Scilab Textbook Companion for Materials Science by R. S. Khurmi and R. S. Sedha¹

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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.

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Chapter 2

structure of atoms

Scilab code Exa 2.1 radius

```
1 //Example 2.1 : radius of the first bohr"s orbit
2 clc;
3 clear;
4 close;
5 //given data :
6 format('v',5)
7 ep=8.854*10^-12;//
8 h=6.626*10^-34;//
9 m=9.1*10^-31;//in Kg
10 e=1.602*10^-19;//
11 r1=((ep*(h^2))/((%pi*m*(e^2))));//
12 disp(r1*10^10,"radius,r1(in angstrom) = ")
```

Scilab code Exa 2.2 radius

```
2 //Example 2.2 : radius of the second bohr"s orbit 3 clc;
```

```
4 clear;
5 close;
6 //given data :
7 format('v',6)
8 r1_h=0.529; // radius for hydrozen atom in Angstrum
9 n1=1; // for the first bohr's orbit of electron in hydrozen atom
10 Z1=1; // for the first bohr's orbit of electron in hydrozen atom
11 k=(r1_h*Z1)/n1^2; // where k is constant
12 n2=2; // for the second bohr orbit
13 Z2=2; //for the second bohr orbit
14 r2_he=k*(n2^2/Z2);
15 disp(r2_he, "radius of the second bohr orbit, r2( Angstrom) = ")
```

Scilab code Exa 2.3 ratio of energy

```
1 // Example 2.3: to prove
2 clc:
3 clear;
4 close;
5 \text{ Z=1;} // \text{assume}
6 n1=1; //orbit 1
7 \text{ n2=2; //orbit } 2
8 \text{ n3=3; //orbit } 3
9 e1=((-13.6*Z)/(n1^2)); //energy for the first orbit
10 e2=((-13.6*Z)/(n2^2));//energy for the second orbit
11 e3=((-13.6*Z)/(n3^2));//energy for the third orbit
12 e31=e3-e1; //energy emitted by an electron jumping
      from orbit nuber 3 to orbit nimber 1
13 e21=e2-e1; //energy emitted by an electron jumping
      from orbit nuber 2 to orbit nimber 1
14 re=e31/e21; //ratio of energy
15 disp(re, "ratio of energy for an electron to jump
```

```
from orbit 3 to orbi 1 and from orbit 2 to orbit 1 is 32/27")
```

Scilab code Exa 2.4 velocity

```
1 //Example 2.4 : velocity
2 clc;
3 clear;
4 close;
5 //given data :
6 format('v',8)
7 h=6.626*10^-34;
8 e=1.6*10^-19;
9 epsilon_o=8.825*10^-12;
10 n=1;
11 Z=1;
12 vn=(Z*e^2)/(2*epsilon_o*n*h);
13 disp(vn,"velocity,vn(m/s) = ")
```

Scilab code Exa 2.5 orbital frequency

```
1 //Example 2.5 : velocity
2 clc;
3 clear;
4 close;
5 //given data :
6 n=1;
7 Z=1;
8 k=6.56*10^15; // k is constant
9 fn=k*(Z^2/n^3);
10 disp(fn, "orbital frequency, fn(Hz) = ")
```

Scilab code Exa 2.6.a energy of photon emitted

```
//Example 2.6.a : the energy of the photon emitted
clc;
clear;
close;
format('v',5);
//given data :
Z=1;//for hydrozen
n1=3;
n2=2;
E3=-(13.6*Z^2)/n1^2;
E2=-(13.6*Z^2)/n2^2;
del_E=E3-E2;
disp(del_E,"the energy of photon emitted, del_E(eV) =
")
```

Scilab code Exa 2.6.b frequency

```
1 //Example 2.6.b : frequency
2 clc;
3 clear;
4 close;
5 //given data :
6 format('v',9)
7 Z=1;//for hydrozen
8 n1=3;
9 n2=2;
10 m=6.626*10^-34;// mass of electron in kg
11 E3=-(13.6*Z^2)/n1^2;
12 E2=-(13.6*Z^2)/n2^2;
13 del_E=E3-E2;
```

```
14 E=del_E*1.6*10^-19; // in joules
15 v=(E/m);
16 disp(v, "frequency of the photon emitted, v(Hz) = ")
```

Scilab code Exa 2.6.c wavelength

```
1 //Example 2.6.c : wave length of the photon emitted
2 clc;
3 clear;
4 close;
5 //given data :
6 format('v',9)
7 Z=1; // for hydrozen
8 n1=3;
9 n2=2;
10 m=6.626*10^-34; // mass of electron in kg
11 C=3*10^8;
12 E3=-(13.6*Z^2)/n1^2;
13 E2 = -(13.6*Z^2)/n2^2;
14 \text{ del}_{\text{E}}=\text{E3}-\text{E2};
15 E=del_E*1.6*10^-19;
16 \text{ v=E/m};
17 lamda=C/v;
18 disp(lamda, "wavelength of the photon emitted, (m) = "
```

Chapter 3

crystal structure

Scilab code Exa 3.1 miller indices

```
1 //Example 3.1: miller indices
2 clc;
3 clear;
4 close;
5 //given data
6 \text{ x1=1;} //
7 x2=1; //
8 x3=2;//
9 h1=1/x1;//
10 h2=1/x2;//
11 h3=1/x3;//
12 disp("Miller indices of the plane (112) are: "+
      string(h1)+","+string(h2)+","+string(h3))
13 x11=0; //
14 x21=0; //
15 \times 31 = 1; //
16 h11=%inf;//
17 h21=%inf;//
18 h31=1/x31; //
19 disp("Miller indices of the plane (001) are : "+"
      string(h11)+","+string(h21)+","+string(h31))
```

```
20 x111=1;//
21 x211=0;//
22 x311=1;//
23 h111=1/x111;//
24 h211=%inf;//
25 h311=1/x311;//
26 disp("Miller indices of the plane (101) are: "+
string(h111)+","+string(h211)+","+string(h311))
```

Scilab code Exa 3.2 miller indices

```
1 //Example 3.2: miller indices
2 clc;
3 clear;
4 close;
5 //given data
6 x1=0; //
7 x2=2; //
8 x3=0; //
9 h1=%inf;//
10 h2=1/x2;//
11 h3=%inf;//
12 disp("Miller indices of the plane (020) are: "+
      string(h1)+","+string(h2)+","+string(h3))
13 x11=1; //
14 x21=2;//
15 \times 31 = 0; //
16 h11=1/x11; //
17 h21=1/x21;//
18 h31=%inf;//
19 disp("Miller indices of the plane (120) are: "+
      string(h11)+","+string(h21)+","+string(h31))
20 \times 111 = 2; //
21 \times 211 = 2; //
22 \times 311 = 0; //
```

Scilab code Exa 3.3 miller indices

```
1 // Example 3.3: miller indices
2 clc;
3 clear;
4 close;
5 x=1/2; //
6 x1=1/x; //
7 r2=0; //
8 r3=0;//
9 \times 10 = -1; //
10 x2=1/-x10;//
11 r4=0;//
12 r5=0;//
13 disp("miller indices (Case 1) of the given plane are
      "+string(x1)+" : "+string(r2)+" : "+string(r3)+"
14 disp("miller indices (Case 2) of the given plane are
      "+string(x2)+" : "+string(r3)+" : "+string(r4)+"
```

