

Scilab Textbook Companion for  
Material Science In Engineering  
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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

# REVIEW OF ATOMIC CONCEPTS ATOMIC MODELS AND PERIODIC TABLE

Scilab code Exa 2.1 Find average atomic weight of Hydrogen

```
1 //Find average atomic weight of Hydrogen.  
2 //Exa:1.1  
3 close;  
4 clc;  
5 clear;  
6 a1=1.0078; //atomic weight of H-1  
7 a2=2.0143; //atomic weight of H-2  
8 p1=99.985; //% of H-1  
9 p2=.015; //% of H-2  
10 a=((a1*p1)+(a2*p2))/100  
11 disp(a,"Average atomic weight of Hydrogen = ");
```

---

### Scilab code Exa 2.2 Find distance

```
1 //Ex:2.2
2 clc;
3 clear;
4 close;
5 z=79; //atomic no. of gold
6 e=7.68*1.6*10^-13; //ke in J
7 e_c=1.6*10^-19; //charge of electron in C
8 e_0=8.854*10^-12; //permittivity F/m
9 d=(2*e_c^2*z)/(4*3.14*e_0*e); //distance in m
10 disp(d,"distance (in m) = ");
```

---

### Scilab code Exa 2.3 Find number of particles scattered at 75 and 135 degree

```
1 //Find number of particles scattered at 75 and 135
  degree
2 //Ex:2.3
3 clc;
4 clear;
5 close;
6 n=44; //no. of particles scattered per minute
7 a=90; //angle in degrees
8 x=sind(a/2);
9 c=n*x^4;
10 disp(c,"Proportionality constant = ");
11 b=75; //angle in degrees
12 y=sind(b/2);
```

```

13 n1=c/y^4;
14 disp(n1,"No. of particles scattered at 75 degree (in
    per minute) = ");
15 d=135;//angle in degrees
16 z=sind(d/2);
17 n2=c/z^4;
18 disp(n2,"No. of particles scattered at 135 degree (
    in per minute) = ");

```

---

Scilab code Exa 2.4 Find radius of first orbit of electron in Hydrogen atom

```

1 //Find radius of first orbit of electron in Hydrogen
    atom
2 //Ex:2.4
3 clc;
4 clear;
5 close;
6 n=1;//first orbit
7 e_0=8.85*10^-12;//permittivity in freee space
8 h=6.62*10^-34;//planck's constant
9 m=9.1*10^-31;//mass of an electron in kg
10 e=1.6*10^-19;//charge of an electron in C
11 z=1;
12 r=n^2*e_0*h^2/(3.14*m*e^2*z);//radius of first orbit
    in m
13 r1=r*10^10;//radius in Angstorm
14 disp(r1,"Radius of first orbit of electron in
    Hydrogen atom (in Angstorm) = ");

```

---

### Scilab code Exa 2.5 Find Ionisation Energy

```
1 //Find Ionisation Energy
2 //Ex:2.5
3 clc;
4 clear;
5 close;
6 e_0=8.85*10^-12;//permittivity in free space in sqC
    /N/sqm
7 h=6.62*10^-34;//planck's constant in Js
8 m=9.1*10^-31;//mass of an electron in kg
9 e=1.6*10^-19;//charge of an electron in C
10 z=1;//for hydrogen
11 n=1;
12 e=m*z^2*e^4/(8*e_0^2*h^2*n^2);//ionisation energy in
    J
13 disp(e,"Ionisation Energy (in J) = ");
14 e1=e/(1.602*10^-19);//in eV
15 disp(e1,"Ionisation Energy (in eV) = ");
```

---

### Scilab code Exa 2.6 Find dimensions of elliptical locus

```
1 //Find dimensionsof elliptical locus
2 //Ex:2.6
3 clc;
4 clear;
5 close;
6 n=4;//fourth orbit
7 a=(0+1)/n;//for s suborbit
8 disp(a,"For s suborbit b/a = ");
9 b=(1+1)/n;//for p suborbit
10 disp(b,"For p suborbit b/a = ");
11 c=(2+1)/n;//for d suborbit
```

```
12 disp(c,"For d suborbit b/a = ");
13 d=(3+1)/n;//for f suborbit
14 disp(d,"For f suborbit b/a = ");
```

---

### Scilab code Exa 2.7 Find Minimum Uncertainty in Position

```
1 //Find Minimum Uncertainty in Position
2 //Ex:2.7
3 clc;
4 clear;
5 close;
6 h=6.62*10^-34;//planck's constant in Js
7 p=10^-27;//uncertainty in momentum in kg m/s
8 x=h/(2*3.14*p);//uncertainty in position in m
9 disp(x,"Minimum Uncertainty in Position (in m) = ")
  ;
```

---

# Chapter 3

## CHEMICAL BONDING

Scilab code Exa 3.4 Find Energy of third and fifth orbit

```
1 //Find Energy of third and fifth orbit
2 //Ex:3.4
3 clc;
4 clear;
5 close;
6 e1=-13.6;//ionization potential of hydrogen in volts
7 n=3;//third orbit
8 e3=e1/(n^2);//Energy of third orbit in volts
9 disp(e3,"Energy of third orbit (in volts) = ");
10 m=5;//fifth orbit
11 e5=e1/(m^2);//Energy of fifth orbit in volts
12 disp(e5,"Energy of fifth orbit (in volts) = ");
```

