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function [] = Q2(numOfAtom, numOfStep)
% ELEC4700 - Assignment 1
% Xiaochen Xin 100989338
C.q_0 = 1.60217653e-19;           % electron charge
C.hb = 1.054571596e-34;          % Dirac constant
C.h = C.hb * 2 * pi;              % Planck constant
C.m_0 = 9.10938215e-31;           % electron mass(kg)
C.kb = 1.3806504e-23;             % Boltzmann constant
C.eps_0 = 8.854187817e-12;        % vacuum permittivity
C.mu_0 = 1.2566370614e-6;         % vacuum permeability
C.c = 299792458;                  % speed of light
C.g = 9.80665; %metres (32.1740 ft) per s^2
C.am = 1.66053892e-27;

mn = 0.26*C.m_0; %Effective Mass
l = 200e-9; %Length of area (m)
w = 100e-9; %Width of area (m)
T = 300; %Kelvin

vth = sqrt(C.kb*T/mn); %thermal velocity(velocity at which the
    particles are travelling at)

tmn = 0.2e-12; %mean time between collision (s)

%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%
%%
rvx = randn(numOfAtom,1)*sqrt(C.kb*T/mn); %random vx
rvy = randn(numOfAtom,1)*sqrt(C.kb*T/mn); %random vY

v = sqrt(rvx.^2+rvy.^2);
figure (1)
plot(hist(v,100))
ylabel("Number of Particles")
xlabel("Bins")
title ("Particle Velocity Distribution")

xr = 200e-9.*rand(numOfAtom,1); %x of 100 random locations
yr = 100e-9.*rand(numOfAtom,1); %y of 100 random locations

%Define two arrays store the previous locations
xrp = xr;
yrp = yr;

MFPx = xr;
MFPy = yr;
MFP = zeros(numOfAtom,1);
MTBC = zeros(numOfAtom,1);
scatter_number = zeros(numOfAtom,1);

t = 1.5e-14; %time interval that captures line
xd = rvx*t; %displacement in x during one time interval

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yd = rvy*t; %displacement in y during one time interval

Pscat = 1-exp(-t/tmn); % Probability that a particle scatters

for p = 1:1:numOfStep
    scatter_prob = rand(numOfAtom,1);
    %%%%%%%%%Calculate Mean Free Path%%%%%%%%%
    MFP(scatter_prob<Pscat) = MFP(scatter_prob<Pscat) +
    sqrt((xd(scatter_prob<Pscat)-
MFPx(scatter_prob<Pscat)).^2+(yd(scatter_prob<Pscat)-
MFPy(scatter_prob<Pscat)).^2);
    MTBC(scatter_prob<Pscat) = MTBC(scatter_prob<Pscat) +
    sqrt((xd(scatter_prob<Pscat)-
MFPx(scatter_prob<Pscat)).^2+(yd(scatter_prob<Pscat)-
MFPy(scatter_prob<Pscat)).^2)./v(scatter_prob<Pscat);
    scatter_number(scatter_prob<Pscat) =
    scatter_number(scatter_prob<Pscat) + 1;
    MFPx(scatter_prob<Pscat) = xr(scatter_prob<Pscat);
    MFPy(scatter_prob<Pscat) = yr(scatter_prob<Pscat);
    %%%%%%%%%Calculate average temperature of all particles%%%%%%%%%
    %%%%%%%%%
    v = sqrt(rvx.^2+rvy.^2);
    TParticles = (0.5*mn*v.^2)/(C.kb);%Tempearture of individual
particles
    Tave (p) = sum(TParticles)/numOfAtom;%Average temperature of all
particles
    figure (3)
    plot (Tave)
    xlabel("Number of steps (1.5e-14s/step)")
    ylabel("Temperature (K)")
    ylim ([0, 500])
    xlim ([0, numOfStep])
    title("average temperature over time")
    %%%%%%%%%Calculate Mean Time Between Collision%%%%%%%%%
    %%%%%%%%%

    rvx_new = randn(numOfAtom,1)*sqrt(C.kb*T/mn); %new random vx
    rvy_new = randn(numOfAtom,1)*sqrt(C.kb*T/mn); %new random vY
    rvx(scatter_prob<Pscat) = rvx_new(scatter_prob<Pscat);
    rvy(scatter_prob<Pscat) = rvy_new(scatter_prob<Pscat);

    xd = rvx*t; %displacement in x during one time interval
    yd = rvy*t; %displacement in y during one time interval

    xr = xr+xd;
    yr = yr+yd;
    %Define the left&right wrap-around
    xrp(xr >2e-7) = - (2e-7 - xrp(xr >2e-7));% changing previous point
to prevent line drawn across canvas
    xr(xr > 2e-7) = xr(xr > 2e-7)-(2e-7);
    xrp(xr <0) = 2e-7 - xrp(xr <0);% changing previous point to
prevent line drawn across canvas
    xr(xr < 0) = xr(xr < 0 )+(2e-7);