Scilab code Exa 3.4 miller indices

```
1 // Example 3.4: miller indices 2 clc;
```

```
3 clear;
4 close;
5 a=0.529;//
6 b=1;//
7 c=0.477;//
8 a1=0.264;//
9 b1=1;//
10 c1=0.238;//
11 r1=round(a/a1);//
12 r2=b/b1;//
13 r3=round(c/c1);//
14 disp("miller indices of the given plane are "+string (r1)+": "+string(r2)+": "+string(r3)+"")
```

Scilab code Exa 3.5 miller indices

```
1 //Example 3.5: miller indices
2 clc;
3 clear;
4 close;
5 //given data
6 \text{ x1=1;} //
7 x2=1; //
8 x3=0; //
9 h1=1/x1//
10 h2=1/x2;//
11 h3=%inf;//
12 disp("Miller indices of the plane (110) are: "+
      string(h1)+","+string(h2)+","+string(h3))
13 x11=1; //
14 x21=1; //
15 x31=1;//
16 h11=1/x11;//
17 h21=1/x21;//
18 h31=1/x31; //
```

```
19 disp("Miller indices of the plane (111) are: "+ string(h11)+","+string(h21)+","+string(h31))
```

Scilab code Exa 3.9 atoms per unit cell

```
1 // Example 3.9: atoms per unit cell
2 clc;
3 clear;
4 close;
5 c=8;//corners
6 f=6;//faces
7 nf=(1/2)*f;//no. of atoms in all six faces
8 nc=(1/8)*c;//no. of atoms in all corners
9 ta=nf+nc;//
10 disp(ta,"total number of atoms are")
```

Scilab code Exa 3.10 diameter

```
//Example 3.10 : largest diameter
clc;
clc;
clear;
close;
//given data :
format('v',6)
a=3.61; // edge length in angstrum
r=(a*sqrt(2))/4;
d=2*r;
disp(d,"largest diameter,d(angstrom) = ")
```

Scilab code Exa 3.11 volume change

```
1 //Example 3.11 : volume change in percentage
2 clc;
3 clear;
4 close;
5 //given data :
6 r_bcc=0.1258; // in nm
7 r_fcc=0.1292; // in nm
8 a_bcc=(r_bcc*4)/sqrt(3);
9 a_fcc=(r_fcc*4)/sqrt(2);
10 v_fcc=(a_fcc)^3; // in nmn^3
11 v_bcc=(a_bcc)^3; // in nmn^3
12 V=((v_fcc-v_bcc)/v_bcc)*100;
13 disp(V,"volume change in percentage, V(%) = ")
```

Scilab code Exa 3.12 number of atoms

```
1
2  //Example 3.12 : number of atom/mm^2
3  clc;
4  clear;
5  close;
6  format('v',8)
7  //given data :
8  a=3.03*10^-7;  // lattice constant in mm
9  A=1/a^2;// for 100 planes
10 B=0.707/a^2;//for(110) planes
11 C=0.58/a^2;// for(111) planes
12 disp(A,"number of atom for (100) plane, = ")
13 disp(B,"number of atoms for (110) plan, = ")
14 disp(C,"number of atoms for (111) plan, = ")
```

Scilab code Exa 3.13 number of atoms

```
1
2  //Example 3.13 : number of atom/mm^2 of planes
3  clc;
4  clear;
5  close;
6  //given data :
7  format('v',9)
8  a=2.87*10^-7; // lattice constant in mm
9  A=1/a^2; // for 100 planes
10  B=1.414/a^2; //for(110) planes
11  C=1.732/a^2; // for(111) planes
12  disp(A,"number of atom for (100) plane, = ")
13  disp(B,"number of atoms for (110) plan, = ")
14  disp(C,"number of atoms for (111) plan, = ")
```

Scilab code Exa 3.14 number of atoms

```
1
2  //Example 3.14 : number of atom/mm^2 surface area
3  clc;
4  clear;
5  close;
6  //given data :
7  a=4.93*10^-7; // lattice constant in mm
8  A=2/a^2;// for 100 planes
9  B=1.414/a^2;//for(110) planes
10  C=2.31/a^2;// for (111) planes
11  disp(A,"number of atoms for (100) plane) = ")
12  disp(B,"number of atoms for (110) plan = ")
13  disp(C,"number of atoms for (111) plan = ")
```

Scilab code Exa 3.15 planar density

```
1 //Example 3.15 : planar density
2 clc;
3 clear;
4 close;
5 //given data :
6 format('v',9)
7 a=0.143*10^-6; // atomic radius in mm
8 A=2.31/(a^2);// for (111) planes
9 disp(A, "atom, A(atoms/mm^2) = ")
10 // answer is wrong in book
```

Scilab code Exa 3.16 volume

```
1 //Example 3.16 : volume
2 clc;
3 clear;
4 close;
5 format('v',7)
6 //given data :
7 a=0.2665; // in mm
8 c=0.4947; // in mm
9 V=(3*sqrt(3)*a^2*c)/2;
10 disp(V,"volume, V(mm^3) = ")
```

Scilab code Exa 3.17 packing efficiency and lattice parameter

```
7 //given data :
8 r=1.22// in angstrum
9 a=(4*r)/sqrt(3);
10 efficiency=(%pi*sqrt(3))/8;
11 disp(efficiency, "efficiency = ")
12 disp(a, "lattice parameter, a(angstrom) = ")
```

Scilab code Exa 3.18 interplanar distance

```
1 //Example 3.18 : interplanar distance
2 clc;
3 clear;
4 close;
5 //given data :
6 h=1;
7 k=1;
8 l=1;
9 //d=a/sqrt(h^2+k^2+l^2)
10 dBYa=1/sqrt(h^2+k^2+l^2);
11 disp("Interplanor distance(Angstrom) is a*"+string(dBYa));
```

Scilab code Exa 3.19 interplanar spacing

```
1 //Example 3.19 : spacing
2 clc;
3 clear;
4 close;
5 //given data :
6 h1=2;
7 k1=0;
8 11=0;
9 h2=2;
```

```
10 k2=2;

11 12=0;

12 h3=1;

13 k3=1;

14 13=1;

15 r=1.246;

16 a=(4*r)/sqrt(2);// in angstrum

17 //d=a/sqrt(h^2+k^2+l^2)

18 d1=a/sqrt(h1^2+k1^2+l1^2);

19 d2=a/sqrt(h2^2+k2^2+l2^2);

20 d3=a/sqrt(h3^2+k3^2+l3^2);

21 disp(d1,"d_200 spacind, d1(angstrom) = ")

22 disp(d2,"d_220 spacind, d2(angstrom) = ")

23 disp(d3,"d_111 spacind, d3(angstrom) = ")
```

Scilab code Exa 3.20 interplanar spacing

```
1 //Example 3.20 : interplaner spacing d_220
2 clc;
3 clear;
4 close;
5 format('v',6)
6 //given data :
7 a=0.316;// in nm
8 h=2;
9 k=2;
10 l=0;
11 d=a/sqrt(h^2+k^2+1^2);
12 disp(d,"inter planer spacing d_220,d(nm) = ")
13 // answer is wrong in book
```

Scilab code Exa 3.21 ratio of cubic lattice sepration between the successive lattice planes

```
1 // Example 3.21: interplanar spacing d220
2 clc;
3 clear;
4 close;
5 format('v',5)
6 a=1; //constant assume
7 a1=[1;0;0];//lattice planes
8 a2=[1;1;0];//lattice planes
9 a3=[1;1;1];//lattice planes
10 d100=a/(sqrt(a1(1,1)^2+a1(2,1)^2+a1(3,1)^2));//
     interplanar distance between (100) planes
11 d110=a/(sqrt(a2(1,1)^2+a2(2,1)^2+a2(3,1)^2));
     interplanar distance between (110) planes
12 d111=a/(sqrt(a3(1,1)^2+a3(2,1)^2+a3(3,1)^2));//
     interplanar distance between (111) planes
13 disp("ratio of interplanar distances is "+string(
     d100)+":"+string(d110)+":"+string(d111)+"")
```