---

Scilab code Exa 3.5 Find dipole moment and percentage ionic character

```
1 //Find Dipole moment and Percentage ionic character
2 //Ex:3.5
```

```
3 clc;
4 clear;
5 close;
6 d=0.9178; //molecular distance in angstorm
7 d1=d*10^-10; //in m
8 e=1.602*10^-19; //in C
9 dm=e*d1; //dipole moment in fully ionic state in Cm
10 disp(dm, "Dipole moment of HF in fully ionic state (
    in Cm) = ");
11 dm_m=6.375*10^-30; //measured dipole moment in Cm
12 p=(dm_m*100)/dm; //Percentage ionic character
13 disp(p, "Percentage ionic character");
```

---

## Chapter 4

# CRYSTALLOGRAPHY

Scilab code Exa 4.3 Find Lattice constant and atomic packing fraction of NaCl having FCC structure

```
1 //Find Lattice constant and atomic packing fraction
  of NaCl having FCC structure
2 //Ex:4.3
3 clc;
4 clear;
5 close;
6 r_na=0.98;//ionic radius of sodium in angstrom
7 r_cl=1.81;//ionic radius of chlorine in angstrom
8 n=4;//in fcc there are 4 Na and 4 Cl ions
9 a=((2*r_na)+(2*r_cl));//lattice constant
10 disp(a,"Lattice constant = ");
11 apf=((n*(4/3)*3.14*r_na^3)+(n*(4/3)*3.14*r_cl^3))/a
   ^3;
12 disp(apf,"atomic packing fraction of NaCl having FCC
   structure = ");
```

---

Scilab code Exa 4.4 Find density of Copper



```

1 //Find density of Copper
2 //Ex:4.4
3 clc;
4 clear;
5 close;
6 r=1.278;//radius of copper in angstorm
7 //copper has FCC structure
8 a=4*r/sqrt(2);//in angstorm
9 disp(a,"a (in Angstorm) = ");
10 a1=a*10^-8;//in cm
11 aw=63.54;//atomic weight of copper
12 ne=4;//fcc
13 na=6.023*10^23;//Avagadro's no.
14 p=aw*ne/(na*a1^3);//in g/cm^3
15 p1=10^3*p;
16 disp(p1,"Density of Copper (in kg/cu m)");

```

---

Scilab code Exa 4.5 Find distance between two adjacent atoms

```

1 //Find Distance between two adjacent atoms
2 //Ex:4.5
3 clc;
4 clear;
5 close;
6 w_na=23;//atomic weight of Na
7 w_cl=35.5;//atomic weight of Cl
8 w=w_na+w_cl;//effective no.of atoms in FCC structure
9 n=4;//FCC
10 na=6.023*10^23;//Avagadro's no.
11 w_4=w*n/na;//weight of 4 molecules in gm
12 p=2.18;//density in gm/cm^3
13 a=(w_4/p)^(1/3);//in cm
14 a1=a*10^8;//in angstorm

```

```
15 disp(a1,"unit cell dimension (in angstorm) = ");
16 d=a1/2;
17 disp(d,"Distance between two adjacent atoms (in
    Angstorm) = ");
```

---

#### Scilab code Exa 4.6 Find atomic radius

```
1 //Find atomic radius
2 //Ex:4.6
3 clc;
4 clear;
5 close;
6 n=2;//BCC
7 p=7.86;//density in gm/cm^3
8 aw=55.85;//atomic weight of iron
9 na=6.023*10^23;//Avagadro's no.
10 a=((aw*n)/(na*p))^(1/3);//in cm
11 a1=a*10^8;//in angstorm
12 disp(a1,"unit cell dimension of iron (in angstorm) =
    ");
13 r=sqrt(3)*a1/4;
14 disp(r,"atomic radius = ");
```

---

## Chapter 5

# MILLER INDICES AND X RAY CRYSTALLOGRAPH TECHNIQUES

Scilab code Exa 5.1 Find MILLER INDICES OF THE PLANE

```
1 //Find MILLER INDICES OF THE PLANE
2 //Ex:5.1
3 clc;
4 clear;
5 close;
6 p=1;
7 q=1/2;
8 r=3;
9 h=1/p;
10 k=1/q;
11 l=1/r;
12 h1=3*h;
13 k1=3*k;
14 l1=3*l;
15 disp(h1,"MILLER INDICES OF THE PLANE are h =");
16 disp(k1,"k = ");
17 disp(l1,"l = ");
```

---

**Scilab code Exa 5.3 Find MILLER INDICES OF THE PLANE**

```
1 //Find MILLER INDICES OF THE PLANE
2 //Ex:5.3
3 clc;
4 clear;
5 close;
6 p=2/4; //intercepts
7 q=3/3;
8 r=4/2;
9 h=1/p;
10 k=1/q;
11 l=1/r;
12 h1=2*h;
13 k1=2*k;
14 l1=2*l;
15 disp(l1,k1,h1,"MILLER INDICES ARE ");
```

