

Scilab Textbook Companion for  
Semiconductor Device Physics And Design  
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# Book Description

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Scilab numbering policy used in this document and the relation to the above book.

**Exa** Example (Solved example)

**Eqn** Equation (Particular equation of the above book)

**AP** Appendix to Example(Scilab Code that is an Appednix to a particular Example of the above book)

For example, Exa 3.51 means solved example 3.51 of this book. Sec 2.3 means a scilab code whose theory is explained in Section 2.3 of the book.

# Contents

List of Scilab Codes	4
1 STRUCTURAL PROPERTIES OF SEMICONDUCTORS	5
2 ELECTRONIC LEVELS IN SEMICONDUCTORS	8
3 CHARGE TRANSPORT IN MATERIALS	11
4 JUNCTIONS IN SEMICONDUCTORS PN DIODES	18
5 SEMICONDUCTOR JUNCTIONS	24
6 BIPOLAR JUNCTION TRANSISTORS	26
7 TEMPORAL RESPONSE OF DIODES AND BIPOLAR TRANSISTORS	31
8 FIELD EFFECT TRANSISTORS	33
9 FIELD EFFECT TRANSISTORS MOSFET	35

# List of Scilab Codes

Exa 1.1	Crystal structure . . . . .	5
Exa 1.2	Surface density . . . . .	6
Exa 1.2.5.1	Planar density of atoms on a surface . . . . .	7
Exa 2.1	density of states in 2D and 3D systems . . . . .	8
Exa 2.2	effective density of states for the conduction and valence bands . . . . .	8
Exa 2.5	Piezoelectric Effect . . . . .	9
Exa 2.7.1	Fermi energy and Fermi velocity . . . . .	10
Exa 3.1	polar optical phonon emission rate . . . . .	11
Exa 3.2	transport under an electric field . . . . .	11
Exa 3.3	relaxation time due to ionized impurity scattering . . . . .	12
Exa 3.4	time between scattering events using the conductivity effective mass . . . . .	12
Exa 3.5	conductivity of doped versus undoped . . . . .	13
Exa 3.6	maximum and minimum conductivity . . . . .	14
Exa 3.7	High field transport velocity field relations . . . . .	14
Exa 3.8	transit time of electron in semiconductor device . . . . .	15
Exa 3.9	band to band tunneling probability . . . . .	16
Exa 3.10	diffusion coefficient using velocity field relation . . . . .	16
Exa 3.11	carrier generation rate for optical radiation . . . . .	17
Exa 4.1	pn junction in equilibrium . . . . .	18
Exa 4.2	effect of the generation recombination current in pn diode . . . . .	19
Exa 4.3	diode current at a forward bias . . . . .	19
Exa 4.4	breakdown voltage of diode . . . . .	20
Exa 4.5	e h recombination time . . . . .	21
Exa 4.6	internal radiative efficiency for diodes . . . . .	21
Exa 4.7	injection efficiency of LED . . . . .	22
Exa 4.8	photon flux and optical power generated by LED . . . . .	23

Exa 5.2	SCHOTTKY BARRIER . . . . .	24
Exa 6.1	emitter efficiency of BJT and HBT . . . . .	26
Exa 6.2	change in the base width and collector current with voltage and Early voltage . . . . .	27
Exa 6.3	maximum base width for a given beta . . . . .	28
Exa 6.4	output conductance and emitter efficiency and gain . . . . .	29
Exa 7.1	cutoff frequency of transistor . . . . .	31
Exa 8.1	gate current density for MESFET . . . . .	33
Exa 8.5.1	2DEG concentration for MODFET . . . . .	34
Exa 9.1	maximum depletion width in a MOS capacitor . . . . .	35
Exa 9.2	MOS capacitor . . . . .	35
Exa 9.3	MOS capacitor threshold voltage . . . . .	36
Exa 9.4	MOS capacitor threshold voltage and channel conductivity . . . . .	37
Exa 9.6	oxide capacitance and capacitance at the flat band and the minimum capacitance at threshold for MOS capacitor . . . . .	38
Exa 9.7	saturation current of MOSFET . . . . .	38
Exa 9.8	drain current for NMOS device . . . . .	39
Exa 9.9	threshold voltage in n channel MOSFET . . . . .	40
Exa 9.10	shift in the threshold voltage arising from source body bias . . . . .	40
Exa 9.11	n channel MOSFET characteristics . . . . .	41

# Chapter 1

## STRUCTURAL PROPERTIES OF SEMICONDUCTORS

Scilab code Exa 1.1 Crystal structure

```
1 a = 5.43*10^-8; //lattice constant for silicon in cm
2 N = 8/8 +6/2; // Silicon has a diamond structure
  which is made up of the fcc lattice with two
  atoms on each lattice point. The fcc unit cube
  has a volume a3. The cube has eight lattice sites
  at the cube edges. However, each of these points
  is shared with eight other cubes. In addition,
  there are six lattice points on the cube face
  centers. Each of these points is shared by two
  adjacent cubes.
3 disp("Silicon has a diamond structure which is made
  up of the fcc lattice with two atoms on each
  lattice point. The fcc unit cube has a volume a3.
  The cube has eight lattice sites at the cube
  edges. However, each of these points is shared
  with eight other cubes. In addition, there are
  six lattice points on the cube face centers. Each
  of these points is shared by two adjacent cubes.
  ")
```

```

4 disp(N, "Thus, number of lattice points per cube of
   volume a^3 = ")
5 disp("In silicon there are two silicon atoms per
   lattice point. The number density is, therefore,")
   )
6 Nsi = N*2/a^3;
7 disp(Nsi,"in atoms per cm cube")
8 l = 50*10^-4;
9 b = 2*10^-4;
10 h = 1*10^-4;
11 vol = l*b*h; //volume of the MOSFET
12 disp(vol, "volume of th MOSFET (in cm cube) = ")
13 nmos = Nsi*vol;
14 disp(nmos,"Number of Si atoms in the MOSFET = ")