Scilab code Exa 3.22 perpendicular distance

```
1 // Example 3.22: perpendicular distance
2 clc;
3 clear;
4 close;
5 a=1;//constant assume
6 a1=[1;1;1];//lattice planes
7 a2=[2;2;2];//lattice planes
8 d1=a/(sqrt(a1(1,1)^2+a1(2,1)^2+a1(3,1)^2));//
        perpendicular distance between origin and (111)
        planes
9 d2=a/(sqrt(a2(1,1)^2+a2(2,1)^2+a2(3,1)^2));//
        perpendicular distance between origin and (222)
        planes
10 d22 = d1-d2;//perpendicular distance between the
        planes (111) and (222)
```

11 disp(d22, "perpendicular distance between the planes (111) and (222)")

Scilab code Exa 3.23 angle

```
1 // Example 3.23: angle between planes (122) and
                         (111)
   2 \text{ clc};
   3 clear;
  4 close;
   5 a=1;// assume
   6 a1=[1;2;2];//lattice planes
   7 a2=[1;1;1];//lattice planes
   8 d1=a/(sqrt(a1(1,1)^2+a1(2,1)^2+a1(3,1)^2));//
                         perpendicular distance between origin and (111)
                         planes
   9 d2=a/(sqrt(a2(1,1)^2+a2(2,1)^2+a2(3,1)^2));/
                         perpendicular distance between origin and (222)
                         planes
10 cphi= ((a1(1,1)*a2(1,1))+(a1(2,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1)*a2(2,1)+(a1(3,1)*a2(2,1))+(a1(3,1)*a2(2,1)*a2(2,1)+(a1(3,1)*a2(2,1)+(a1(3,1)*a2(2,1)+(a1(3,1)*a2(2,1)+(a1(3,1)*a2(2,1)+(a1(3,1)*a2(2,1)+(a1(3,1)*a2(2,1)+(a1(3,1)*a2(2,1)+(a1(3,1)*a2(2,1)+(a1(3,1)*a2(2,1)+(a1(3,1)*a2(2,1)+(a1(3,1)*a2(2,1)+(a1(3,1)*a2(2,1)+(a1(3,1)*a2(2,1)+(a1(3,1)*a2(2,1)+(a1(3,1)*a2(2,1)+(a1(3,1)*a2(2,1)+(a1(3,1)*a2(2,1)+(a1(3,1)*a2(2,1)+(a1(3,1)*a2(2,1)+(a1(3,1)*a2(2,1)+(a1(3,1)*a2(2,1)+(a1(3,1)*a2(2,1)*a2(2,1)+(a1(3,1)*a2(2,1)+(a1(3,1)*a2(2,1)+(a1(3,1)*a2(2,1)+(a1(3,1)*a2(2,1)+(a1(
                        a2(3,1)))*(d1*d2);//
11 d=acosd(cphi);// in degree
12 d1=floor(d);//
13 d2=d-d1; //
14 disp("angle between planes (122) and (111) is "+
                         string(d1)+" degree "+string(round(60*d2))+"
                        minutes")
```

Scilab code Exa 3.24 concentration of iron atoms

```
1 //Example 3.24 : concentration of iron
2 clc;
3 clear;
```

```
4 close;
5 format('v',9)
6 //given data:
7 d=7.87;
8 N=6.023*10^23; // avogadro's number
9 A=55.85; // atomic weight
10 I=A/N; // mass of iron atom
11 atom=d/I;
12 disp(atom,"number of atoms(atoms/cm^3) = ")
```

Scilab code Exa 3.25 lattice constants

```
1 //Example 3.25 : lattice constant
2 clc;
3 clear;
4 close;
5 //given data :
6 format('v',6)
7 n=2;
8 A=55.8;
9 N=6.023*10^26; // avogadro's number in /kg-mole
10 b=7.87*10^3; // in kg/m^3
11 a=((A*n)/(N*b))^(1/3);
12 disp(a*10^10," lattice constant, a(angstrom)")
```

Scilab code Exa 3.26 density

```
1 //Example 3.26 : density
2 clc;
3 clear;
4 close;
5 //given data :
6 format('v',5)
```

```
7 n=4;
8 N=6.023*10^23; // avogadro's number
9 r=1.278*10^-8; // in cm
10 A=63.5;
11 a=(r*4)/sqrt(2); // in cm
12 b=(A*n)/(a^3*N);
13 disp(b,"density of copper, b(g/cc) = ")
```

Scilab code Exa 3.27 number of atoms

```
//Example 3.27 : number of atoms
clc;
clc;
clear;
close;
//given data :
n=4;
N=6.023*10^23; // avogadro's number
A=55.85;
a=2.9*10^-8;
b=7.87; // density in g/cc
//a^3=(A*n)/(N*b)
n=round((a^3*N*b)/A);
disp(n,"number of atoms, n = ")
```

Scilab code Exa 3.28 lattice constants

```
1 //Example 3.28 : lattice constant
2 clc;
3 clear;
4 close;
5 //given data :
6 d=6250;//density
7 N=6.02*10^23;//avogadro's number
```

```
8     n=4;
9     m=60.2*10^-3; // atomic mass
10     M=(n*m)/N;
11     V=M/d;
12     a=V^(1/3)*10^9;
13     disp(a,"the lattice constant, a(nm) = ")
14     //ANSWER IS WRONG IN THE TEXT BOOK
```

Scilab code Exa 3.29 number of atoms

```
1 //Example 3.29 : the number of atoms
2 clc;
3 clear;
4 close;
5 //given data :
6 d=7.87; //in g/cm^3
7 A=55.85;
8 a=2.9*10^-8; // in cm
9 N=6.02*10^23; //avogadro's number
10 n=(d*a^3*N)/A;
11 disp(round(n), "the number of atom, n = ")
```

Scilab code Exa 3.30 number of vacancies in copper

```
// Example 3.30: calculate the number of vacancies
in the copper
clc;
clear;
close;
B=1.38*10^-23;//boltzman constant in J/atom-K
B1=8.62*10^-5;// bolzman constant in ev/atom-K
Qv=0.9;// eV/atom
t=27;// room temperatyre in degree celsius
```

```
9 pcu=8.4; //in g/cm^3
10 Acv=63.5; // in g/mol
11 T=t+273; //temperture in kelvin
12 Nv=6.023*10^23; //
13 P=8.4; //
14 Ns=(Nv*P)/Acv; // number of regular lattice sites
15 Nv1=Ns*exp(-Qv/(B1*T)); //
16 disp(Nv1, "number of vacancies in copper in vacancies /cm^3")
17 //answer is wrong in the textbook
```

Scilab code Exa 3.31 interplanar spacing

```
//Example 3.31 : interplanar spacing
clc;
clcar;
close;
//given data :
format('v',5)
theta=20.3;//in degree
lamda=1.54;// in angstrum
n=1;
a=sind(theta)
d=lamda/(2*a);
disp(d,"interplanar spacing,d(angstrom) = ")
```

Scilab code Exa 3.32 interatomic spacing

```
1 //Example 3.32 : interatomic spacing
2 clc;
3 clear;
4 close;
5 //given data :
```

```
6 format('v',9)
7 theta=30;//in degree
8 lamda=1.54;// in angstrum
9 n=1;
10 a=sind(theta)
11 d=lamda/(2*a);
12 disp(d,"interatomic spacing,d(angstrom) = ")
```