---

**Scilab code Exa 5.5 Find interplanar Spacing**

```
1 //Find Interplanar Spacing
2 //Ex:5.5
3 clc;
4 clear;
5 close;
6 r=1.246; //radius in angstorm
7 h=2;
```

```

8 k=0;
9 l=0;
10 x=sqrt(h^2+k^2+l^2);
11 a=2*sqrt(2)*r;//in angstorm
12 d_200=a/x;//interplanar spacing in angstorm
13 disp(d_200,"Interplanar Spacing (200) (in Angstorm)
    = ");
14 h1=2;
15 k1=2;
16 l1=0;
17 x1=sqrt(h1^2+k1^2+l1^2);
18 d_220=a/x1;//interplanar spacing in angstorm
19 disp(d_220,"Interplanar Spacing (220) (in Angstorm)
    = ");
20 h2=1;
21 k2=1;
22 l2=1;
23 x2=sqrt(h2^2+k2^2+l2^2);
24 d_111=a/x2;//interplanar spacing in angstorm
25 disp(d_111,"Interplanar Spacing (111) (in Angstorm)
    = ");

```

---

Scilab code Exa 5.6 Find Linear Density per unit length

```

1 //Find Linear Density per unit length
2 //Ex:5.6
3 clc;
4 clear;
5 close;
6 a=3.61*10^-10;//unit cell in m
7 r_110=2/(sqrt(2)*a);//in atoms/m
8 r_a=r_110/10^3;//in atoms/mm
9 disp(r_a,"Linear Density per unit length along

```

```

        direction [110] (in atoms/mm) =");
10 r_111=1/(sqrt(3)*a); //in atoms/m
11 r_b=r_111/10^3; //in atoms/mm
12 disp(r_b,"Linear Density per unit length along
        direction [111] (in atoms/mm) =");

```

---

### Scilab code Exa 5.7 Find Planar Density

```

1 //Find Planar Density
2 //Ex:5.7
3 clc;
4 clear;
5 close;
6 r_po=1.7*10^-10; //radius of polonium in m
7 r_rh=1.34*10^-10; //radius of rhodium in m
8 r_cr=1.25*10^-10; //radius of chromium in m
9 a_po=2*r_po; //in m
10 a_rh=2*sqrt(2)*r_rh; //in m
11 a_cr=4*r_cr/sqrt(3);
12 p_po=1/a_po^2; // /sqm
13 disp(p_po,"Planar Density on [100] in Polonium (per
        sqm) = ");
14 p_rh=1.414/a_rh^2; // /sqm
15 disp(p_rh,"Planar Density on [110] in Rhodium (per
        sqm) = ");
16 p_cr=1.732/a_cr^2; // /sqm
17 disp(p_cr,"Planar Density on [111] in Chromium (per
        sqm) = ");

```

---

Scilab code Exa 5.8 Find Glancing angle and Interplanar spacing of the crystal

```
1 //Find Glancing angle and Interplanar spacing of the
   crystal
2 //Ex:5.8
3 clc;
4 clear;
5 close;
6 w=0.824; //wavelength in angstorm
7 a1=8.58; //angle at n=1 in degrees
8 n1=1;
9 n3=3;
10 a3=asind((n3/n1)*sind(a1)); //angle at n=3 in degrees
11 disp(a3,"Glancing angle for third order diffraction
   = ");
12 d=w/(2*sind(a1)); //in angstorm
13 disp(d,"Interplanar spacing of the crystal (in
   Angstorm) = ");
```

---

Scilab code Exa 5.9 Find Glancing angle and lattice parameter

```
1 //Find Glancing angle and lattice parameter
2 //Ex:5.9
3 clc;
4 clear;
5 close;
6 a=17.03; //in degrees
7 w=0.71; //in angstorm
8 n=1;
9 d=n*w/(2*sind(a)); //interplanar spacing in angstorm
10 disp(d,"Interplanar Spacing (in angstorm) = ");
11 // given that  $h^2+k^2+l^2=8$ 
12 a=sqrt(8)*d; //in angstorm
```

```
13 disp(a,"Lattice parameter of the crystal (in  
    Angstorm) = ");
```

---

**Scilab code Exa 5.10 Find Dimension of unit cell**

```
1 //Find Dimension of unit cell  
2 //Ex:5.10  
3 clc;  
4 clear;  
5 close;  
6 w=0.0708;//wavelength in nm  
7 h=1;  
8 k=0;  
9 l=0;  
10 s=0.0132;//a common divisor i.e. sin^2(theta)=0.0132  
11 a=sqrt((w^2*(h^2+k^2+l^2))/(4*s));//in nm  
12 a1=10^3*a;//in pm  
13 disp(a1,"Dimension of unit cell (in Picometer) = ");
```