```

---

### Scilab code Exa 1.2 Surface density

```

1 a = 5.65*10^-8; //lattice constant in cm
2 disp("In the (001) surfaces, the top atoms are
   either Ga or As leading to the terminology Ga
   terminated (or Ga stabilized) and As terminated (
   or As stabilized), respectively. A square of area
   a2 has four atoms on the edges of the square and
   one atom at the center of the square. The atoms
   on the square edges are shared by a total of four
   squares. The total number of atoms per square is
   ")
3 N = 4/4 +1;
4 disp(N)
5 density = N/a^2; //Surface density
6 disp(density, "The surface density (in per cm square
   ) of Ga atoms on a Ga terminated (001) GaAs
   surface")

```

---

**Scilab code Exa 1.2.5.1** Planar density of atoms on a surface

```
1 a = 5.63*10^-8; //lattice constant in cm
2 disp("We can form a triangle on the (111) surface of
      GaAs. There are three atoms on the tips of the
      triangle. These atoms are shared by six other
      similar triangles. There are also 3 atoms along
      the edges of the triangle which are shared by two
      adjacent triangles. Thus the number of atoms in
      the triangle are")
3 N = 3/6 +3/2;
4 disp(N)
5 area = 3^0.5/2*a^2; //area of triangle
6 density = N/area; //The density of GaAs atoms on the
      surface
7 disp(density,"The density (in per square cm) of GaAs
      atoms on the surface = ")
```

---

## Chapter 2

# ELECTRONIC LEVELS IN SEMICONDUCTORS

Scilab code Exa 2.1 density of states in 2D and 3D systems

```
1 m0 = 0.91 * 10^-30; //in kg
2 h = 1.05*10^-34; //in J.s
3 E = 1; // energy in eV
4 q = 1.6*10^-19;
5 N3 = 2^0.5*(m0*q)^1.5*E^0.5/((%pi)^2*h^3);
6 disp(N3*10^-6,"The density of states for a 3D system
   in per eV per cm cube")
7 N2 = m0*q/(%pi*h^2);
8 disp(N2*10^-4,"The density of states for a 2D system
   in per eV per cm square")
```

---

Scilab code Exa 2.2 effective density of states for the conduction and valence bands

```
1 kBT = 26*10^-3; //in eV
2 m0 = 0.91 * 10^-30; //in kg
```

```

3 me = 0.067*m0;
4 q = 1.6*10^-19;
5 h = 1.05457*10^-34; //in J.s
6 N = 2*(me*q*kBT/2/%pi/h^2)^1.5;
7 disp(N*10^-6, "The effective density of states (in
   per cm cube) is")
8 disp("In silicon , the density of states mass is to
   be used in the effective density of states , which
   is")
9 kdos = 6^(2/3)*(0.98*0.19*0.19)^(1/3);
10 disp(kdos,"m0 times");
11 mdos = kdos*m0;
12 disp(mdos,"In silicon , the density of states mass is
   to be used in the effective density of states ,
   which is")

```

---

### Scilab code Exa 2.5 Piezoelectric Effect

```

1 //a = 0.3*aAlN + 0.7*aGaN = 3.111
2 a = 3.111*10^-8; //in cm
3 q = 1.6*10^-19;
4 Exx = 0.006; //strain tensor
5 disp(Exx,"Exx = ")
6 Ezz = -0.6*0.006;
7 disp(Ezz,"Ezz = ")
8 Ppz = 0.0097; //in C/m2
9 disp("The pizoelectric effect induced polar charge
   then becomes Ppz = 0.0097 C/m2")
10 disp(Ppz/q*10^-4,"The pizoelectric effect induced
   polar charge (in electron charge per cm2)=")
11 Psp= 0.3*0.089 + (0.7-1)*0.029;
12 disp(Psp,"The piezoelectric charge the spontaneous
   polarization charge(in C/m2) is")
13 disp(Psp/q*10^-4,"The piezoelectric charge the
   spontaneous polarization charge(in electron

```

charge per cm2) =”)

---

**Scilab code Exa 2.7.1** Fermi energy and Fermi velocity

```
1 n = 10^22; // electrons per cm cube
2 h = 1.05457*10^-34; //in J.s
3 m0 = 0.91 * 10^-30; //in kg
4 q = 1.6*10^-19;
5 E= 0.5*h^2*(3*pi*pi*n)^(2/3)/(m0*q);
6 disp(E*10^4, "The Fermi energy (in eV) is the highest
   occupied energy state at 0 K is")
7 vel = h*(3*pi*pi*n*10^6)^(1/3)/m0;
8 disp(vel, "The fermi velocity (in m/s) is")
```

---

## Chapter 3

# CHARGE TRANSPORT IN MATERIALS

Scilab code Exa 3.1 polar optical phonon emission rate

```
1 opeGaAs = 36; //optical phonon energies in GaAs in
   meV
2 opeGaN = 90; //optical phonon energies in GaN in meV
3 disp(opeGaAs,"The optical phonon energies in GaAs (
   in meV)")
4 disp(opeGaN,"The optical phonon energies in GaN (in
   meV)")
5 disp ("If the electron energies are below these
   values , there is no phonon emission.The phonon
   occupation number in GaAs at 300 K is 0.33 and in
   GaN is 0.032. Thus above threshold , the emission
   to absorption ratios are approximately 4:1 and
   32:1 respectively.")
```

---

Scilab code Exa 3.2 transport under an electric field

```

1 m0 = 0.91 * 10^-30; //in kg
2 m = 0.26*m0; //effective mass
3 E = 50*10^-3; //optical phonon energy in eV
4 t = 10^-13; //carrier scattering relaxation time at
    300K
5 q = 1.6*10^-19;
6 kBT = 26*10^-3; //in eV
7 vd = (2*q*(E-1.5*kBT)/m)^0.5;
8 disp(vd, "Drift velocity (in m/s) = ")
9 ef = vd*m/t/q; //electric field in V/cm
10 disp(ef, "Electric field(in V/m) =")

