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```

clear; clc;

e = 1.6e-19; eVtoJ = e; JtoEv = e^(-1);
me = 9.11*1e-31; nm = 1e-9;

T = 800; % K

checkTime = [0, 1, 5, 10]; %years

% atoms' radius
dx = 0.56; %nm

% Count layers
a = 10; % monolayers
b = 6;
c = 6;

sizeHS = a + b + c + b + a;

% Fermi Energy
EFermi = 1.51*1e-20; % J

% Applyied voltage
dU = 0:0.01:0.5;

% grid of Al conentration
grid_x_Al = [zeros(1, a), ...
    0.44*ones(1, b), ...
    zeros(1, c), ...
    0.44*ones(1, b), ...
    zeros(1, a)
];

% Get profile Ec
[grids_Ec, grids_meff, grids_C_Al] = getDiffOpenAlGaAs( grid_x_Al,
    checkTime, dx*nm, T );

% get J from V
for j = 1 : length(checkTime)
    J(j, :) = getJ(dx*nm, ...
        grids_meff(j, :)*me, ...
        grids_Ec(j, :)*eVtoJ, ...
        dU*eVtoJ, ...
        EFermi...
    );
end

showResult(grids_C_Al, grids_Ec, checkTime, J, sizeHS, dx, T, dU); %
    Img 4.3, Img 4.4

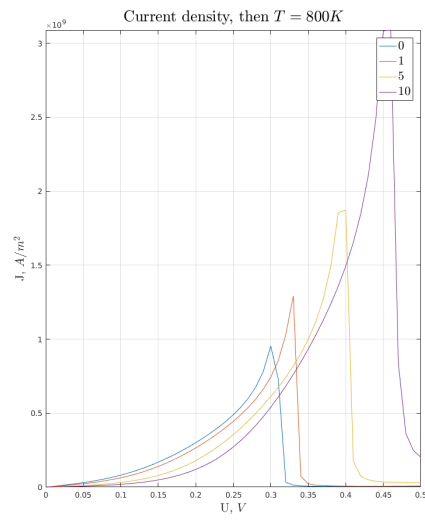
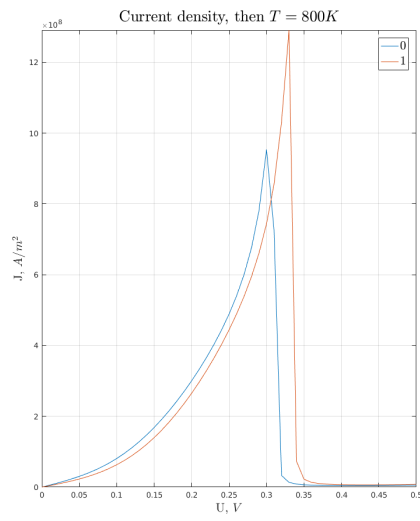
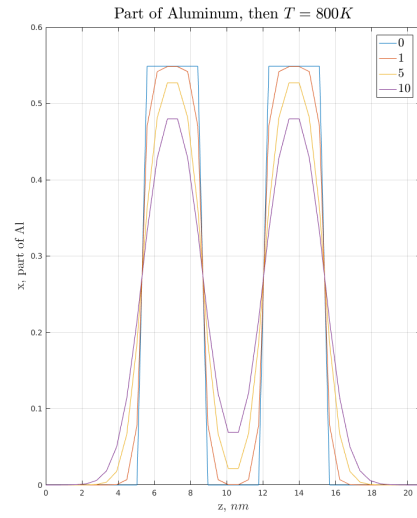
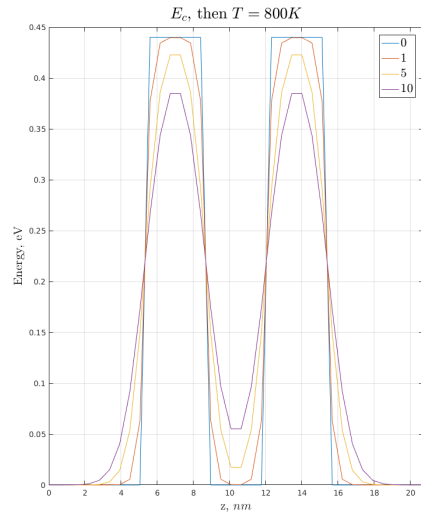
```

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`outResult =`

1



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