---



## Chapter 6

# IMPERFECTIONS DEFECTS AND DISLOCATIONS IN SOLIDS

Scilab code Exa 6.1 Find No of Frenkel defect

```
1 //Find No. of Frenkel defect
2 //Ex:6.1
3 clc;
4 clear;
5 close;
6 na=6.023*10^23;//Avagadro's no.
7 p=3170;//density in kg/m^3
8 mw=7.9*10^-2;//molecular weight of CaF2
9 nl=na*p/mw;//calcium ions/cubic m
10 ni=2*nl;// /cubic m
11 t=1300;//in K
12 ef=2.7*1.6*10^-19;//energy of formation of one
    frenkel defect
13 k=1.38*10^-23;//boltzmann constant
14 nf=sqrt(nl*ni)*exp(-ef/(2*k*t));//in /cubic m
15 disp(nf,"No. of Frenkel defect per unit volume of
    Calcium Fluoride (in /m^3) = ");
```

---

**Scilab code Exa 6.2 Calculate Ratio of no of vacancies**

```
1 //Calculate Ratio of no. of vacancies
2 //Ex:6.2
3 clc;
4 clear;
5 close;
6 r=8.314; // J/mol K
7 t1=300; //in K
8 ent=168*103; //enthalpy of formation of vacancy in J
   /mol
9 x1=exp(-ent/(r*t1)); //x1=n/Na
10 t2=1000; //in K
11 x2=exp(-ent/(r*t2)); //x2=n/Na
12 rt=x1/x2; //ratio
13 disp(rt,"Ratio of no. of vacancies = ");
```

---

**Scilab code Exa 6.4 Find Elastic Strain Energy**

```
1 //Find Elastic Strain Energy
2 //Ex:6.4
3 clc;
4 clear;
5 close;
6 v=0.31; //poisson 's ratio
7 bv=.25*10-9; //burger 's vector in m
8 ri=1.1*10-9; //in m
```

```

 9 r0=10^5*bv; //in m
10 sm=45*10^9; //shear modulus in n/sqm
11 gb_2=sm*bv^2;
12 u_ed=(gb_2/(4*3.14*(1-v)))*log(r0/ri);
13 disp(u_ed,"Elastic Strain Energy of Edge dislocation
      (in J/m) = ");
14 u_sd=(gb_2/(4*3.14))*log(r0/ri);
15 disp(u_sd,"Elastic Strain Energy of Screw
      dislocation (in J/m) = ");
16 r=u_ed/u_sd; //ratio
17 disp(r,"Ratio of energies of edge dislocation over
      screw dislocation = ");

```

---

**Scilab code Exa 6.5** Calculate Total number of created vacancies

```

1 //Calculate Total no. of created vacancies
2 //Ex:6.5
3 clc;
4 clear;
5 close;
6 r=1.7*10^-10; //atomic radius in m
7 n1=10^-3; //lmm=10^-3m
8 a=2*r; //in m
9 n=n1/a;
10 ed=2*10^-6; //edge dislocation in m
11 ns=ed/a;
12 nv=n*ns;
13 disp(nv,"Total no. of created vacancies = ");

```

---

# Chapter 7

## MECHANICAL PROPERTIES

Scilab code Exa 7.3 Find various Modulous of elasticity

```
1 //Find various Modulous of elasticity
2 //Ex7.3
3 clear;
4 close;
5 a1=222*10^9; //in N
6 a2=168*10^9; //in N
7 e1=1.90; //in sqm
8 e2=1.42; //in sqm
9 da=a1-a2; //in N
10 de=e1-e2; //in sqm
11 e_tan=da/de;
12 e_tann=e_tan*10^-9; //in Gpa
13 disp(e_tann,"Tangent Modulous of elasticity (in Gpa)
    = ");
14 a3=180*10^9; //in N
15 e3=1.46; //in sqm
16 e_sec=10^-9*a3/e3; //in Gpa
17 disp(e_sec,"Secant modulous of elasticity (in Gpa) =
    ");
18 a=85*10^6;
19 e=.68*10^-3;
```

```
20 e_y=10^-9*a/e; //in Gpa
21 disp(e_y,"Youngs modulus (in Gpa) = ");
```

---

### Scilab code Exa 7.5 Find stress

```
1 //Find Stress
2 //Ex:7.5
3 clc;
4 clear;
5 close;
6 n=3;
7 a=300;
8 v_cr=2.8*10^-8; // in cm/cm/hour creep rate
9 x=log(v_cr)-n*log(a);
10 a1=exp(x);
11 t=365*24; //in hours
12 e=2*10^6; //kgf/sqcm
13 ai=750; //in kgf/sqcm
14 a_tf=sqrt(1/((1/ai^(n-1))+(a1*e*(n-1)*t)));
15 disp(a_tf,"Stress Remaining (in kgf/sq cm) = ");
```

---

# Chapter 8

## MECHANICAL TESTING

Scilab code Exa 8.1 Find Flexural Strength Shear Strength and Modulus of Rupture

```
1 //Find Flexural Strength Shear Strength and
  Modulus of Rupture
2 //Ex:8.1
3 clc;
4 clear;
5 close;
6 b=225;//in mm
7 h=10;//in mm
8 l=1100;//in mm
9 f1=250;//in N
10 m=f1*l/4;//in N-mm
11 f=f1/2;//in N
12 a=(6*m)/(b*h^2);//in N/mm^2
13 disp(a,"Flexural Strength (in N/sqmm) = ");
14 t=(3*f)/(2*b*h);//in N/sqmm
15 disp(t,"Shear Strength (in N/sqmm) = ");
16 f2=350;//in N at which glass breaks
17 r=f2*l/4;//in N-mm
18 i=(b*h^3)/12;//in mm^4
19 y=h/2;//in mm
20 mr=r*y/i;//in n/sqmm
```

```
21 disp(mr,"Modulous of Rupture (in N/sqmm) = ");
```