Scilab code Exa 3.33 order of Braggs reflection

```
1 //Example 3.33 : number of per order
2 clc;
3 clear;
4 close;
5 //given data :
6 format('v',5)
7 theta=90;//in degree
8 lamda=1.54;// in angstrum
9 a=sind(theta)
10 d=1.181;
11 n=(2*d*a)/lamda;
12 disp(n,"number of order, n = ")
```

Scilab code Exa 3.34 size of unit cell

```
// Example 3.34: size of unit cell
clc;
clear;
close;
n=1;//
a=1;//assume
h=0.58;//wavelnegth in armstrong
th=9.5;//reflection angle in degree
```

```
9 a1=[2;0;0]; // miller indices
10 d200=a/(sqrt(a1(1,1)^2+a1(2,1)^2+a1(3,1)^2)); //
        interplanar distance between (200) planes
11 a=((n*h)/(2*d200*sind(th))); // zsize of unit cell
12 disp(a,"size of unit cell in ")
13 // amswer is wrong in the textbook
```

Scilab code Exa 3.35 Bragg angle

```
1 // Example 3.35: bragg angle
2 clc;
3 clear;
4 close;
5 n=1; //
6 a=3.57; //in
7 h=0.54; //wavelnegth in
8 a1=[1;1;1];//miller indices
9 d111=a/(sqrt(a1(1,1)^2+a1(2,1)^2+a1(3,1)^2));//
      interplanar distance between (111) planes
10 \operatorname{snd} = ((n*h)/(2*d111)); //
11 th=asind(snd);// bragg angle in degree
12 d1=floor(th);//
13 d2=th-floor(d1);//
14 disp("angle between planes (122) and (111) is "+
      string(d1)+" degree "+string(round(60*d2))+"
      minutes")
15 //wavelength is given wrong in example it is 0.54
      and it is taken as 1.54
```

Scilab code Exa 3.36 interplanar spacing and Miller Indices

```
1 // Example 3.36: interplanner spacing and miller indices
```

```
2 clc;
3 clear;
4 close;
5 a=3.16;// in
6 h=1.54;// in
7 n=1;//
8 th=20.3;// in degree
9 d=((n*h)/(2*sind(th)));// interplanner spacing in
10 x=a/d;//
11 y=x^2;//
12 disp(d,"interplanner spacing in is")
13 disp("miller indices are (110), (011) or (101)")
```

Scilab code Exa 3.37 interplanar spacing and diffraction angke

```
1 // Example 3.36: interplanner spacing and
      diffraction angle
2 clc;
3 clear;
4 close;
5 a = .2866; // in
6 h=0.1542; // in nm
7 n=1; //
8 a1=[2;1;1];//miller indices
9 d211=a/(sqrt(a1(1,1)^2+a1(2,1)^2+a1(3,1)^2));//
     interplanar distance between (211) planes
10 snd=((n*h)/(2*d211));//
11 th=asind(snd);// bragg angle in degree
12 d1=floor(th);//
13 d2=th-floor(d1);//
14 disp("angle between planes (122) and (111) is "+
     string(d1)+" degree "+string(round(60*d2))+"
     minutes")
15 disp(d211,"interplanner spacing in
                                          is")
16 //answer is wrong in the textbook
```

Chapter 5

Electron Theory of Metals

Scilab code Exa 5.1.a probability

```
1 // Example 5.1.a: probability for diamond
2 clc, clear
3 // given :
4 format('v',9)
5 Eg=5.6; // in eV
6 k=86.2*10^-6; // in eVk^-1
7 T=273+25; // in K
8 E_Ef=Eg/2;
9 f_E=1/(1+exp(E_Ef/(k*T)));
10 disp(f_E," probability for diamond, f_E = ")
```

Scilab code Exa 5.1.b probability

```
1 // Example 5.1.b: probability for silicon
2 clc, clear
3 // given :
4 Eg=1.07; // in eV
5 k=86.2*10^-6; // in eVk^-1
```

```
6 T=273+25; // in K
7 E_Ef=Eg/2;
8 f_E=1/(1+exp(E_Ef/(k*T)));
9 disp(f_E, "probability for diamond, f_E = ")
10 // answer is wrong in book
```

Scilab code Exa 5.2 resistance

```
1 // Example 5.2: resistance
2 clc, clear
3 // given :
4 l=1; // length in m
5 A=4*10^-4; // area of cross section in m^2
6 p=0.01*10^-2; // resistivity in ohm—m
7 R=p*(1/A);
8 disp(R,"resistance of wire,R(ohm) = ")
```

Scilab code Exa 5.3 resistance

```
1 // Example 5.3: resistance
2 clc, clear
3 // given :
4 format('v',5)
5 p=1.7*10^-8; // resistivity i ohm—m
6 d=0.0005; // diameter of the wire in m
7 l=31.4; // length in m
8 A=(%pi*d^2)/4;
9 R=p*(1/A);
10 disp(R,"resistance of wire,R(ohm) = ")
```

Scilab code Exa 5.4 conductivity

```
// Example 5.4: conductivity
clc, clear
// given :
format('v',8)
V=.432; // voltage drop across the wire in volts
I=10; // current through the wire in A
I=1; // length in m
d=1*10^-3; // diameter in m
R=V/I;
A=(%pi*d^2)/4;
p=(R*A)/1;
b=1/p;
disp(b,"conductivitty,b(ohm^-1.m^-1) = ")
```

Scilab code Exa 5.5 drift velocity

```
1 // Example 5.5: drift velocity
2 clc, clear
3 // given :
4 format('v',5)
5 n=10^19; // in m^3
6 b=0.01; // conductivity in ohm^-1. m^-1
7 V=0.17; // in volts
8 d=.27*10^-3; // in m
9 e=1.602*10^-19; // in C
10 m=9.1*10^-31; // in kg
11 E=V/d; // in volt/m
12 v=((b*E)/(n*e));
13 disp(v,"drift velocity of electron, v (m/sec) = ")
```

Scilab code Exa 5.6 conductivity

```
1 // Example 5.6: conductivity
2 clc, clear
3 // given :
4 e=1.6*10^-19; // in C
5 T=300; // temerature in K
6 t=2*10^-14; // time in sec
7 c=63.54; // atomic weight of copper in a.m.u
8 \text{ m=9.1*10}^{-31}; // mass in kg
9 // we know that 63.45 grams of copper contains
     6.023*10^23 free electrons since one atom
     contributes one electron the volume of 63.54 gram
      of copper is 8.9 cubic centimetre(c.c).
10 n=6.023*10^23/(c/8.9); //number of electrons per
      unit volume (c.c)
11 n1=n*10^6; // the number of electrons per m^3
12 b=(e^2*n1*t)/m;
13 disp(b, "conductivity, b(mho/m) = ")
```

Scilab code Exa 5.7 mobility of electrons

Scilab code Exa 5.8 mobility of electrons

```
1 // Example 5.8: mobility
2 clc, clear
3 // given :
4 format('v',6)
5 d=10.5; // density of silver in gm/c.c
6 w=107.9; // atomic weight
7 b=6.8*10^5; // conductivity in mhos/cm
8 e=1.602*10^-19; // in C
9 N=6.023*10^23;
10 n=(N*d)/w;
11 mu=b/(e*n);
12 disp(mu," mobility of electron ,mu(m^2/volt-sec) = ")
```

