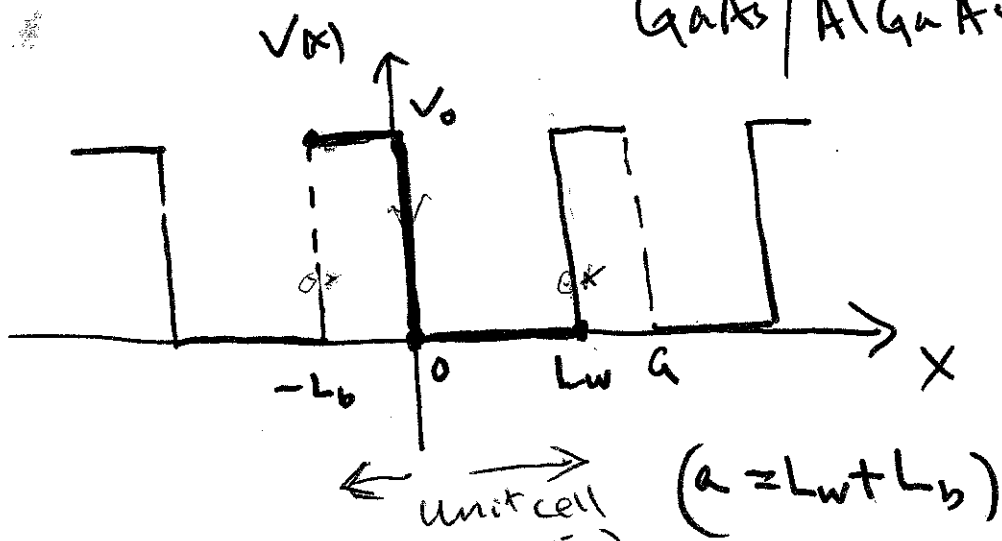


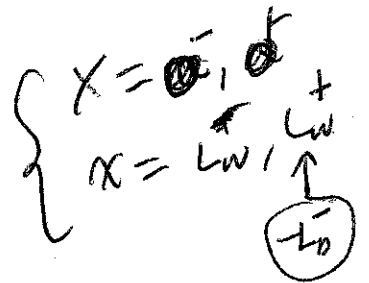
1D Superlattice (Semiconductor GaAs/AlGaAs)



for $0 < x < L_w =$

$$\psi(x) = A e^{ik_w x} + B e^{-ik_w x}$$

$$k_w = \left(\frac{2m_w E}{\hbar^2} \right)^{1/2}$$



for $-L_b < x < 0 =$

$$\psi(x) = C e^{k_b x} + D e^{-k_b x}$$

$$k_b = \left[\frac{2m_b (V_0 - E)}{\hbar^2} \right]^{1/2}$$

BCs:

$$A + B = C + D \quad \text{--- (1)}$$

$$ik_w (A - B) = k_b (C - D) \quad \text{--- (2)}$$

$$\psi(L_w) = \psi(L_w) = \psi(-L_b) e^{ika} = \left[C e^{-k_b L_b} + D e^{k_b L_b} \right] e^{ika} \quad \text{--- (3)}$$

$$\psi'(L_w) = \psi'(L_w) = \psi'(-L_b) e^{ika} = ik_w (A e^{ik_w L_w} - B e^{-ik_w L_w}) = k_b (C e^{-k_b L_b} - D e^{k_b L_b}) e^{ika} \quad \text{--- (4)}$$

Only wavefunctions used in $-L_b \leq x \leq L_w$ within the unit cell

```

%while k<10
%k=k+1;
k=5
x=0;
while x<300
x=x+1;
y11=x-k^2;
y12=0;
y21=0;
y22=x-3*k^2;
y=[y11 y12;y12 y22];
%z=abs(det(y));
%if z<40
%a=x
%plot(k,a,'ro')
%hold on
%end
semilogy(x,abs(det(y)),'r+')
hold on
end
end
end