---

**Scilab code Exa 8.2 FIND BRINELL HARDNESS NUMBER**

```
1 //FIND BRINELL HARDNESS NUMBER
2 //Ex:8.2
3 clc;
4 clear;
5 close;
6 d=5;//in mm
7 id=32.5/10;//indentation diameter in mm
8 p=30*d^2;//load for steel specimen in kgf
9 disp(p,"Load P for steel specimen (in kgf) = ");
10 bhn=p/((3.14*d/2)*(d-sqrt(d^2-id^2)));//in kgf/sqmm
11 disp(bhn,"BRINELL HARDNESS NUMBER of the steel
    specimen = ");
```

---

**Scilab code Exa 8.3 Find Rupture Energy Modulous Of Rupture and Notch Imapct Stren**

```
1 //Find Rupture Energy Modulous Of Rupture and Notch
    Imapct Strength
2 //Ex:8.3
3 clc;
4 clear;
5 close;
6 l=0.1;//frictinal and windage losses in kgf-m
7 dr=5.9;//dial reading in kgf-m
8 u=dr-l;//in kgf-m
9 disp(u,"Rupture Energy (in kgf-m) = ");
```

```

10 t=10; //in mm
11 d=t/5; //depth of V-notch in mm
12 te=t-d; //effective thickness in mm
13 ve=75*10*te; //effective volume in cu. mm
14 vem=ve*10^-9; //in cu. m
15 mr=u/vem; //in kgf/sqm
16 disp(mr, "Modulus Of Rupture (in kgf/sqm) = ");
17 ae=t*te; //effective area of cross section in sqmm
18 aem=ae*10^-6; //in sqm
19 is=u/aem; //in kg/m
20 disp(is, "Notch Impact Strength (in kg/m) = ");
21 ui=30; //in kgf-m
22 a=160; //angle in degrees
23 r=0.8; //swing radius in m
24 uf=ui-u; //in kgf-m
25 w=19.33; //weight of hammer in kgf-m
26 hf=uf/w; //in m
27 disp(hf, "Height risen by Hammer (in m) = ");
28 //hf=r*(1-cos(b))
29 b=acosd((r-hf)/r); //in degrees
30 disp(b, "Angle after Breaking the specimen (in
    degress) = ");

```

---

#### Scilab code Exa 8.4 Find Stress Ratio and range

```

1 //Find Stress Ratio and range
2 //Ex:8.4
3 clc;
4 clear;
5 close;
6 a_m=70; //mean stress in Mpa
7 a_r=210; //stress amplitude in Mpa
8 a_max=((2*a_m)+a_r)/2; //maximum stress in MPa

```



```

9 disp(a_max,"Maximum Stress Level (in MPa) = ");
10 a_min=2*a_m-a_max;//Minimum stress in MPa
11 disp(a_min,"Minimum Stress Level (in MPa) = ");
12 s=a_min/a_max;//stress ratio
13 disp(s,"Stress Ratio = ");
14 sr=a_max-a_min;//stress range in MPa
15 disp(sr,"Stress Range (in MPa) = ");

```

---

Scilab code Exa 8.5 Calculate ENDURANCE STRESS FROM using various relations

```

1 //Calculate ENDURANCE STRESS FROM using various
  relations
2 //Ex:8.5
3 clc;
4 clear;
5 close;
6 p_min=20;//in kN
7 p_max=50;//in kN
8 l=500;//in mm
9 d=60;//in mm
10 a_u=650;//in MPa
11 a_y=520;//in MPa
12 fos=1.8;//factor of safety
13 m_max=p_max*l/4;//maximum bending moment in kN mm
14 m_min=p_min*l/4;//minimum bending moment in kN mm
15 m_m=(m_max+m_min)/2;//mean bending moment in kN mm
16 m_a=(m_max-m_min)/2;//alternating bending moment in
  kN mm
17 z=3.14*d^3/32;
18 a_m=(m_m/z)*1000;//mean bending stress in MPa
19 a_a=(m_a/z)*1000;//alternating bending stress in MPa
20 a_e1=a_a/((1/fos)-(a_m/a_u)^2*fos);//in MPa
21 disp(a_e1,"ENDURANCE STRESS FROM Gerbers Parabolic

```

```
Function (in MPa) = ");
22 a_e2=a_a/((1/fos)-(a_m/a_u)); //in MPa
23 disp(a_e2,"ENDURANCE STRESS FROM Goodman Straight
Line Relation (in MPa) = ")
24 a_e3=a_a/((1/fos)-(a_m/a_y)); //in MPa
25 disp(a_e3,"ENDURANCE STRESS FROM Soderberg Straight
Line Relation (in MPa) = ")
```

