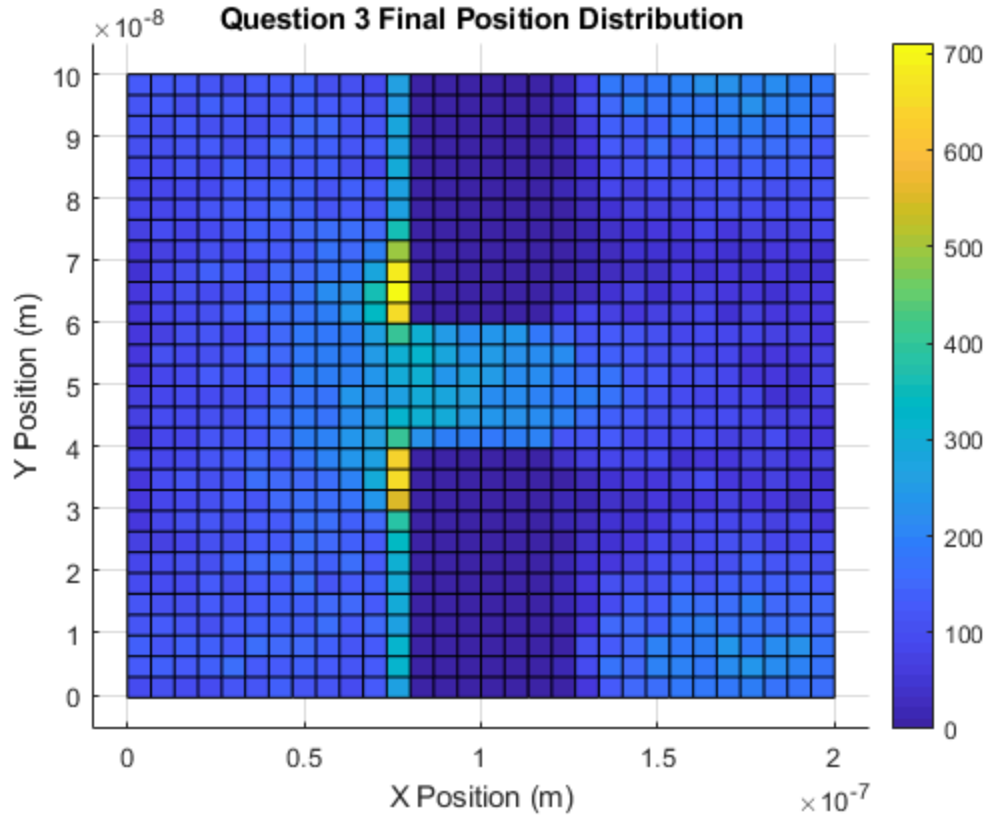


## B - Density Plot

```

figure;
Z = [transpose(x(1,:)), transpose(y(1,:))];
hist3(Z, 'Nbins',[30 30], 'CdataMode', 'auto');
hold on;
colorbar;
view(2);
grid on;
title('Question 3 Final Position Distribution');
xlabel('X Position (m)');
ylabel('Y Position (m)');
hold off;

```



The electrons have a dominate flow in one direction. This causes the electrons to be accelerated, the electrons will continue to accelerate in one direction and be stuck at the bottleneck until they pass through the small hole this causes a build up on one side of the bottleneck and also causes the passage way of the bottleneck to be more densely populated by electrons. On the other side of the bottleneck there is a lower density as the electrons are allowed to spread out and not be restricted.

end

## C - More Accurate Simulation

Allowing electrons to flow through the bottleneck would be more accurate as it has higher resistivity but doesn't stop all electrons from passing through it

The probability of diffusion is set to be a given value by the user, having a more accurate diffusion rate would also produce more accurate simulations

In materials there are always impurities, allowing there to be random impurities that cause greater scattering at that specific point, or diffusion or a higher restivity would provide a more accurate simulation.

Currently this is a two dimensional example, using a three dimensional example would be more accurate.

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