```

---

**Scilab code Exa 3.3** relaxation time due to ionized impurity scattering

```

1 u = 8500*10^-4; //in m2/V.s
2 Nd = 10^17;
3 new_u = 5000*10^-4;
4 m0 = 0.91 * 10^-30; //in kg
5 m = 0.067*m0;
6 q = 1.6*10^-19;
7 t1 = m*u/q;
8 disp(t1, "relaxation time(in s) = ")
9 t2 = m*new_u/q;
10 disp(t2, "If the ionized impurities are present, the
    time (in s) =")
11 t_imp = t2*t1/(t1-t2);
12 disp (t_imp, "The impurity-related time (in s) = ")

```

---

**Scilab code Exa 3.4** time between scattering events using the conductivity effective mass

```

1 u = 1500*10^-4; // in m2/V.s
2 m0 = 0.91 * 10^-30; //in kg

```

```

3 q = 1.6*10^-19;
4 kt = 0.19;
5 kl = 0.98;
6 ks = 3*kt*kl/(2*kl+kt);
7 t = u*ks*m0/q;
8 disp(t,"The scattering time(in s) =")

```

---

### Scilab code Exa 3.5 conductivity of doped versus undoped

```

1 Nd = 10^17;
2 Ni = 1.5*10^10;
3 Ni2 = 1.84 * 10^6;
4 Pi2 = 1.84 * 10^6;
5 Pi = 1.5*10^10;
6 un1 = 1000;
7 up1 = 350;
8 un2 = 8000;
9 up2 = 400;
10 Nn = 0.5*Nd;
11 Pn = Ni^2/Nn;
12 q = 1.6*10^-19;
13 disp(Pn,"Hole density for Si(in per cm cube) =")
14 s_n1 = Nn*q*un1+Pn*q*up1;
15 s_un1 = Ni*q*un1+Pi*q*up1;
16 s_n2 = Nn*q*un2+Pn*q*up2;
17 s_un2 = Ni2*q*un2+Pi2*q*up2;
18 disp(s_n1,"The conductivity of Si (in per ohm per cm
) = ")
19 disp(s_un1,"The conductivity of undoped Si (in per
ohm per cm) = ")
20 disp(s_n2,"The conductivity of GaAs (in per ohm per
cm) = ")
21 disp(s_un2,"The conductivity of undoped GaAs (in per
ohm per cm) = ")

```

---

**Scilab code Exa 3.6** maximum and minimum conductivity

```
1 d1 = 2.78 * 10^19; //max density for Si
2 d2 = 7.72 * 10^18; // max density for GaAs
3 Nd = 10^17;
4 Ni = 1.5*10^10;
5 Ni2 = 1.84 * 10^6;
6 Pi2 = 1.84 * 10^6;
7 Pi = 1.5*10^10;
8 un1 = 1000;
9 up1 = 350;
10 un2 = 8000;
11 up2 = 400;
12 Nn = 0.5*Nd;
13 Pn = Ni^2/Nn;
14 q = 1.6*10^-19;
15 s1 = d1*q*un1;
16 s2 = d2*q*up2;
17 //To find the minimum we take the derivative with
    respect to p and equate the result to zero, which
    gives the below expression
18 p = Ni*(un1/up1)^0.5;
19 smin1 = Ni*q*(un1*(up1/un1)^0.5 + up1*(un1/up1)^0.5)
    ;
20 smin2 = Ni2*q*(un2*(up2/un2)^0.5 + up2*(un2/up2)
    ^0.5);
21 disp(s1,"maximum conductivity for Si = ")
22 disp(s2,"maximum conductivity for GaAs = ")
23 disp(smin1,"minimum conductivity for Si = ")
24 disp(smin2,"minimum conductivity for GaAs = ")
```

---

**Scilab code Exa 3.7** High field transport velocity field relations

```

1 E1 = 10^3; //in V/cm
2 E2 = 10^5; //in V/cm
3 v1 = 1.4*10^6; //in cm/s
4 v2 = 1*10^7; //in cm/s
5 m0 = 0.91 * 10^-30; //in kg
6 m = 0.26*m0;
7 q = 1.6*10^-19;
8 u1 = v1/E1;
9 u2 = v2/E2;
10 disp(u1,"mobility (in cm square per V.s) at 1 kV/cm
    = ")
11 disp(u2,"mobility (in cm square per V.s) at 100 kV/
    cm = ")
12 t1 = m*u1*10^-4/q;
13 t2 = m*u2*10^-4/q;
14 disp(t1,"relaxation time (in s) at 1 kV/cm = ")
15 disp(t2,"relaxation time (in s) at 100 kV/cm = ")

```

---

**Scilab code Exa 3.8** transit time of electron in semiconductor device

```

1 E = 50*10^3; //in V/cm
2 v = 10^7; // in cm/s
3 L = 10^-5; //in cm
4 m0 = 0.91 * 10^-30; //in kg
5 m = 0.067*m0;
6 q = 1.6*10^-19;
7 a = q*E/m;
8 t = (2*L/a)^0.5;
9 disp(t,"If the transport is ballistic , transit time
    (in s) = ")
10 t2 = L/v;
11 disp(t2,"If the saturation velocity is used, transit
    time (in s) = ")

```

---

**Scilab code Exa 3.9** band to band tunneling probability

```
1 E = 2*10^7; //in V/m
2 m0 = 0.91 * 10^-30; //in kg
3 q = 1.6*10^-19;
4 h = 1.05*10^-34; //in J.s
5 m1 = 0.065*m0; //for GaAs
6 m2 = 0.02*m0; // for InAs
7 E1 = 1.5; //in eV
8 E2 = 0.4; //in eV
9 p1 = -4*(2*m1)^0.5*(E1*q)^1.5/(3*q*h*E);
10 disp(p1,"Tunneling probability is exponent to the
    power")
11 tp1 = %e^p1;
12 disp(tp1,"Tunneling probability = ")
13 p2 = -4*(2*m2)^0.5*(E2*q)^1.5/(3*q*h*E);
14 disp(p2,"Tunneling probability is exponent to the
    power")
15 tp2 = %e^p2;
16 disp(tp2,"Tunneling probability = ")
17 disp("In InAs the band-to-band tunneling will start
    becoming very important if the field is      2
    105 V/cm.")
```