---

## Chapter 9

# MICROSTRUCTURAL EXAMINATION AND NON DESTRUCTIVE TESTING

Scilab code Exa 9.1 Find grain diameter

```
1 //Find grain diameter
2 //Exa:9.1
3 clc;
4 clear;
5 close ;
6 n=2^(12-1); //astm no.=12
7 //1 sq inch=645mm^2
8 d=1/sqrt((n/645)*10^4); //grain diameter in mm
9 disp(d,"grain diameter for ASTM no. 12 (in mm) = ");
```

---

Scilab code Exa 9.2 Find Average and Boundary area

```

1 //Find Average and Boundary area
2 //Exa:9.2
3 clc;
4 clear;
5 close ;
6 n=2^(5-1); //astm no. = 5 in grain/inch^2
7 //Lineal and Areal magnifications are related as
   *100 Lineal = *10000 Areal
8 x=n/(.01*.01); //in grain/inch^2 at 1 x.
9 a=1/x; //average area in inch^2
10 a1=2.54*2.54*a; //average area in cm.^2
11 disp(a1,"Average area of one grain (in sq cm) = ");
12 l=sqrt(x); //grains/inch of length
13 s=(1/l)^2; //surface area in sq inch
14 s6=6*s; //surface area of 6 surfaces of cubic grain
   in sq inch
15 b=0.5*s6*(1^3); //total boundary area in sq inch
16 b1=b/(2.54); //total boundary area in sq cm
17 disp(b1,"Boundary Area per cubic centimetre of steel
   (in sq cm) = ");

```

---

## Chapter 10

# PHASE DIAGRAM AND EQUILIBRIUM DIAGRAM

Scilab code Exa 10.1 Prove Two component system cannot have more than 4 phases in

```
1 //Prove Two component system cannot have more than 4
   phases in an equilibrium
2 //Exa:10.1
3 close;
4 clc;
5 clear;
6 c=2; //for 2 component system
7 disp("D=C-P+2");
8 disp("Total no. of variables = P*(C-1)+2");
9 p=4;
10 d=c-p+2; //degree of freedom
11 t=p*(c-1)+2; //no. of total variables
12 disp(d,"Degree of freedom = ");
13 disp(p,"when p = ");
14 disp("Two component system cannot have more than 4
   phases in an equilibrium");
```

---

Scilab code Exa 10.2 Calculate Total weight of lead and tin

```
1 //Calculate Total weight of lead and tin
2 //Exa:10.2
3 close;
4 clc;
5 clear;
6 p_pb=11364.1;//density of lead in kg/m^3
7 p_sn=7220.14;//density of tin in kg/m^3
8 p_e=100/((38/p_pb)+(62/p_sn));//density of eutectic
    composition at point D
9 disp(p_e,"Density of eutectic composition at point D
    (in kg/m3) = ");
10 w=.88*p_e;//in kgf
11 w_pb=.38*w;//of lead in kgf
12 w_sn=.62*w;//of tin in kgf
13 p_b=7300;//density in beta phase in kg/m^3
14 w1=.12*p_b;//in kgf
15 w1_pb=.03*w1;//of lead in kgf
16 w1_sn=.97*w1;//of tin in kgf
17 w2_pb=w_pb+w1_pb;//Total weight of lead in kgf
18 disp(w2_pb,"Total weight of lead (in kgf) = ");
19 w2_sn=w_sn+w1_sn;//Total weight of tin in kgf
20 disp(w2_sn,"Total weight of tin (in kgf) = ");
21 sn=(w2_sn/(w2_sn+w2_pb))*100;
22 disp(sn,"% of Sn = ");
```

---

Scilab code Exa 10.4 Find weight fractions

```

1 //Find weight fractions
2 //Ex:10.4
3 clc;
4 clear;
5 close;
6 c_be=100;
7 c_e=1.65;
8 c_o=10;
9 w=(c_be-c_o)/(c_be-c_e);
10 disp(w,"weight fractions = ");

```

---

Scilab code Exa 10.5 Find Maximum weight of tin

```

1 //Find Maximum weight of tin
2 //Ex:10.5
3 clc;
4 clear;
5 close;
6 //m be amount of tin
7 w_sn=900;//weight of tin
8 w_pb=1000;//weight of lead
9 m=((w_pb*0.97)-w_sn)/(1-.97);//in grams
10 m1=m/1000;//maximum mass of tin in kgm
11 disp(m1,"Maximum weight of tin that can be added
    without changing systems temperature (in kgm) = "
    );

```

---

Scilab code Exa 10.6 Find Weight fraction of errite and Cementite

```
1 //Find Weight fraction of errite and Cementite
2 //Ex:10.6
3 clc;
4 clear;
5 close;
6 c=0.83;//carbon
7 f=0;//ferrite
8 ce=6.67;//cementite
9 w_a=(ce-c)/(ce-f);
10 disp(w_a,"Weight fraction of errite = ");
11 w_b=(c-f)/(ce-f);
12 disp(w_b,"Weight Fraction of Cementite = ");
```