Scilab code Exa 5.9 mobility of electrons and drift velocity

```
1 // Example 5.9: mobility and drift velocity
2 clc, clear;
3 // given :
4 b=6.5*10^7; // conductivity in ohm^-1.m^-1
5 e=1.602*10^-19; // in C
6 n=6*10^23; //
7 E=1; // in V/m
8 mu=b/(e*n);
9 v=mu*E;
10 disp(mu," mobility ,mu(m^2/volt-sec) = ")
11 disp(v," drift velocity ,v(m/sec) = ")
12 // mobility and drift is calculated wrong in book
```

Scilab code Exa 5.10 density and drift velocity

```
1 //Example 5.10 : density and drift velocity
2 clc;
3 clear;
```

```
4 close;
5 //given data:
6 format('v',9)
7 e=1.602 *10^-19;
8 b=58*10^6; // in ohm^-1 m^-1
9 mu_n=3.5*10^-3; // in m^2/V s
10 E=0.5; // in V/m
11 n=b/(e*mu_n);
12 disp(n,"density,n(m^-3) = ")
13 v=mu_n*E;
14 disp(v,"drift velocity,v(m/s) = ")
```

Scilab code Exa 5.11 velocity

```
1 //Example 5.11 : velocity
2 clc;
3 clear;
4 close;
5 //given data :
6 m=9.109*10^-31; // in kg
7 e=1.602 *10^-19;
8 Ef=2.1// in ev
9 Wf=e*Ef;// in J
10 vf=sqrt((2*Wf)/m);
11 disp(vf," velocity, vf(m/s) = ")
```

Scilab code Exa 5.12.a velocity

```
1 //Example 5.12.a : velocity
2 clc;
3 clear;
4 close;
5 //given data :
```

```
6 m=9.1*10^-31; // in kg
7 e=1.602 *10^-19;
8 Ef=3.75; // in ev
9 Wf=(e*Ef); // in J
10 vf=sqrt(((2*Wf)/m));
11 disp(vf,"velocity, vf(m/s) = ")
12 // answer is wrong in book
```

Scilab code Exa 5.12.b mobility of electrons

```
1
2  //Example 5.12.b : mobility of electron
3  clc;
4  clear;
5  close;
6  //given data :
7  m=9.1*10^-31; // in kg
8  e=1.602 *10^-19;
9  Ef=3.75; // in ev
10  t=10^-14; // in sec
11  mu=(e*t)/m;
12  disp(mu," mobility ,mu(m^2/V-sec) = ")
```

Scilab code Exa 5.13 mean free path

```
1 //Example 5.13 : the mean free path
2 clc;
3 clear;
4 close;
5 //given data :
6 format('v',5)
7 t=10^-9; // in sec
8 m=9.109*10^-31; // in kg
```

```
9 e=1.602 *10^-19;
10 Ef=7// in ev
11 Wf=e*Ef;// in J
12 vf=sqrt((2*Wf)/m);
13 lamda=vf*t*10^3;
14 disp(lamda,"the mean free path, lamda(mm) = ")
```

Scilab code Exa 5.14 mobility and average time

```
1 //Example 5.14 : mobility and average time
2 clc;
3 clear;
4 close;
5 //given data:
6 format('v',6)
7 \text{ m=9.109*10}^{-31}; // in kg
8 e=1.602 *10^-19;
9 d=8.92*10^3;// in kg/m^3
10 p=1.73*10^-8; // ohm-m
11 A=63.5; //atomic weight
12 N=6.023*10^22; // avogadro's number
13 n=(N*d)/A;
14 b=1/p;// conductivity
15 mu=b/(n*e);
16 disp(mu, "mobility, mu(m^2/V-s) = ")
17 t=(mu*m)/e;
18 disp(t*10^9, "average time, t(ns) = ")
```

Scilab code Exa 5.15 electrical resistivity

```
1 //Example 5.15 : electrical resistivity
2 clc;
3 clear;
```

```
4 close;
5 //given data :
6 format('v',8)
7 r=1.86*10^-10; // in m
8 t=3*10^-14; // in sec
9 a=2;
10 m=9.1*10^-31; // in kg
11 e=1.602 *10^-9;
12 A = 23 * 10^{-3}; // in kg/m
13 N=6.023*10^23; // avogadro's number
14 M = (a*A)/N;
15 V=((4/sqrt(3))*r)^3;
16 \quad d=M/V;
17 mu = ((e*t)/m);
18 n = (N*d)/A;
19 b=1.602 *10^-19*n*mu;
20 p = (1/b);
21 disp(p, "resistivity, p(ohm-m) = ")
```

Mechanical Tests of Metals

Scilab code Exa 7.1 shear modulus

```
//Example 7.1 : shear modulus of the material
clc;
clear;
close;
//given data :
format('v',6)
E=210; // youngs's modulus in GN/m^2
v=0.3; // poisson ratio
G=E/(2*(1+v));
disp(G,"shear modulus,G(GN/m^2) = ")
```

Scilab code Exa 7.2 young modulus of elasticity yield point uttimate stress and percentage elongation

```
1 //Example 7.2 : young's modulus of elasticity, yield
    point stress, ultimate stress and percentage
    elongation
2 clc;
```

```
3 clear;
4 close;
5 format('v',9)
6 //given data :
7 d=40*10^-3; //in m
8 W=40*10^3; // load in N
9 del_l=3.04*10^-5;// in m
10 L=200*10^-3; // in m
11 load_max = 242*10^3; //in N
12 1=249*10^-3; // length of specimen in m
13 10=(d+L); // in m
14 A = (\%pi*d^2)/4;
15 b=W/A;
16 epsilon=del_l/L;
17 E=(b/epsilon);
18 disp(E, "young modulus, E(N/m^2) = ")
19 Y_load=161*10^3;
20 Y_stress=Y_load/A;
21 disp(Y_stress," yield point stress, Y_stress(N/m^2) =
22 U_stress=load_max/A;
23 disp(U_stress,"ultimate stress, U_stress(N/m^2) = ")
24 p_elongation=((1-10)/10)*100;
25 disp(p_elongation, percentage elongation,
      p_{elongation}(\%) = ")
26 //percentage elongation is calculated wrong in
     textbook
```

Scilab code Exa 7.3.a yield point stress

```
1 // Example 7.3.a: yield point stress
2 clc;
3 clear;
4 close;
5 format('v',10)
```

```
6 yl=40; // yeild load in kN
7 ml=71.5; // maximum load in kN
8 fl=50.5; // fracture load in kN
9 glf=79.5; // gauge length of fratture in mm
10 st=7.75*10^-4; // strain at load of 20kN
11 d=12.5; // specimen diamtere in mm
12 sl=62.5; // specimen length in mm
13 A=(%pi*(d*10^-3)^2)/4; // in meter square
14 ylp=((yl*10^3)/(A)); // yeild point stress in N/m^2
15 disp(ylp," yeild point stress in N/m^2")
```

Scilab code Exa 7.3.b ultimate tensile strength

```
1 // Example 7.3.b: ultimate tensile strength
2 clc;
3 clear;
4 close;
5 format('v',10)
6 y1=40; // yeild load in kN
7 ml=71.5; //maximum load in kN
8 fl=50.5; // fracture load in kN
9 glf=79.5;//gauge length of fratture in mm
10 st=7.75*10^-4; // strain at load of 20kN
11 d=12.5; //specimen diamtere in mm
12 sl=62.5; //specimen length in mm
13 A = (\%pi*(d*10^-3)^2)/4; // in meter square
14 ylp=((yl*10^3)/(A));//yeild point stress in N/m^2
15 uts=((ml*10^3)/(A));//ultimate tensile strangth in N
      /\mathrm{m}^2
16 disp(uts," ultimate tensile strangth in N/m<sup>2</sup>")
```