---

**Scilab code Exa 3.10** diffusion coefficient using velocity field relation

```
1 E1 = 10^5; //in V/m
2 E2 = 10^6; //in V/m
3 v1 = 1.4*10^4; //in m/s
4 v2 = 7*10^4; //in m/s
5 kBT = 26*10^-3; //in eV
6 q = 1.6*10^-19;
```

```
7 D1 = v1*kBT/E1;
8 D2 = v2*kBT/E2;
9 disp(D1,"diffusion constant (in m square per.s) at 1
    kV/cm = ")
10 disp(D2,"diffusion constant (in m square per.s) at
    10 kV/cm = ")
```

---

**Scilab code Exa 3.11** carrier generation rate for optical radiation

```
1 P = 10^3; //in W per cm square
2 E = 1.5; //in eV
3 ab = 3*10^3; //in per cm
4 t = 10^-9; //in s
5 q = 1.6*10^-19;
6 G = ab*P/E/q;
7 disp(G,"The carrier generation rate (in per cm cube
    per sec) at the surface of the sample = ")
8 d = G*t;
9 disp(d,"The excess carrier density (in per cm cube)
    = ")
```

---

## Chapter 4

# JUNCTIONS IN SEMICONDUCTORS PN DIODES

Scilab code Exa 4.1 pn junction in equilibrium

```
1 Nd = 10^16; //in per cm cube
2 p = 10^18; //in per cm cube
3 Na = 10^18; //in per cm cube
4 Nc = 2.8 * 10^19; //in per cm cube
5 Nv = 10^19; //in per cm cube
6 kT = 26*10^-3; //in eV
7 eps0 = 8.84*10^-12; //in F/m
8 eps = 11.9*eps0;
9 Eg = 1.1; //in eV
10 q = 1.6*10^-19;
11 En = kT*log(Nd/Nc);
12 disp(En,"The Fermi level positions in the n-region
        relative to the conduction band (in eV) = ")
13 Ep = -kT*log(p/Nv);
14 disp(Ep,"The Fermi level positions in the p-region
        relative to the valence band (in eV) = ")
15 Vbi = Eg + En - Ep;
```

```

16 disp(Vbi,"built-in potential = ")
17 Wp = (2*eps*Vbi*Nd/(q*Na*10^6*(Na+Nd)))^0.5;
18 disp(Wp,"depletion width on the p-side (in m) = ")

```

---

**Scilab code Exa 4.2** effect of the generation recombination current in pn diode

```

1 kT = 26*10^-3; //in eV
2 t = 10^-6;
3 q = 1.6*10^-19;
4 A = 10^-3; //in square cm
5 ni = 1.5*10^10;
6 eE = 3.2*10^4;
7 Igr0 = q*ni*A*%pi*kT/(2*t*eE);
8 disp(Igr0,"prefactor of the generation-recombination
   current (in Ampere) =")
9 V1 = 0.2;
10 Igr1 = Igr0*(exp(0.5*V1/kT)-1);
11 disp(Igr1,"generation-recombination current (in A)
   at bias of 0.2 V =")
12 V2 = 0.6;
13 Igr2 = Igr0*(exp(0.5*V2/kT)-1);
14 disp(Igr2,"generation-recombination current (in A)
   at bias of 0.6 V =")

```

---

**Scilab code Exa 4.3** diode current at a forward bias

```

1 n = 10^17;
2 p = 10^17;
3 Tn = 10^-7;
4 Tp = 10^-7;
5 Dn = 30;
6 Dp = 10;

```

```

7 A = 10^-4;
8 t = 10^-8;
9 V1 = 0.5;
10 V2 = 0.6;
11 E1 = 6.94 * 10^4;
12 E2 = 5.74 * 10^4;
13 np = 2.25*10^3;
14 pn = 2.25*10^3;
15 Ln = 17.32*10^-4;
16 Lp = 10*10^-4;
17 Vbi = 0.817;
18 q = 1.6*10^-19;
19 I0 = q*A*(Dp*pn/Lp + Dn*np/Ln);
20 disp(I0,"prefactor in the ideal diode equation = ")
21 Igr01 = q*ni*A*%pi*kT/(2*t*E1);
22 Igr02 = q*ni*A*%pi*kT/(2*t*E2);
23 disp(Igr01,"prefactor to the recombination-
      generation current at 0.5V = ")
24 disp(Igr02,"prefactor to the recombination-
      generation current at 0.6V")
25 I1 = I0*exp(V1/kT)+Igr01*exp(0.5*V1/kT);
26 disp(I1,"Current at 0.5 V = ")
27 I2 = I0*exp(V2/kT)+Igr02*exp(0.5*V2/kT);
28 disp(I2,"Current at 0.6 V = ")
29 n = (log(I2/I1)*kT/(V2-V1))^-1;
30 disp(n,"ideality factor of the diode in the given
      range =")

```

---

#### Scilab code Exa 4.4 breakdown voltage of diode

```

1 Na = 10^19;
2 Nd = 10^16;
3 E1 = 4*10^5;
4 E2 = 10^7;
5 eps0 = 8.84*10^-14; //in F/m

```

```

6 eps = 11.9*eps0;
7 q = 1.6*10^-19;
8 V1 = eps*E1^2/(2*Nd*q);
9 disp(V1,"breakdown voltage for Si = ")
10 V2 = eps*E2^2/(2*Nd*q);
11 disp(V2,"breakdown voltage for diamond = ")

```

---

#### Scilab code Exa 4.5 e h recombination time

```

1 p = 10^21;
2 t0 = 0.6*10^-9;
3 kT = 26*10^-3; //in eV
4 m0 = 0.91 * 10^-30; //in kg
5 m1 = 0.067*m0;
6 m2 = 0.45*m0;
7 q = 1.6*10^-19;
8 h = 1.05*10^-34;
9 tri = (0.5*p/t0)*(2*%pi*h*h/kT/q/(m1+m2))^1.5;
10 tr = 1/tri;
11 disp(tr)