---



# Chapter 12

## HEAT TREATMENT

Scilab code Exa 12.1 calculate percentage change in volume

```
1 //calculate percentage change in volume
2 //Ex:12.1
3 clc;
4 clear;
5 close;
6 n_bcc=1/2;
7 n_fcc=1/4;
8 r_fcc=1.26;//in Angstorm
9 r_bcc=1.24;//in Angstorm
10 a_bcc=4*r_bcc/sqrt(3);//in Angstorm
11 a_fcc=2*sqrt(2)*r_fcc;//in Angstorm
12 v_fcc=a_fcc^3;
13 v_bcc=a_bcc^3;
14 v=100*((n_fcc*v_fcc)-(n_bcc*v_bcc))/(n_fcc*v_fcc);
15 disp(v,"Percentage change in volume = ");
```

---

Scilab code Exa 12.3 Estimate free energy change during recrystallization

```

1 //Estimate free energy change during
  recrystallization
2 //Ex:12.3
3 clc;
4 clear;
5 close;
6 e=51; //Young modulus in GPa
7 v=0.22; //poisson ratio
8 g=e/(2*(1-v)); //shear modulus in GPa
9 b=2*10^-10;
10 ue=(1/2)*g*10^9*b^2*10^12; //in J/m^3
11 disp(ue,"Change in free energy during
  recrysatlization (in J/m^3) = ");

```

---

**Scilab code Exa 12.5** Find contribution of particles

```

1 //Find contribution of particles
2 //Ex:12.5
3 clc;
4 clear;
5 close;
6 g=41*10^9; //in N/m^2
7 b=0.64*10^-9; //in m
8 l=20*10^-6; //in m
9 t=g*b/l; //in N/m^2
10 T=t*10^-6; //in MPa
11 disp(T,"Contribution of these particles (in MPa) = "
  );

```

---

## Chapter 14

# MAGNETIC PROPERTIES AND MATERIALS

Scilab code Exa 14.1 Find relative permeability and Intensity of magnetisation

```
1 //Find relative permeability and Intensity of
  magnetisation
2 //Ex:14.1
3 clc;
4 clear;
5 close;
6 x=1500; //susceptibility
7 h=2400; //magnetic field in A/m
8 u_r=1+x;
9 disp(u_r,"relative permeability = ");
10 m=x*h; //in A/m
11 disp(m,"Intensity of magnetisation (in A/m)");
12 u_0=4*3.14*10^-7;
13 b=u_0*u_r*h; //in T
14 disp(b,"Remanance (in T) = ")
```

---

**Scilab code Exa 14.2** Estimate Hysteresis energy loss per unit volume

```
1 //Estimate Hysteresis energy loss per unit volume
2 //Ex:14.2
3 clc;
4 clear;
5 close;
6 w=80+80;//width of loop in A/m from graph
7 h=0.15+.15;//height of loop in Wb/sqm
8 a=w*h;//area of the loop in J
9 disp(a,"Hysteresis energy loss per unit volume of
      magnetic material during one cycle (in J) = ");
```

---

**Scilab code Exa 14.3** Find Power loss due to Hysteresis

```
1 //Find Power loss due to Hysteresis
2 //Ex:14.3
3 clc;
4 clear;
5 close;
6 a=600;//loop area in J/sqm
7 f=50;//in Hz
8 v=0.01//volume in cu. m
9 w=a*f*v;//in W
10 disp(w,"Power loss due to Hysteresis (in W) = ");
```

---

#### Scilab code Exa 14.4 Find Loss at 40 Hz

```
1 //Find Loss at 40 Hz
2 //Ex:14.4
3 clc;
4 clear;
5 close;
6 w_h1=300;
7 b_1=0.9;
8 y=b_1^1.7;
9 b_2=1.1;
10 x=b_2^1.7;
11 f1=50;
12 f2=40;
13 w_h2=(w_h1*x*f2)/(y*f1);
14 w_h22=w_h2*1.22; //in W
15 disp(w_h22,"Loss at 40 Hz (in W) = ");
```

---

#### Scilab code Exa 14.5 Find Magnetic strength

```
1 //Find Magnetic strength
2 //Ex:14.5
3 clc;
4 clear;
5 close;
6 m=6000;
7 u_r=200000;
8 x=u_r-1;
```

```
9 h=m/x;
10 u_0=4*3.14*10^-7;
11 b=u_0*u_r*h;
12 disp(b,"Magnetic strength (in T) = ");
```

---

#### Scilab code Exa 14.6 Estimate Saturation Magnetisation

```
1 //Estimate Saturation Magnetisation
2 //Ex:14.6
3 clc;
4 clear;
5 close;
6 b=9.27*10^-24; //Bohr Magneton in A/sqm
7 m=0.6*b;
8 a=0.35*10^-9;
9 n=4; //FCC
10 m_g=n*m/a^3; //in A/m
11 disp(m_g,"Saturation Magnetisation (in A/m) =");
```