```

$$y_{21} = y_{12} = 0$$

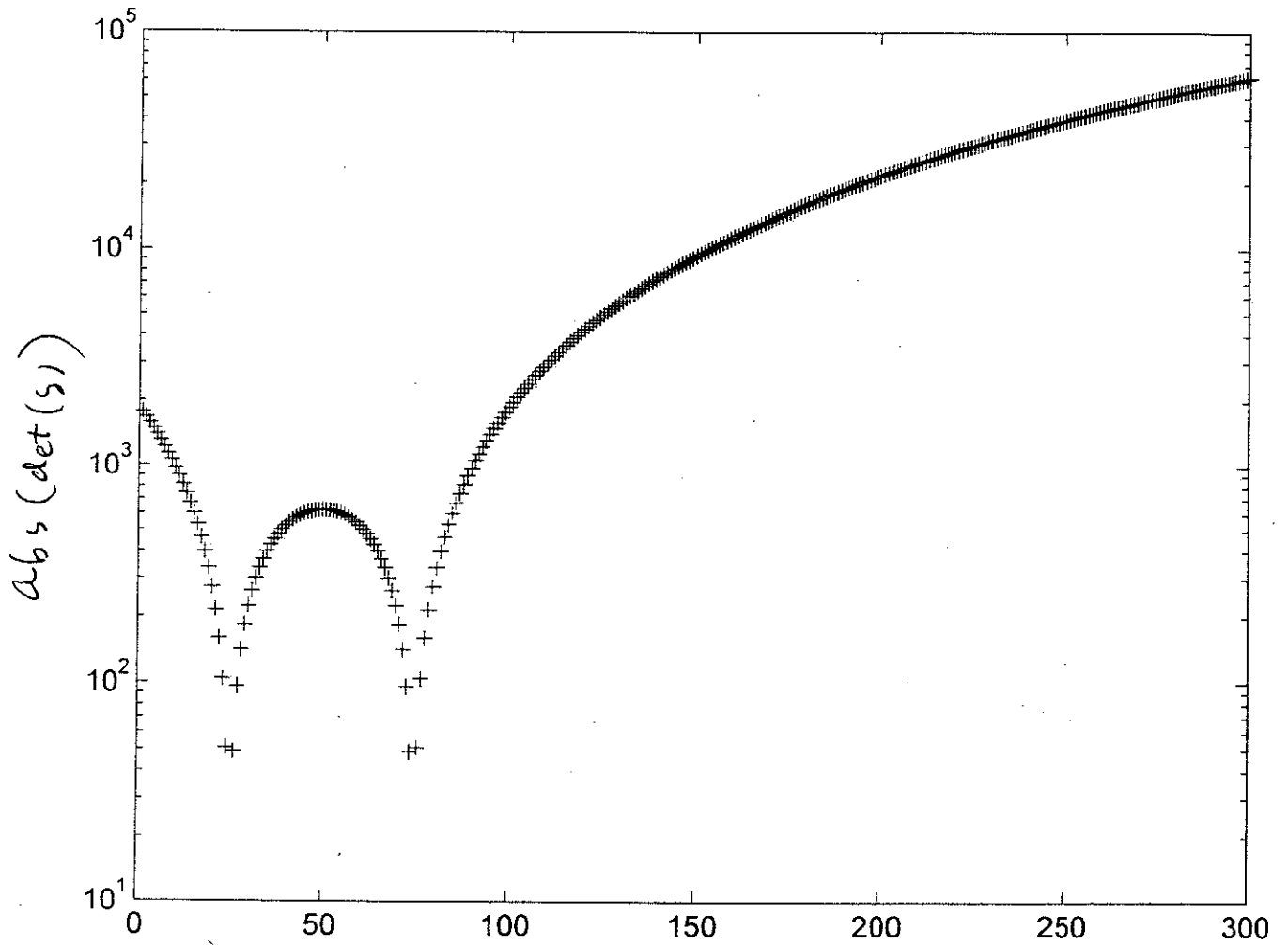
test function 3/25/00

$$f = \begin{vmatrix} (x - k^2) & 0 \\ 0 & (x - 3k^2) \end{vmatrix}$$

$$= (x - k^2)(x - 3k^2) = 0$$

→ Fig. 1

$$\begin{cases}
 \Rightarrow A + B - C - D = 0 & \text{--- (1)} \\
 kA - kB - kC + kD = 0 & \text{--- (2)} \\
 A \quad B \quad C \quad D = 0 & \text{--- (3)} \\
 A \quad B \quad C \quad D = 0 & \text{--- (4)}
 \end{cases}$$



x

Fig. 1

```
k=0;
while k<10
k=k+1;
x=0;
while x<300
x=x+1;
y11=x-k^2;
y12=0;
y21=0;
y22=x-3*k^2;
y=[y11 y12;y12 y22];
z=abs(det(y));
if z<40
a=x
plot(k,a,'ro')
hold on
end
%semilogy(x,abs(det(y)))
%hold on
end
```