```

---

#### Scilab code Exa 4.6 internal radiative efficiency for diodes

```

1 t0 = 0.6*10^-9;
2 tnr = 10^-7;
3 p = 10^22;
4 kT = 26*10^-3; //in eV
5 m0 = 0.91 * 10^-30; //in kg
6 m1 = 0.067*m0;
7 m2 = 0.45*m0;
8 tri = (0.5*p/t0)*(2*%pi*h*h/kT/q/(m1+m2))^1.5;
9 tr1 = 1/tri;

```

```

10 disp(tr1,"When the p-type doping is  $10^{16}$  cm3 ,
    the hole density is low and the e-h recombination
    time (in s) for the injected electrons = ")
11 tr2 = t0*((m1+m2)/m1)^1.5;
12 disp(tr2,"In the case where the p-doping is high ,
    the recombination time(in s) = ")
13 nqr1 = (1+tr1/tnr)^-1;
14 disp(nqr1,"For the low-doping case , the internal
    quantum efficiency for the diode = ")
15 nqr2 = (1+tr2/tnr)^-1;
16 disp(nqr2,"For the more heavily doped p region , the
    internal quantum efficiency for the diode = ")

```

---

#### Scilab code Exa 4.7 injection efficiency of LED

```

1 Dn = 30;
2 Dp =15;
3 Na = 5*10^16;
4 Nd = 5*10^17;
5 q = 1.6*10^-19;
6 tn = 10^-8;
7 tp = 10^-7;
8 ni = 1.84*10^6;
9 np = ni^2/Na;
10 pn = ni^2/Nd;
11 Ln = (Dn*tn)^0.5;
12 Lp = (Dp*tp)^0.5;
13 disp(Ln,"diffusion length , Ln (in cm) = ")
14 disp(Lp,"diffusion length , Lp (in cm) = ")
15 n = (q*Dn*np/Ln)/((q*Dn*np/Ln)+(q*Dp*pn/Lp));
16 disp(n,"injection efficiency (assuming no
    recombination via traps) = ")

```

---

Scilab code Exa 4.8 photon flux and optical power generated by LED

```
1 V1 = 1;
2 nqr = 0.5;
3 A = 10^-2;
4 Dn = 30;
5 Dp = 15;
6 Na = 5*10^16;
7 Nd = 5*10^17;
8 q = 1.6*10^-19;
9 tn = 10^-8;
10 tp = 10^-7;
11 ni = 1.84*10^6;
12 np = ni^2/Na;
13 pn = ni^2/Nd;
14 Ln = (Dn*tn)^0.5;
15 Lp = (Dp*tp)^0.5;
16 kT = 26*10^-3; //in eV
17 In = A*q*Dn*np/Ln*(exp(V1/kT)-1);
18 disp(In,"The electron current injected into the p-
    region will be responsible for the photon
    generation. This current (in A) = ")
19 Iph = In*nqr/q;
20 disp(Iph,"photons generated per second = ")
21 E = 1.41; //in eV
22 P = Iph*E*q;
23 disp(P,"Each photon has an energy of 1.41 eV (=
    bandgap of GaAs). The optical power (in W) = ")
```

---

# Chapter 5

## SEMICONDUCTOR JUNCTIONS

Scilab code Exa 5.2 SCHOTTKY BARRIER

```
1 A = 10^-3;
2 Na = 10^19;
3 Nd = 10^16;
4 q = 1.6*10^-19;
5 Tp = 10^-6;
6 Tn = 10^-6;
7 Dp = 10.5;
8 kT = 26*10^-3; //in eV
9 T = 300;
10 Vs = -0.67;
11 Vf = 0.3;
12 eRc = 110;
13 Is = A*eRc*T^2*exp(Vs/kT);
14 disp(Is,"Reverse saturation current (in Ampere) = ")
15 I = Is*exp(Vf/kT);
16 disp(I,"For a forward bias of 0.3 V, the current(in
    Ampere) = ")
17 Lp = (Dp*Tp)^0.5;
18 disp(Lp, "Lp(in cm) = ")
```

```
19 pn = 2.25*10^4;
20 I0 = A*q*Dp*pn/Lp;
21 disp(I0," Saturation current (in Ampere) = ")
22 disp("For the p-n diode to have the same current
      that the Schottky diode has at 0.3 V, the voltage
      required is 0.71 V.")
```

---

## Chapter 6

# BIPOLAR JUNCTION TRANSISTORS

Scilab code Exa 6.1 emitter efficiency of BJT and HBT

```
1 Nde = 5*10^17;
2 Nab = 10^17;
3 Db = 100;
4 De = 15;
5 Wb = 0.5*10^-4;
6 dEg = 0.36;
7 Le = 1.5*10^-4;
8 ni = 2.2*10^6;
9 kT = 26*10^-3; //in eV
10 peo = ni^2/Nde;
11 nbo = ni^2/Nab;
12 disp(peo,"emitter minority carrier concentrations (
    in per cm cube)= ")
13 disp(nbo,"base minority carrier concentrations (in
    per cm cube) = ")
14 gammae = 1- (peo*De*Wb)/(nbo*Db*Le);
15 disp(gammae,"emitter efficiency = ")
16 peo1 = peo*exp(-dEg/kT);
17 disp(peo1,"In the HBT, the value of peo is greatly
```

```

    suppressed. The new value(in per cm cube) = ")
18 gamma1 = 1- (peo1*De*Wb)/(nbo*Db*Le);
19 disp(gamma1,"emitter efficiency = ")
20 disp("In this case the emitter efficiency is
    essentially unity")

```

---

**Scilab code Exa 6.2** change in the base width and collector current with voltage and Early voltage

```

1 Wb = 10^-4;
2 Vcb1 = 1;
3 Vcb2 = 5;
4 q = 1.6*10^-19;
5 Db = 20;
6 Vbe = 0.7;
7 kT = 26*10^-3; //in eV
8 ni = 1.5*10^10;
9 Nab = 5*10^16;
10 Nde = 5*10^15;
11 eps0 = 8.84*10^-14; //in F/m
12 eps = 11.9*eps0;
13 Vbi = kT*log(Nab*Nde/ni^2)
14 disp(Vbi,"Built in voltage (in V) = ")
15 dW2 = 2*eps*(Vbi+Vcb1)*Nde/(q*Nab*(Nab+Nde));
16 dW = sqrt(dW2);
17 disp(dW,"The extent of depletion into the base side
    (in cm) = ")
18 Wbn = Wb - dW;
19 disp(Wbn,"neutral base width (in cm) = ")
20 dW1 = (2*eps*(Vbi+Vcb2)*Nde/(q*Nab*(Nab+Nde)))^0.5;
21 disp(dW1,"When the collector-base voltage increases
    to 5 V, extent of depletion into the base side (
    in cm)")
22 Wbn1 = Wb - dW1;
23 disp(Wbn1,"neutral base width (in cm) = ")