Scilab code Exa 7.3.c percentage elongation

```
// Example 7.3.c: percentage elongation
clc;
clear;
close;
format('v',10)
fyl=40;//yeild load in kN
ml=71.5;//maximum load in kN
fl=50.5;//fracture load in kN
glf=79.5;//gauge length of fratture in mm
st=7.75*10^-4;//strain at load of 20kN
d=12.5;//specimen diamtere in mm
sl=62.5;//specimen length in mm
a=(%pi*d*10^-3)^2/4;// in meter square
pel=((glf-sl)/sl)*100;//percentage elongation
disp(pel,"percentage elongation is")
```

Scilab code Exa 7.3.d modulus of elasticity

```
1 // Example 7.3.d:modulus of elasticity
2 clc;
3 clear;
4 close;
5 format('v',8)
6 y1=40; // yeild load in kN
7 ml=71.5; //maximum load in kN
8 fl=50.5; //fracture load in kN
9 glf=79.5; //gauge length of fratture in mm
10 st=7.75*10^-4; //strain at load of 20kN
11 d=12.5; //specimen diamtere in mm
12 sl=62.5; //specimen length in mm
13 A = (\%pi*(d*10^-3)^2)/4; // in meter square
14 ylp=((yl*10^3)/(A));//yeild point stress in N/m^2
15 uts=((ml*10^3)/(A));//ultimate tensile strangth in N
16 pel=((glf-sl)/sl)*100;//percentage elongation
```

```
17 strss=((20*10^3)/A);//stress at 20kN in N/m^2 18 mel=strss/st;//modulus of elasticity in N/m^2 19 disp(mel,"modulus of elasticity in N/m^2")
```

Scilab code Exa 7.3.e modulus of resilience

```
1 // Example 7.3.e: yield point stress
2 clc;
3 clear;
4 close;
5 format('v',6)
6 y1=40; // yeild load in kN
7 ml=71.5; //maximum load in kN
8 fl=50.5; //fracture load in kN
9 glf=79.5; //gauge length of fratture in mm
10 st=7.75*10^-4; //strain at load of 20kN
11 d=12.5; //specimen diamtere in mm
12 sl=62.5; //specimen length in mm
13 A = (\%pi*(d*10^-3)^2)/4; // in meter square
14 ylp=((yl*10^3)/(A)); // yeild point stress in N/m^2
15 uts=((ml*10^3)/(A));//ultimate tensile strangth in N
      /\mathrm{m}^2
16 pel=((glf-sl)/sl)*100;//percentage elongation
17 strss=((20*10^3)/A); //stress at 20kN in N/m^2
18 mel=strss/st;//modulus of elasticity in N/m<sup>2</sup>
19 mrs=((ylp*10^-3)^2/(2*mel)); //modulus of resilience
20 disp(mrs, "modulus of resilience is")
```

Scilab code Exa 7.3.f fracture stress

```
1 // Example 7.3.f: fracture stress
2 clc;
3 clear;
```

```
4 close;
5 format('v',10)
6 y1=40; // yeild load in kN
7 ml=71.5; //maximum load in kN
8 fl=50.5; //fracture load in kN
9 glf=79.5; //gauge length of fratture in mm
10 st=7.75*10^-4; //strain at load of 20kN
11 d=12.5; //specimen diamtere in mm
12 sl=62.5; //specimen length in mm
13 A = (\%pi*(d*10^-3)^2)/4; // in meter square
14 ylp=((yl*10^3)/(A));//yeild point stress in N/m^2
15 uts=((ml*10^3)/(A));//ultimate tensile strangth in N
      /\mathrm{m}^2
16 pel=((glf-sl)/sl)*100;//percentage elongation
17 strss=((20*10^3)/A); //stress at 20kN in N/m^2
18 mel=strss/st;//modulus of elasticity in N/m<sup>2</sup>
19 mrs=((ylp*10^-3)^2/(2*mel)); //modulus of resilience
20 fs=((f1*10^3)/(A));//fracture stress in N/m^2
21 disp(fs, "fracture stress in N/m^2")
```

Scilab code Exa 7.3.g modulus of toughness

```
1
2 // Example 7.3.g: modulus of toughness
3 clc;
4 clear;
5 close;
6 format('v',10)
7 yl=40;//yeild load in kN
8 ml=71.5;//maximum load in kN
9 fl=50.5;//fracture load in kN
10 glf=79.5;//gauge length of fratture in mm
11 st=7.75*10^-4;//strain at load of 20kN
12 d=12.5;//specimen diamtere in mm
13 sl=62.5;//specimen length in mm
```

Scilab code Exa 7.4 true breaking and nominal breaking stress

```
2 //Example 7.4 : true breaking stress and nominal
     breaking stress
3 clc;
4 clear;
5 close;
6 format('v',4)
7 //given data :
8 d1=12.7; // in mm
9 B_load=14; // in K-N
10 A1=(\%pi*d1^2)/4;// original cross section area
11 d2=7.87; // in mm
12 A2=(\%pi*d2^2)/4; // final cross sction area
13 T_stress=B_load/A2;
14 disp(T_stress*1000,"true breaking stress, T_stress(N/
     mm^2 = "
15 N_stress=B_load/A1;
16 disp(N_stress*1000, "nominal breaking stress, N_stress
```

```
\rm (N/mm^2) = ") 17 //true breaking stress unit is wrong in the textbook
```

Scilab code Exa 7.5.a yield stress

```
// Example 7.5.a: yield point stress
clc;
clear;
close;
format('v',10)
yl=34;//yeild load in kN
ul=61;//ultimate load in kN
fl=78;//final length in mm
glf=60;//gauge length of fratture in mm
fd=7;//final diamtere in mm
cl=12;//specimen diamtere in mm
sl=62.5;//specimen length in mm
A=(%pi*(d)^2)/4;// in meter square
ylp=((yl*10^3)/(A));//yeild point stress in N/mm^2
disp(floor(ylp), "yeild point stress in N/mm^2")
```

Scilab code Exa 7.5.b ultimate tesnile stress

```
1 // Example 7.5.b: ultimate tensile stress
2 clc;
3 clear;
4 close;
5 format('v',6)
6 yl=34;//yeild load in kN
7 ul=61;//ultimate load in kN
8 fl=78;//final length in mm
9 glf=60;//gauge length of fratture in mm
10 fd=7;//final diamtere in mm
```

```
11 d=12; //specimen diamtere in mm
12 sl=62.5; //specimen length in mm
13 A=(%pi*(d)^2)/4; // in meter square
14 uts=((ul*10^3)/(A)); // ultimate tensile strangth in N /mm^2
15 disp(uts," ultimate tensile strangth in N/mm^2")
```

Scilab code Exa 7.5.c percentage reduction

```
1 // Example 7.5.c: percentage reduction
2 clc;
3 clear;
4 close;
5 format('v',4)
6 y1=34; // yeild load in kN
7 ul=61; // ultimate load in kN
8 fl=78; // final length in mm
9 glf=60; //gauge length of fratture in mm
10 fd=7; // final diamtere in mm
11 d=12; //specimen diamtere in mm
12 sl=62.5; //specimen length in mm
13 A = (\%pi*(d)^2)/4; // in mm square
14 A1=(\%pi*(fd)^2)/4;// in mm square
15 pr = (A - A1)/A; //
16 disp(pr*100, "percentage reduction is")
```