→ Fig. 2

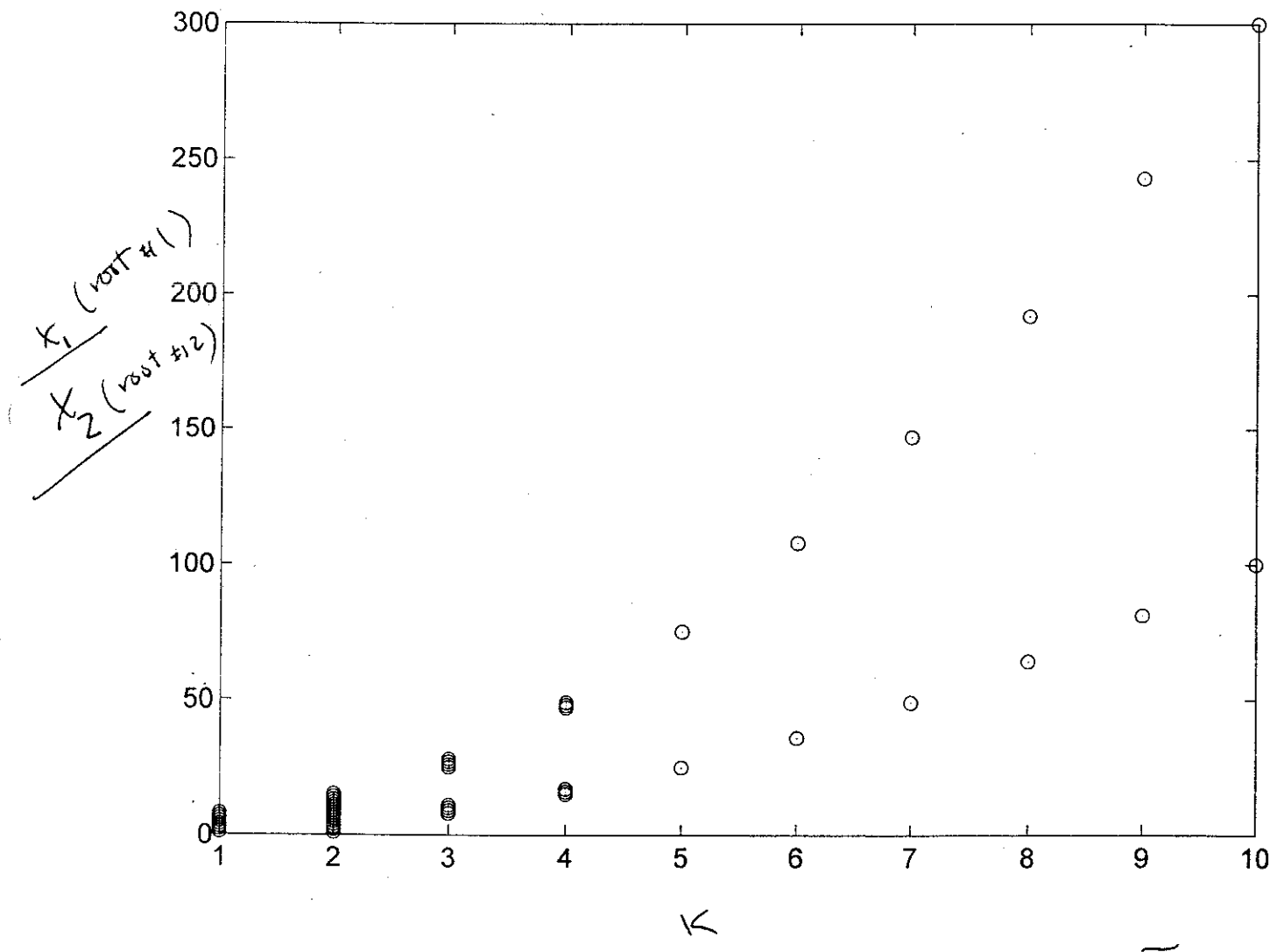


Fig. 2

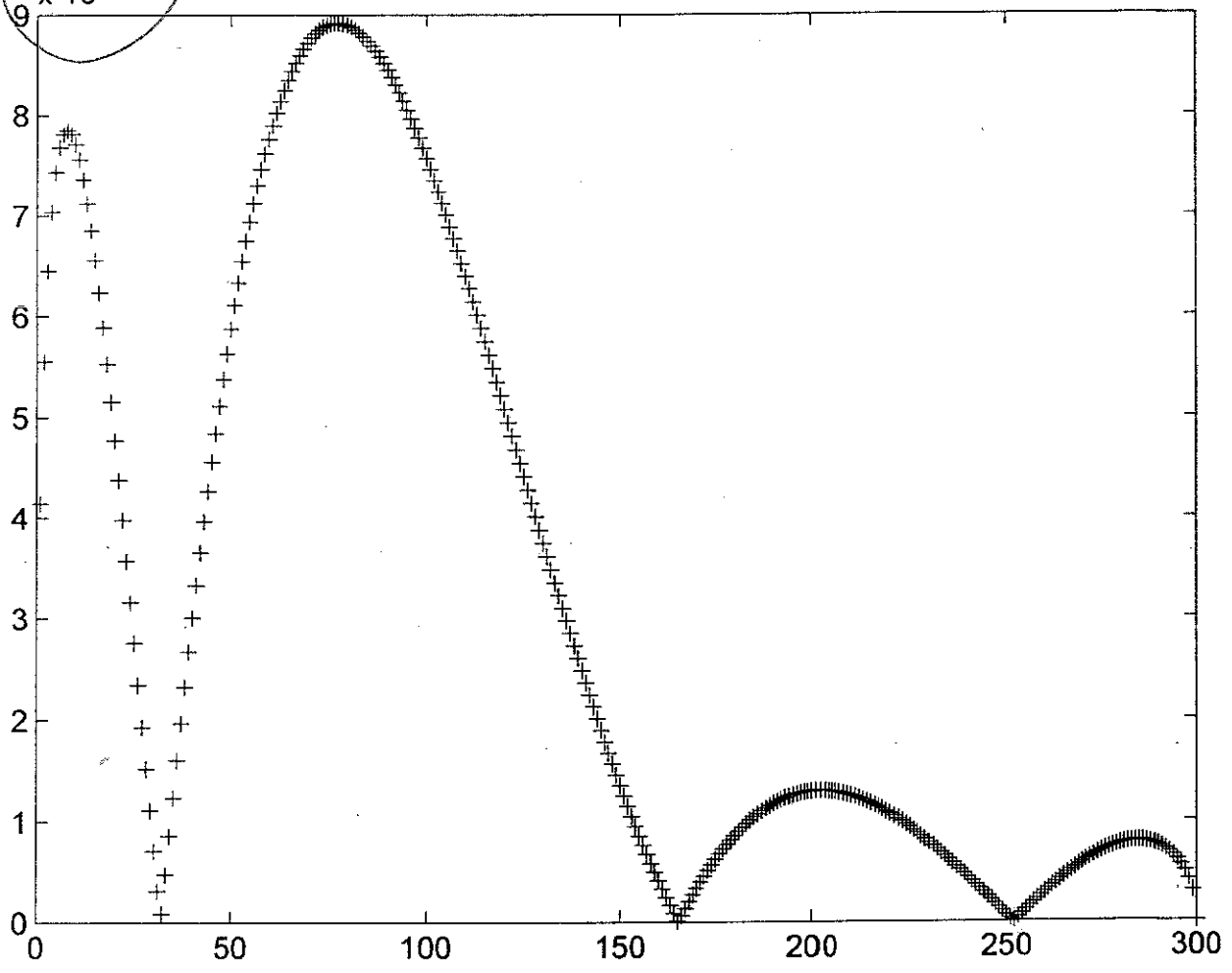
```

% this m.file is to calculate the band structure of *
% 1D superlattice
% The following is a set of constants used
%
i=sqrt(-1);          % define imaginary unit
m0=9.1e-31;         % free electron mass, kg
hbar=1.05459e-34;   % Planck constant/2pi in unit of J.s
mw=0.067*m0;        % electron effective mass in well
mb=0.1*m0;          % electron effective mass in barrier
mev=1.6e-22;        % energy conversion factor
v0=300*mev;         % barrier height
angs=1e-10;         % length conversion factor
lw=100*angs;        % well width
lb=20*angs;         % barrier width
a=lb+lw;            % period, be considered as spacing bet ions
%
% 1st. for a given k value
k=pi/a/50;
% for given k, you can plot the abs(det(z)) vs e and find out
% where are the roots, where z is a matrix by 4 boundary conditions
e=0;
while e<300
e=e+1;
kw=sqrt(2*mw*e*mev/hbar/hbar);
kb=sqrt(2*mb*(v0-e*mev)/hbar/hbar);
z11=1;
z12=1;
z13=-1;
z14=-1;
z21=i*kw;
z22=-i*kw;
z23=-kb;
z24=kb;
z31=exp(i*kw*lw);
z32=exp(-i*kw*lw);
z33=-exp(-kb*lb+i*k*a);
z34=-exp(kb*lb+i*k*a);
z41=i*kw*exp(i*kw*lw);
z42=-i*kw*exp(-i*kw*lw);
z43=-kb*exp(-kb*lb+i*k*a);
z44=kb*exp(kb*lb+i*k*a);
x=[z11 z12 z13 z14;z21 z22 z23 z24;z31 z32 z33 z34;z41 z42 z43 z44];
dx=abs(det(x));
plot(e,dx,'r') → f.g.3
hold on
end