```

```

24 nbo = ni^2/Nab;
25 disp(nbo,"base minority carrier concentrations (in
    per cm cube) = ")
26 Jc1 = q*Db*nbo/Wbn*exp(Vbe/kT);
27 disp(Jc1,"For the base-collector bias of 1 V,
    collector current density (in A/square cm)")
28 Jc2 = q*Db*nbo/Wbn1*exp(Vbe/kT);
29 disp(Jc2,"For the base-collector bias of 5 V,
    collector current density (in A/square cm)")
30 slope = (Jc2-Jc1)/(Vcb2-Vcb1);
31 disp(slope,"The slope of the Jc vs. VCE curve = ")
32 Va = Jc2/slope - (Vcb2+Vbe);
33 disp(Va, "Early voltage (in V) = ")

```

---

**Scilab code Exa 6.3** maximum base width for a given beta

```

1 kT = 26*10^-3; //in eV
2 ni = 1.5*10^10;
3 eps0 = 8.84*10^-14; //in F/m
4 eps = 11.9*eps0;
5 q = 1.6*10^-19;
6 Nde = 10^18;
7 Nab = 10^17;
8 Ndc = 5*10^16;
9 Db = 30;
10 Lb = 15*10^-4;
11 De = 10;
12 Le = 5*10^-4;
13 Vf = 1;
14 B = 100;
15 Vr = 5;
16 Wbn = Db*Nde*Le/(De*Nab*B);
17 disp(Wbn,"neutral base width (in cm) = ")
18 Vbi = kT*log(Nab*Ndc/ni^2);
19 disp(Vbi,"built in voltage (in V) = ")

```

```

20 //dW = (2*eps*(Vbi+Vr)*Nde/(q*Nab*(Nab+Nde)))^0.5;
21 dW = (2*eps*(Vbi+Vr)*Ndc/(q*Nab*(Nab+Ndc)))^0.5;
22 disp(dW,"the depletion width (in cm) on the base
    side of the EBJ for a 5 volt bias at the base
    collector junction")
23 Wb = Wbn + dW;
24 disp(Wb, "base width (in cm) = ")

```

---

**Scilab code Exa 6.4** output conductance and emitter efficiency and gain

```

1 kT = 26*10^-3; //in eV
2 ni = 1.5*10^10;
3 Vbe = 0.7;
4 q = 1.6*10^-19;
5 Nde = 10^18;
6 Nab = 10^17;
7 Ndc = 10^16;
8 Db = 30;
9 Lb = 10*10^-4;
10 Wb = 10^-4;
11 De = 10;
12 Le = 10*10^-4;
13 We = 10^-4;
14 A = 4*10^-6;
15 Vf = 1;
16 Vr1 = 5;
17 Vr2 = 6;
18 Vbi = kT*log(Nab*Ndc/ni^2);
19 disp(Vbi,"built in voltage (in V) = ")
20 dW1 = (2*eps*(Vbi+Vr1)*Ndc/(q*Nab*(Nab+Ndc)))^0.5;
21 dW2 = (2*eps*(Vbi+Vr2)*Ndc/(q*Nab*(Nab+Ndc)))^0.5;
22 disp(dW1,"depletion width (in cm) on the base side
    of the BCJ at 5 V = ")
23 disp(dW2,"depletion width (in cm) on the base side
    of the BCJ at 6 V = ")

```

```

24 Wbn1 = Wb - dW1;
25 disp(Wbn1,"neutral base width (in cm) at 5V = ")
26 Wbn2 = Wb - dW2;
27 disp(Wbn2,"neutral base width (in cm) at 6V = ")
28 gammae = 1- (Nab*De*Wbn1)/(Nde*Db*We)
29 disp(gammae,"emitter efficiency (for a narrow
    emitter of width We) = ")
30 B = 1 - Wbn1^2/2/Lb^2;
31 disp(B,"the base transport factor = ")
32 alpha = B*gammae;
33 disp(alpha,"alpha = ")
34 betae = alpha/(1-alpha);
35 disp(betae,"current gain = ")
36 Nbo = ni^2/Nab;
37 Ic1 = q*A*Db*Nbo/Wbn1*(exp(Vf/kT)-1);
38 disp(Ic1,"collector current (in A) at 5V = ")
39 Ic2 = q*A*Db*Nbo/Wbn2*(exp(Vf/kT)-1);
40 disp(Ic2,"collector current (in A) at 6V = ")
41 g0 = Ic2-Ic1/(6-5);
42 disp(g0,"The output conductance = ")

```

---

## Chapter 7

# TEMPORAL RESPONSE OF DIODES AND BIPOLAR TRANSISTORS

Scilab code Exa 7.1 cutoff frequency of transistor

```
1 Ie = 1.5*10^-3;
2 Cje = 2*10^-12;
3 Wb = 0.4*10^-4;
4 Db = 60;
5 Wdc = 2*10^-4;
6 Rc = 30;
7 Ct = 0.4*10^-12;
8 vs = 5*10^6;
9 kT = 26*10^-3; //in eV
10 q = 1.6*10^-19;
11 Re = kT/Ie;
12 disp(Re,"The emitter resistance (in ohm) = ")
13 Te = Re*Cje;
14 disp(Te,"Te (in s) = ")
15 Tt = Wb^2/2/Db;
16 disp(Tt,"base transit time (in s) = ")
17 Tc = 0.5*Wdc/vs;
```

```
18 disp(Tc," collector transit time (in s) = ")
19 Tcc = Rc*Ct;
20 disp(Tcc," collector charging time (in s) = ")
21 Tec = Te+Tt+Tc+Tcc;
22 disp(Tec," total time (in s) = ")
23 ft = (2*%pi*Tec)^-1;
24 disp(ft," cutoff frequency (in Hz) = ")
```