Scilab code Exa 7.5.d percentage elongation

```
1 // Example 7.5.d: percentage elonagtion
2 clc;
3 clear;
4 close;
5 format('v',4)
```

```
6 yl=34; // yeild load in kN
7 ul=61; // ultimate load in kN
8 fl=78; // final length in mm
9 glf=60; // gauge length of fratture in mm
10 fd=7; // final diamtere in mm
11 d=12; // specimen diamtere in mm
12 sl=62.5; // specimen length in mm
13 A=(%pi*(d)^2)/4; // in mm square
14 A1=(%pi*(fd)^2)/4; // in mm square
15 pr=(fl-glf)/glf; //
16 disp(pr*100, "percentage elonagtion is")
```

Scilab code Exa 7.6 strain

```
1 //Example 7.6 : strain
2 clc;
3 clear;
4 close;
5 format('v',10)
6 //given data :
7 b=44.5*10^3;//force
8 E=1.1*10^5;// in N/mm^2
9 A=15.2*19.1// in mm^2
10 epsilon=b/(A*E);
11 disp(epsilon, "strain, epsilon (mm) = ")
```

Scilab code Exa 7.7 true stress and true strain

```
1 //Example 7.7 :stress and strain
2 clc;
3 clear;
4 close;
5 format('v',6)
```

```
6 //given data :
7 sigma=450; //in MPa
8 epsilon=0.63;
9 sigma_t=sigma*(1+epsilon);
10 disp(sigma_t,"true stress, sigma_t (MPa) = ")
11 epsilon_t=log(1+epsilon);
12 disp(epsilon_t,"true strain, epsilon_t = ")
```

Scilab code Exa 7.8 greater stress

```
1 // Example 7.8: which part has a greater stress
2 clc;
3 clear;
4 close;
5 l=24;//length in mm
6 b=30;//breadth in mm
7 ld=7000;//load in kg
8 sd=10;//steel bar diamtere in mm
9 sl=5000;//load in kg
10 al=ld/(l*b);//stress on aluminium bar in kg/mm^2
11 a=((%pi*sd^2)/4);//area in mm^2
12 slb=sl/a;//stress on steel bar in kg/mm^2
13 disp("stress on aluminium bar is "+string(al)+" kg/mm^2 is less than stress on steel bar "+string(slb)+" kg/mm^2")
```

Mechanical Tests of Metals

Scilab code Exa 8.1 critical resolved shear stress

```
1 // Example 8.1: critical resolved shear stress of
      silver
2 clc;
3 clear;
4 close;
5 format('v',5)
6 Ts=15; // tensile stress in Mpa
7 d=[1;1;0];
8 d1=[1;1;1];
9 csda=((d(1,1)*d1(1,1))+(d(2,1)*d1(2,1))+(d(3,1)*d1
     (3,1))/((sqrt(d(1,1)^2+d(2,1)^2+d(3,1)^2))*sqrt(
     d1(1,1)^2+d1(2,1)^2+d1(3,1)^2));//angle degree
10 d2 = [0;1;1];
11 csdb = ((d(1,1)*d2(1,1))+(d(2,1)*d2(2,1))+(d(3,1)*d2(2,1))
     (3,1))/((sqrt(d(1,1)^2+d(2,1)^2+d(3,1)^2))*sqrt(
     d2(1,1)^2+d2(2,1)^2+d2(3,1)^2);/angle degree
12 t=Ts*csda*csdb;//critical resolved shear stress in
13 disp(t, "critical resolved shear stress in MPa")
```

Scilab code Exa 8.2 yield strength

```
1 // Example 8.2: yield strength of material
2 clc;
3 clear;
4 close;
5 format('v',6)
6 ys1=115; // yeild strength in MN/mm^2
7 ys2=215;// yeild strength in MN/mm^2
8 d1=0.04; //diamtere in mm
9 d2=0.01; // diamtere in mm
10 A = [2 \ 10; \ 1 \ 10];
11 B=[230;215];
12 x = A \setminus B;
13 si=x(1,1);// in MN/mm^2
14 k=x(2,1);//
15 d3=0.016; //in mm
16 sy = si +(k/sqrt(d3)); //yeild strength for a grain
      size in MN/mm<sup>2</sup>
17 disp(sy," yeild strength for a grain size in MN/mm<sup>2</sup>"
      )
```

Scilab code Exa 8.3 yield stress

```
// Example 8.3: yield strength of material
clc;
clear;
close;
ys1=120;// yeild strength in MN/mm^2
ys2=220;// yeild strength in MN/mm^2
d1=0.04;//diamtere in mm
d2=0.01;//diamtere in mm
```

```
9 A=[2 10; 1 10];
10 B=[240;220];
11 x=A\B;
12 si=x(1,1);// in MN/mm^2
13 k=x(2,1);//
14 d3=0.025;//in mm
15 sy= si +(k/sqrt(d3));//yeild strength for a grain size in MN/mm^2
16 disp(sy,"yeild strength for a grain size in MN/mm^2")
```

Scilab code Exa 8.4 grain diameter

```
1 //Example 8.4 : grain diameter
2 clc;
3 clear;
4 close;
5 format('v',6)
6 //given data :
7 N=9;
8 m=8*2^N;
9 grain=1/sqrt(m);
10 disp(grain,"the grain diameter(mm) = ")
```

Fracture of Metals

Scilab code Exa 9.1 fracture strength

```
1 //Example 9.1 : difference
2 clc;
3 clear;
4 close;
5 //given data :
6 E=200*10^9; // in N/m^2
7 C=(4*10^-6)/2; // in m
8 gama=1.48; // in J/m^2
9 sigma=sqrt((2*E*gama)/(%pi*C));
10 disp(sigma*10^-6, "fracture strength, sigma(MN/m^2) = ")
```

Scilab code Exa 9.2 fracture strength

```
1 //Example 9.2 : the fracture strength and compare
2 clc;
3 clear;
4 close;
```

```
5 format('v',10)
6 //given data:
7 E=70*10^9; // in N/m^2
8 C=(4.2*10^-6)/2; // in m
9 gama=1.1; // in J/m^2
10 sigma=sqrt((2*E*gama)/(%pi*C));
11 disp(sigma, "fracture strength, sigma(N/m^2) = ")
```

Scilab code Exa 9.3 maximum length

```
//Example 9.3 : maximum length of surfacef
clc;
clear;
close;
format('v',7)
//given data :
sigma=36;//in MN/m^2
gama=0.27;// in J/m^2
C=((2*E*gama)/(sigma^2*%pi))*10^-6;
C2=2*C;
disp(C2,"maximum length of surface flow,C2(micro-m) = ")
```

Scilab code Exa 9.4.a temperture

```
1 // Example 9.4.a: Temperature
2 clc;
3 clear;
4 close;
5 format('v',6)
6 E=350; // in GN/m<sup>2</sup>
7 Y=2; // in J/m<sup>2</sup>
```

```
8 C=2; // in micro meter
9 sg=sqrt((2*E*10^9*Y)/(%pi*C*10^-6)); // IN mn/M^2
10 e=10^-2; // per second
11 T=173600/(round(sg*10^-6)-20.6-61.3*(log10(e))); //
        in kelvin
12 disp(T,"temperture in kelvin for ductile to brittle
        transition at a strain rate of 10^-2 per second")
```

