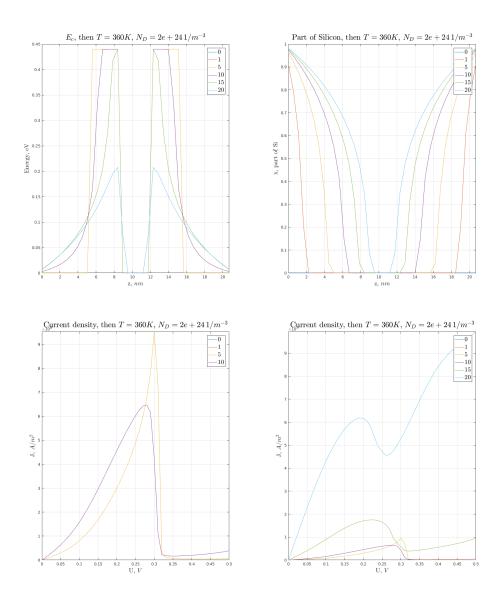
```
%Img 4.7, 4.8
clear; clc;
e = 1.6e-19; eVtoJ = e; JtoEv = e^{(-1)};
me = 9.11*1e-31; nm = 1e-9;
T = 360; % K
checkTime = [0, 1, 5, 10, 15, 20]; %years
% atoms' radius
dx = 0.56; %nm
% Count layers
a = 10; % monolayers
b = 6;
c = 6;
sizeHS = a + b + c + b + a;
% Fermi Energy
EFermi = 1.51*1e-20; % 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)
];
Nd = 2*1e18*1e6;
% Get profile Ec
[grids_Ec, grids_meff, grids_C_Al, Six] = getDiffAlGaAs_Si( grid_x_Al,
checkTime, dx*nm, T, Nd);
% 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, Six/Nd, grids_Ec, checkTime, J, sizeHS, dx, T,
dU, Nd); %Img 4.7, 4.8
```





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