---

# Chapter 8

## FIELD EFFECT TRANSISTORS

Scilab code Exa 8.1 gate current density for MESFET

```
1 phiB = 0.8;
2 N = 10^17;
3 Dp = 20;
4 Lp = 10^-4;
5 A = 8;
6 kT = 26*10^-3; //in eV
7 T = 300;
8 q = 1.6*10^-19;
9 Js = A*T^2*exp(-phiB/kT);
10 disp(Js,"For the Schottky case , Js (in A/cm2) = ")
11 ni = 1.84*10^6;
12 pn = ni^2/N;
13 disp(pn)
14 J0 = q*Dp*pn/Lp;
15 disp(J0,"For the p+-gate we have from p-n diode
    theory , J0 (in A/cm2) = ")
```

---

### Scilab code Exa 8.5.1 2DEG concentration for MODFET

```
1 phiB = 0.9;
2 Nd = 10^18;
3 dEc = 0.24;
4 eps0 = 8.84*10^-14; //in F/m
5 epsb = 12.2*eps0;
6 ds = 30*10^-8;
7 d = 350*10^-8;
8 VG1 = 0;
9 VG2 = -0.5;
10 q = 1.6*10^-19;
11 Vp2 = q*Nd*(d-ds)^2/epsb;
12 disp(Vp2,"The parameter Vp2 (in V) of this structure
    = ")
13 Voff = phiB - Vp2 - 0.24;
14 disp(Voff,"threshold voltage (in V)")
15 Ns1 = -epsb*Voff/q/d;
16 Ns2 = epsb*(VG2-Voff)/q/d;
17 disp(Ns1,"The 2DEG carrier concentration (in per
    sqaure cm) at 0 V = ")
18 disp(Ns2,"The 2DEG carrier concentration (in per
    sqaure cm) at -0.5 V = ")
```

---

## Chapter 9

# FIELD EFFECT TRANSISTORS MOSFET

Scilab code Exa 9.1 maximum depletion width in a MOS capacitor

```
1 kT = 26*10^-3; //in eV
2 T = 300;
3 q = 1.6*10^-19;
4 Na = 10^16;
5 ni = 1.5*10^10;
6 eps0 = 8.84*10^-14; //in F/m
7 eps = 11.9*eps0;
8 phiF = kT*log(Na/ni);
9 disp(phiF,"the potential F (in V) = ")
10 W = (4*eps*phiF/(q*Na))^0.5;
11 disp(W,"The corresponding space charge width(in cm)
    = ")
```

---

Scilab code Exa 9.2 MOS capacitor

```
1 kT = 26*10^-3; //in eV
```

```

2 T = 300;
3 q = 1.6*10^-19;
4 Wf = 4.1;
5 EA1 = 0.9;
6 EA2 = 4.15;
7 Na = 10^14;
8 Eg = 1.11;
9 Efi = Eg/2;
10 Ef = Efi + kT*log(Na/ni);
11 disp(Ef,"Ef = ")
12 disp("below the conduction band")
13 Vfb = Wf - (EA2+Ef);
14 disp(Vfb,"V(fb) = ")

```

---

### Scilab code Exa 9.3 MOS capacitor threshold voltage

```

1 Na = 3*10^16;
2 t = 500*10^-8;
3 Vfb = -1.13;
4 T = 300;
5 kT = 26*10^-3; //in eV
6 q = 1.6*10^-19;
7 ni = 1.5*10^10;
8 eps0 = 8.85*10^-14; //in F/m
9 eps = 11.9*eps0;
10 eps1 = 3.9*eps0;
11 c = 10^11;
12 phiF = kT*log(Na/ni);
13 disp(phiF,"The position of the Fermi level (in V) is
    given by (measured from the intrinsic Fermi
    level)")
14 Qs = (4*eps*phiF*q*Na)^0.5;
15 disp(Qs,"Under the assumption that the charge Qs is
    simple NaW where W is the maximum depletion width
    , we get Qs (in C per cm2)= ")

```

```

16 Vt = Vfb+2*phiF+(Qs*t/eps1);
17 disp(Vt,"In the absence of any oxide charge, the
    threshold voltage (in V) = ")
18 dVt = c*q*(t/eps1);
19 disp(dVt,"In the case where the oxide has trap
    charges, the threshold voltage is shifted by VT
    (in V)= ")

```

---

**Scilab code Exa 9.4** MOS capacitor threshold voltage and channel conductivity

```

1 Na = 5*10^16;
2 phiMS = -0.5;
3 un = 600;
4 up = 200;
5 T = 300;
6 kT = 26*10^-3; //in eV
7 q = 1.6*10^-19;
8 ni = 1.5*10^10;
9 eps0 = 8.85*10^-14; //in F/m
10 eps = 11.9*eps0;
11 eps1 = 3.9*eps0;
12 psiS = 2*phiF;
13 w = 200*10^-8;
14 sigma_fb= Na*q*up;
15 disp(sigma_fb," (fb) (in per ohm-cm) = ")
16 sigma_inv = Na*q*un;
17 disp(sigma_inv," (inv) (in per ohm-cm) = ")
18 phiF = kT*log(Na/ni);
19 disp(phiF," F (in V) = ")
20 Vt = phiMS + psiS + 1.637;
21 disp(Vt,"the threshold voltage (in V) = ")

```

---

**Scilab code Exa 9.6** oxide capacitance and capacitance at the flat band and the minimum capacitance at threshold for MOS capacitor

```

1 Na = 10^16;
2 t = 500*10^-8;
3 kT = 26*10^-3; //in eV
4 q = 1.6*10^-19;
5 ni = 1.5*10^10;
6 eps0 = 8.85*10^-14; //in F/m
7 eps = 11.9*eps0;
8 eps1 = 3.9*eps0;
9 Cox = eps1/t;
10 disp(Cox,"The oxide capacitance (in F/cm2) = ")
11 phiF = kT*log(Na/ni);
12 disp(phiF," F (in V) = ")
13 Wmax = (4*eps*phiF/(q*Na))^0.5;
14 disp(Wmax,"The maximum depletion width (in cm) = ")
15 Cmin = eps1/(t+(eps1*Wmax/eps));
16 disp(Cmin,"The minimum capacitance (in F/cm2) = ")
17 Cfb = eps1/(t+(eps1/eps*(kT*eps/q/Na)^0.5));
18 disp(Cfb,"The capacitance (in F/cm2) under flat band
           conditions = ")
19 disp("Note that Cfb is      80% of Cox and Cmin is
           33% of Cox.")