Scilab code Exa 9.4.b temperature

```
1 // Example 9.4.b: Temperature
2 clc;
3 clear;
4 close;
5 format('v',5)
6 E=350; // in GN/m^2
7 Y=2; // in J/m^2
8 C=2; // in micro meter
9 sg=sqrt((2*E*10^9*Y)/(%pi*C*10^-6)); // IN mn/M^2
10 e=10^-5; // per second
11 T=173600/(round(sg*10^-6)-20.6-61.3*(log10(e))); // in kelvin
12 disp(T,"temperture in kelvin for ductile to brittle transition at a strain rate of 10^-5 per second")
```

Composite Materials and Ceramics

Scilab code Exa 15.1 volume ratio of aluminium and boron

```
//Example 15.1 : colume ratio of aluminium and boron
clc;
clear;
close;
format('v',6)
yal=715;// in GN/,^2
yfe=210;// in GN/,^2
yb=440;// in GN/,^2
B=[71 71;71 440];//
B=[71;210];//
X=A\B;//
disp(X(1,1),"volume ratio of aluminium is")
disp(X(2,1),"volume ratio of boron is")
```

Semiconductors

Scilab code Exa 16.1 concentration of conductive electrons

```
1 //Example 16.1 : concentration
2 clc;
3 clear;
4 close;
5 format('v',9)
6 //given data :
7 e=1.602*10^-19;
8 sigma_i=5*10^-4; // in ohm/m
9 mu_n=0.14; // in m^2/V-sec
10 mu_p=0.05; // in m^2/V-sec
11 n_i=sigma_i/(e*(mu_n+mu_p));
12 disp(n_i*10^6, "the concentration, n_i(/cm^3) = ")
```

Scilab code Exa 16.2 intrinsic carrier density

```
1 //Example 16.2 : intrinsic carrier
2 clc;
3 clear;
```

```
4 close;
5 format('v',9)
6 //given data:
7 e=1.602*10^-19;
8 p_i=2*10^-4;// in ohm-m
9 mu_n=6;// in m^2/V-sec
10 mu_p=0.2;// in m^2/V-sec
11 n_i=1/(e*(mu_n+mu_p)*p_i);
12 disp(n_i,"the intrinsic carrier, n_i(/m^3) = ")
```

Scilab code Exa 16.3 concentration of N type impurity

```
//Example 16.3 : neglect the intrinsic conductivity
clc;
clear;
close;
format('v',9)
//given data :
e=1.6*10^-19;
sigma=10^-12;// in mhos/m
mu_n=0.18;// in m^2/V-sec
n=sigma/(e*mu_n);
N=n;
disp(N,"in(/m^3) = ")
```

Scilab code Exa 16.4 concentration number of electrons carrier

```
//Example 16.4 : number of electron carriers
clc;
clear;
close;
format('v',9)
//given data :
```

```
7 e=1.6*10^-19;
8 p=20*10^-2; // in ohm-m
9 mu_n=100*10^-4; // in m^2/V-sec
10 n=1/(e*mu_n*p);
11 disp(n,"number of electrons carrier, n(/m^3) = ")
```

Scilab code Exa 16.5 concentration of impurity

```
//Example 16.5 : concentration of impurity
clc;
clear;
close;
format('v',9)
e=1.6*10^-19;//
1=10;//in mm
d=1;//in mm
up=0.19;//mobilty of electrons in V-sec
a=(%pi*((d*10^-3)^2))/4;//area in m^2
p=((r*a))/(1*10^-3);//resistivity in Ohm-cm
n=((1/(p*e*up)));//concentration in per m^3
disp(n,"impurity concentration is in per m^3")
```

Scilab code Exa 16.6 intrinsic carrier density

```
1 //Example 16.6 : intrinsic carrier density
2 clc;
3 clear;
4 close;
5 //given data :
6 format('v',10)
7 e=1.602*10^-19;
8 p=3000; // in ohm/m
```

```
9 sigma=1/p;// in ohm/m
10 mu_n=0.14;// in m^2/V-sec
11 mu_p=0.05;// in m^2/V-sec
12 n_i=sigma/(e*(mu_n+mu_p));
13 disp(n_i,"the concentration, n_i(/m^3) = ")
```

Scilab code Exa 16.7 conductivity

```
1 //Example 16.7 : conductivity
2 clc;
3 clear;
4 close;
5 //given data :
6 e=1.602*10^-19;
7 n_i=5.021*10^15; // in m^-3
8 mu_n=0.48; // in m^2/V-sec
9 mu_p=0.013; // in m^2/V-sec
10 sigma=n_i*(e*(mu_n+mu_p));
11 disp(sigma,"the conductivity, sigma(ohm^-1 m^-1) = ")
```

Insulating Materials

Scilab code Exa 17.1 greater charge

```
1 // Example 17.1: greater chanrge
2 clc;
3 clear;
4 close;
5 format('v',10)
6 er1=6;//
7 d1=0.25;// in mm
8 a=1;// assume
9 er2=2.6;//
10 d2=0.1;// in mm
11 c1=(er1/d1);// in ampere
12 c2=(er2/d2);// in amperes
13 disp(" C1 "+string(c1)+"A will hold the more charge than C2 "+string(c2)+"A")
```

Magnetic Materials

Scilab code Exa 18.1 magnetization and flux density

```
//Example 18.1 : magnetization and flux density
clc;
clcar;
close;
//given data :
mu0=4*%pi*10^-7;
H=10^4;// in A/m
Xm=3.7*10^-3;// room temperature
mu_r=1+Xm;
B=mu0*mu_r*H;
M=Xm*H;
disp(B,"the flux density,B(Wb/m^2) = ")
disp(M,"magnetization,M(A/m) = ")
```

Scilab code Exa 18.2.a saturation magnetisation

```
1 //Example 18.2.a : saturation magnetization
2 clc;
```

```
3 clear;
4 close;
5 //given data :
6 mu_b=9.27*10^-24; // A.m^2
7 p=8.9; // in g/cm^3
8 Na=6.023*10^23; // avogadro's number
9 A=58.71; // in g/mol
10 n=((p*Na)/A)*10^6;
11 Ms=0.60*mu_b*n;
12 disp(Ms,"saturation magnetization, Ms(A/m) = ")
```

Scilab code Exa 18.2.b saturation flux density

```
//Example 18.2.b : saturation flux density
clc;
clcar;
close;
//given data :
format('v',5)
mu0=4*%pi*10^-7;
mu_b=9.27*10^-24;// A.m^2
p=8.9; // in g/cm^3
Na=6.023*10^23;// avogadro's number
A=58.71; // in g/mol
n=((p*Na)/A)*10^6;
Ms=0.60*mu_b*n;
Bs=mu0*Ms;
disp(Bs,"saturation flux density, Bs(tesla) = ")
```

Scilab code Exa 18.3 magnetic moments

```
1 //Example 18.3 : magnetic moment
2 clc;
```

Scilab code Exa 18.4 power loss

```
1 //Example 18.4 : power loss
2 clc;
3 clear;
4 close;
5 //given data :
6 V=0.01; //in m^3
7 f=50; // in Hz
8 area=600; //in jm^-1
9 Wh=area*V*f;
10 disp(Wh,"power loss, Wh(watts) = ")
```

Scilab code Exa 18.5 loss of energy

```
1 //Example 18.4 : los of energy 2 clc;
```

```
3 clear;
4 close;
5 //given data :
6 mass=10; // in kg
7 energy_loss=250; // in J/m^2
8 //energy loss at the rate of 50 cycles/s
9 E=energy_loss*50; // in J/m^3
10 E_loss=E*3600; // in J/m^3
11 D=7500; //density in kg/m^3
12 Volume=mass/D;
13 energy_loss_per_hour=E_loss/Volume;
14 disp(energy_loss_per_hour, "energy_loss_per_hour(J/hour) = ")
```