```

965
(electrons)

9×10^{18}



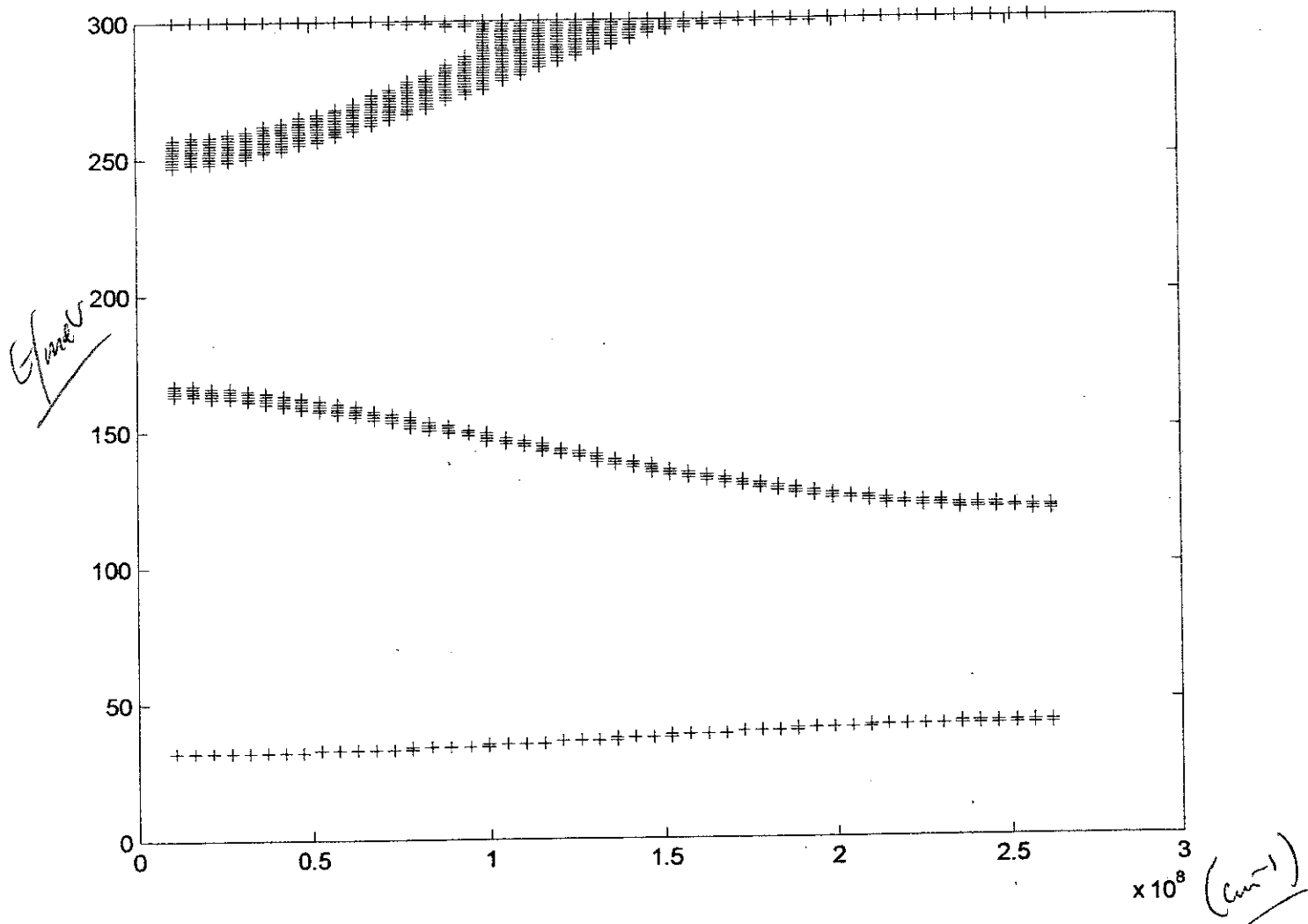
E (mev)