```

---

**Scilab code Exa 9.7** saturation current of MOSFET

```

1 L = 1.5*10^-4;
2 Z = 25*10^-4;
3 un = 600;
4 Na = 10^16;
5 dox = 500*10^-8;
6 Qss = 10^11;
7 phiMS = -1.13;
8 V = 5;

```

```

 9 kT = 26*10^-3; //in eV
10 q = 1.6*10^-19;
11 ni = 1.5*10^10;
12 eps0 = 8.85*10^-14; //in F/m
13 eps = 11.9*eps0;
14 eps1 = 3.9*eps0;
15 phiF = kT*log(Na/ni);
16 disp(phiF,"The Fermi level position (in V) for the
    device = ")
17 Cox = eps/dox;
18 Vfb = phiMS - 0.23;
19 disp(Vfb,"The flat band voltage (in V) = ")
20 Qs = (4*eps*phiF*q*Na)^0.5;
21 Vt = Vfb + 2*phiF + (Qs*dox/eps1);
22 disp(Vt,"The threshold voltage (in V) = ")
23 Id = Z*un*eps1*(V-Vt)^2/(2*L*dox);
24 disp(Id,"The saturation current (in A) = ")

```

---

### Scilab code Exa 9.8 drain current for NMOS device

```

1 kT = 26*10^-3; //in eV
2 q = 1.6*10^-19;
3 ni = 1.5*10^10;
4 eps0 = 8.85*10^-14; //in F/m
5 eps = 11.9*eps0;
6 eps1 = 3.9*eps0;
7 phiMS = 0;
8 Na = 4*10^14;
9 dox = 200*10^-8;
10 L = 10^-4;
11 Z = 10*10^-4;
12 Vgs = 5;
13 Vd = 4;
14 un = 700;
15 phiF = kT*log(Na/ni);

```

```

16 disp(phiF,"The Fermi level position (in V) for the
    device = ")
17 Qs = (4*eps*phiF*q*Na)^0.5;
18 Vt = 2*phiF + (Qs*dox/eps1);
19 disp(Vt,"The threshold voltage (in V) = ")
20 Vds = Vgs - Vt;
21 disp(Vds,"saturation voltage (in V) for a gate bias
    of 5 V = ")
22 Id = Z*un*eps1*(Vds)^2/(2*L*dox);
23 disp(Id,"The saturation current (in A) = ")

```

---

**Scilab code Exa 9.9** threshold voltage in n channel MOSFET

```

1 Z = 10*10^-4;
2 L = 2*10^-4;
3 Cox = 10^-7;
4 Vds = 0.1;
5 Vgs1 = 1.5;
6 Id1 = 50*10^-6;
7 Vgs2 = 2.5;
8 Id2 = 80*10^-6;
9 slope = Id2-Id1/(Vgs2-Vgs1);
10 Vt = -Id2/slope + Vgs2;
11 disp(Vt,"the threshold voltage (in V) = ")

```

---

**Scilab code Exa 9.10** shift in the threshold voltage arising from source body bias

```

1 kT = 26*10^-3; //in eV
2 q = 1.6*10^-19;
3 ni = 1.5*10^10;
4 eps0 = 8.85*10^-14; //in F/m
5 eps = 11.9*eps0;

```

```

6 eps1 = 3.9*eps0;
7 Na = 2*10^16;
8 dox = 500*10^-8;
9 Vsb = 1;
10 phiF = kT*log(Na/ni);
11 disp(phiF,"The Fermi level position (in V) for the
    device = ")
12 Cox = eps1/dox;
13 disp(Cox,"The oxide capacitance (in F/cm2) = ")
14 dVt = (2*q*eps*Na)^0.5/Cox*((2*phiF+1)^0.5 - (2*phiF
    )^0.5);
15 disp(dVt,"The change in the threshold voltage (in V)
    = ")

```

---

#### Scilab code Exa 9.11 n channel MOSFET characteristics

```

1 kT = 26*10^-3; //in eV
2 q = 1.6*10^-19;
3 ni = 1.5*10^10;
4 eps0 = 8.85*10^-14; //in F/m
5 eps = 11.9*eps0;
6 eps1 = 3.9*eps0;
7 Na = 10^14;
8 dox = 500*10^-8;
9 phiMS = -0.83;
10 t = 0.1*10^-4;
11 dVt = 0.5;
12 phiF = kT*log(Na/ni);
13 disp(phiF,"The Fermi level position (in V) for the
    device = ")
14 Qs = (4*eps*phiF*q*Na)^0.5;
15 Vt = phiMS+2*phiF + (Qs*dox/eps1);
16 disp(Vt,"The threshold voltage (in V) = ")
17 disp("In this device there is an inversion layer
    formed even at zero gate bias and the device is

```

in the depletion mode. To increase the threshold voltage by + 0.5 V, i.e., to convert the device into an enhancement-mode device, we need to place more negative charge in the channel. If we assume that the excess acceptors are placed close to the semiconductor-oxide region (i.e., within the distance  $W_{max}$ ), the shift in threshold voltage is simply ( $Na2D$  is the areal density of the acceptors implanted)"

```
18 Na2D = dVt/dox*eps1/q;  
19 disp(Na2D, "the areal density (in per cm2) of the  
    acceptors implanted = ")  
20 Na1 = Na2D/t;  
21 disp(Na1, "The dopants are distributed over a  
    thickness of 0.1 m, the dopant density (in per  
    cm3) = ")
```

---