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%Define the specular top&bottom
rvy(yr > 1e-7) = -rvy(yr > 1e-7);
yr(yr > 1e-7) = (1e-7)-(yr(yr > 1e-7)-(1e-7));
rvy(yr < 0) = -rvy(yr < 0 );
yr(yr < 0) = -yr(yr < 0 );

figure (2)
plot([xrp(1), xr(1)], [yrp(1), yr(1)], 'r')
plot([xrp(2), xr(2)], [yrp(2), yr(2)], 'b')
plot([xrp(3), xr(3)], [yrp(3), yr(3)], 'k')
plot([xrp(4), xr(4)], [yrp(4), yr(4)], 'g')
plot([xrp(5), xr(5)], [yrp(5), yr(5)], 'y')
plot([xrp(6), xr(6)], [yrp(6), yr(6)], 'c')
xlabel("Semiconductor Dimension (m)")
ylabel("Semiconductor Dimension (m)")
title ("Particles Trajectory")
xlim ([0, 2e-7])
ylim([0,1e-7])
grid on
hold on
pause(0.05)

xrp = xr;
yrp = yr;
end
%Display Overall MFP
overallMFP = sum(MFP./scatter_number)/numOfAtom
overallMTBC = sum(MTBC./scatter_number)/numOfAtom
end

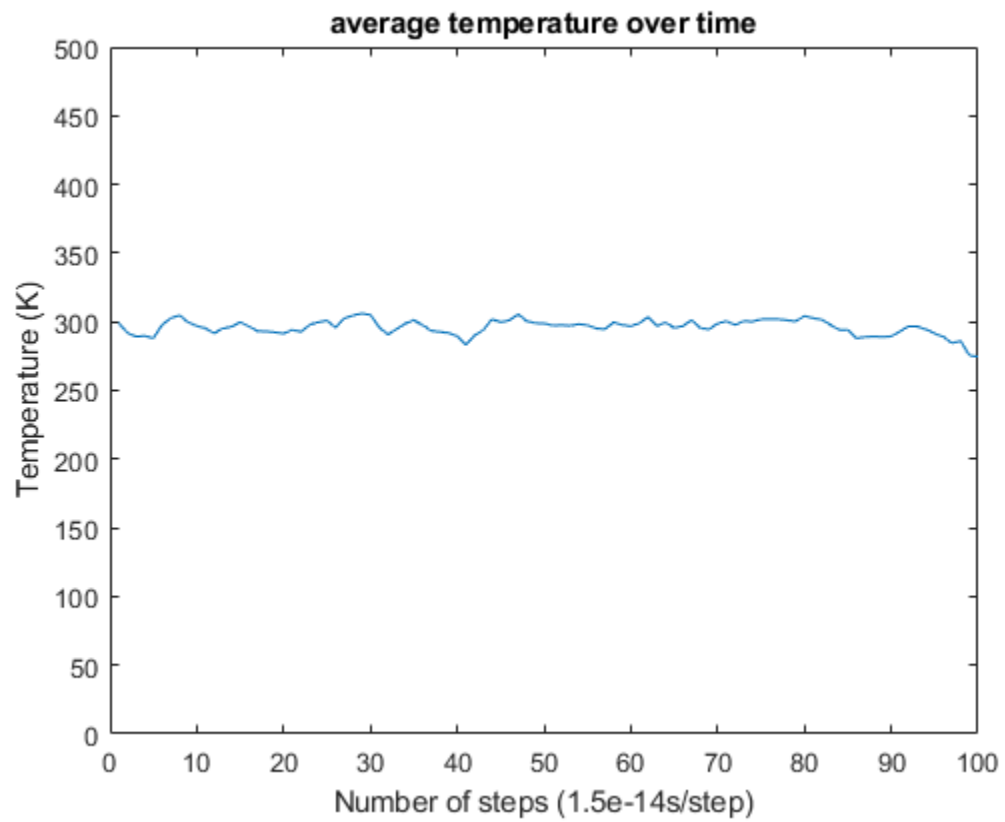
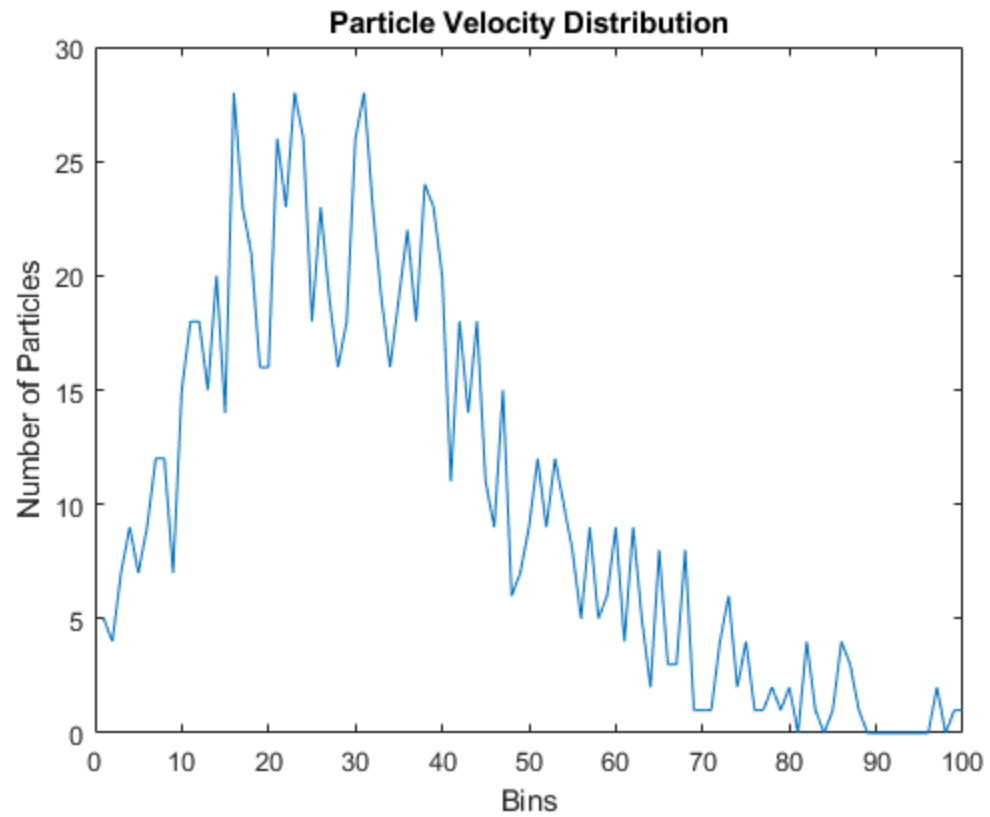
overallMFP =