Fig. 3

```

% this m.file is to calculate the band structure of
% 1D superlattice
% The following is a set of constants used
%
i=sqrt(-1);      % define imaginary unit
m0=9.1e-31;      % free electron mass, kg
hbar=1.05459e-34; % Planck constant/2pi in unit of J.s
mw=0.067*m0;    % electron effective mass in well
mb=0.1*m0;      % electron effective mass in barrier
mev=1.6e-22;    % energy conversion factor
v0=300*mev;     % barrier height
angs=1e-10;     % length conversion factor
lw=100*angs;    % well width
lb=20*angs;     % barrier width
a=lb+lw;        % period, be considered as spacing bet ions
%
% 1st, for a given k value
k=pi/a/50;
while k<pi/a
    k=k+pi/a/50;
% for given k, you can plot the abs(det(z)) vs e and find out
% where are the roots, where z is a matrix by 4 boundary conditions
e=0;
while e<300
e=e+1;
kw=sqrt(2*mw*e*mev/hbar/hbar);
kb=sqrt(2*mb*(v0-e*mev)/hbar/hbar);
z11=1;
z12=1;
z13=-1;
z14=-1;
z21=i*kw;
z22=-i*kw;
z23=-kb;
z24=kb;
z31=exp(i*kw*lw);
z32=exp(-i*kw*lw);
z33=-exp(-kb*lb+i*k*a);
z34=-exp(kb*lb+i*k*a);
z41=i*kw*exp(i*kw*lw);
z42=-i*kw*exp(-i*kw*lw);
z43=-kb*exp(-kb*lb+i*k*a);
z44=kb*exp(kb*lb+i*k*a);
x=[z11 z12 z13 z14;z21 z22 z23 z24;z31 z32 z33 z34;z41 z42 z43 z44];
dx=abs(det(x));
if dx<2e17 ← note the big a for dx
    ee=e;
plot(k, ee, 'r+') → Fig. 4
hold on
end
end
end
end

```

$$\begin{aligned}
 \frac{h}{m} &= \frac{h}{120A} \\
 &= \frac{h}{120 \times 10^8 \text{ m}} = \frac{h}{120} \times 10^{-10} \text{ m}^{-1} \\
 &= \frac{h}{120} \text{ cm}^{-1}
 \end{aligned}$$

Fig. 4 (cm^{-1})