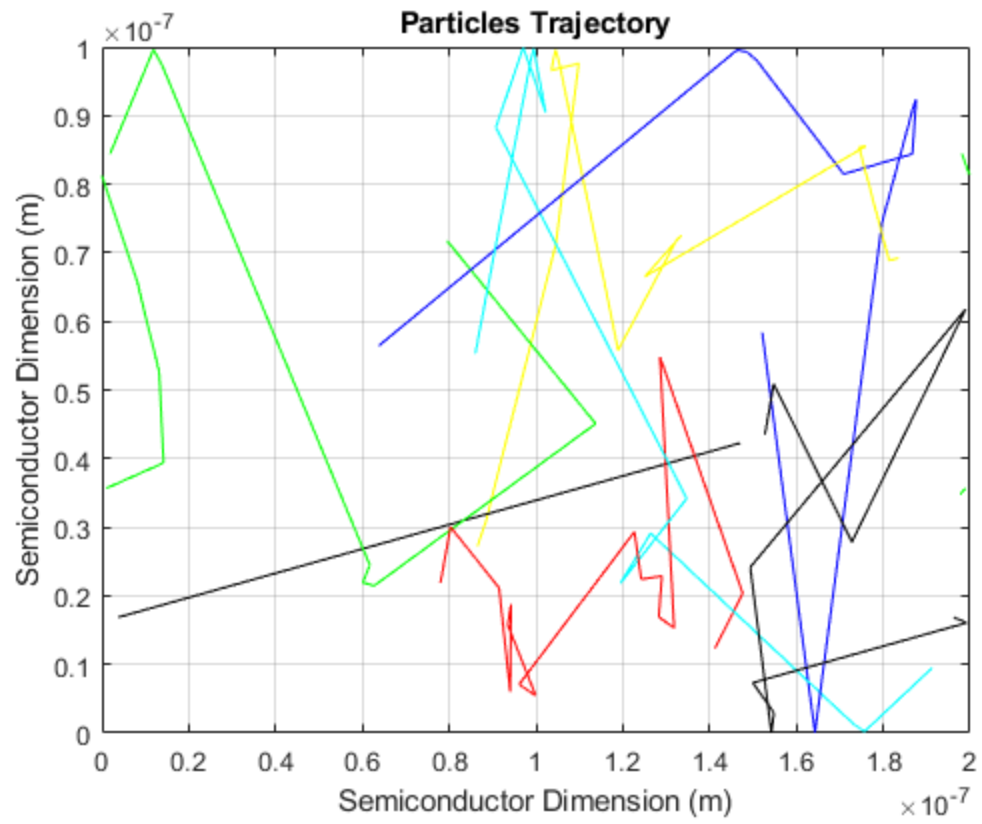
1.1934e-07

overallMTBC =

1.1399e-12

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