---

#### Scilab code Exa 14.9 Find Eddy current loss at 60 and 100 Hz

```
1 //Find Eddy current loss at 60 and 100 Hz
2 //Ex:14.9
3 clc;
4 clear;
5 close;
6 f=50; //in Hz
7 L=100; //Eddy current loss in transformer in W
8 f1=60; //in Hz
```

```

9  w_e=L*(f1/f)^2; //in W
10 disp(w_e,"Eddy current loss at 60 Hz (in W) = ");
11 f2=100; //in Hz
12 w_ee=L*(f2/f)^2; //in W
13 disp(w_ee,"Eddy current loss at 100 Hz (in W) = ");

```

---

**Scilab code Exa 14.13** Find Magnetic field strength and Flux density

```

1 //Find Magnetic field strength and Flux density
2 //Ex:14.13
3 clc;
4 clear;
5 close;
6 l=.25; //in m
7 n=400; //turns
8 i=15; //in A
9 u_0=1.257*10^-6; //in H/m
10 h=n*i/l; //in AT/m
11 disp(h,"Magnetic field strength (in AT/m) = ");
12 u_r=1; //relative permeability
13 b=u_0*u_r*h; //in wB/sqm
14 disp(b,"Flux density (in Wb/sq m) = ");

```

---

# Chapter 15

## ELECTRIC PROPERTIES

Scilab code Exa 15.1 Find Electric field

```
1 //Find Electric field
2 //Ex:15.1
3 clc;
4 clear;
5 close;
6 v=230;//in volts
7 d=0.005;//in m
8 E=-v/d;//in V/m
9 disp(E,"Electric field between pair of conducting
   plates (in V/m) = ");
```

---

Scilab code Exa 15.2 Find Drift Velocity

```
1 //Find Drift Velocity
2 //Ex:15.2
3 clc;
4 clear;
```



```

5  close;
6  n=1019; //no. of electrons per unit volume
7  e=1.602*10-19; //charge of an electron in C
8  a=0.018; //conductivity in ohm/m
9  m=9.1*10-31; //mass of an electron in kg
10 v=0.16; //in volts
11 t=0.29; //thickness in mm
12 efg=v/t; //electric field gradient in V/m
13 vd=a*efg/(n*e);
14 vd1=103*vd; //in m/s
15 disp(vd1," Drift Velocity (in m/sec) = ");

```

---

### Scilab code Exa 15.3 Find Specific Resistance

```

1  //Find Specific Resistance
2  //Ex:15.3
3  clc;
4  clear;
5  close;
6  l=200; //in m
7  r=21; //in ohm
8  d=0.44*10-3; //in m
9  a=3.14*(d/2)2; //area in sq m
10 p=r*a/l; //in ohm-m
11 disp(p," Specific Resistance (in ohm-m) = ");

```

---

### Scilab code Exa 15.4 Find Resistivity

```

1  //Find Resistivity

```

```

2 //Ex:15.4
3 clc;
4 clear;
5 close;
6 p_cu=0.015*10^-6; //resistivity of copper in ohm-m
7 p_ni=0.012*10^-6; //resistivity of nickel in ohm-m
8 p_ag=0.016*10^-6; //resistivity of silver in ohm-m
9 c1=0.25; //atomic % of nickel
10 c2=0.4; //atomic % of silver
11 p=p_cu+(c1*p_ni)+(c2*p_ag);
12 disp(p,"Resistivity of Cu-Ni-Ag alloy at 300 K (in
    ohm-m) = ");

```

---

#### Scilab code Exa 15.5 Calculate Intrinsic Carrier density

```

1 //Calculate Intrinsic Carrier density
2 //Ex:15.5
3 clc;
4 clear;
5 close;
6 m=0.14; //mobility of electron
7 u_h=0.05; //mobility of holes
8 p=3000; //resistivity in ohm-m
9 e=1.602*10^-19; //charge of an electron in C
10 a=1/p; //conductivity
11 n=a/(e*(m+u_h));
12 disp(n,"Intrinsic Carrier density in pure silicon (
    in per cu m) = ");

```

---

**Scilab code Exa 15.6** Estimate Drift Velocity and time taken by electrons

```
1 //Estimate Drift Velocity and time taken by
   electrons
2 //Ex:15.6
3 clc;
4 clear;
5 close;
6 id=1000; //in A/sqm
7 p=0.05; //resistivity in ohm-m
8 l=100*10^-6; //in m
9 m_e=0.4; //in sqm/Vsec
10 e=1.602*10^-19; //charge of electron in C
11 a=1/p; //conductivity
12 n_e=a/(e*m_e); //in per cubic m
13 v_d=id/(n_e*e); //in m/s
14 disp(v_d," Drift Velocity (in m/s) = ");
15 t=l/v_d; //in sec
16 t1=t*10^6; //in msec
17 disp(t1,"Time taken by electrons (in msec) = ");
```

---

**Scilab code Exa 15.7** Find Impurity concentration

```
1 //Find Impurity concentration
2 //Ex:15.7
3 clc;
4 clear;
5 close;
6 d=1*10^-3; //diameter in m
7 a=3.14*(d/2)^2; //area of cross section of rod in sq
   m
8 r=100; //in ohm
9 l=10*10^-3; //in m
```

```

10 p=a*r/l; //in ohm-m
11 c=1/p; //conductivity
12 e=1.602*10^-19; //charge of electron in C
13 u_h=0.19; //mobility of holes in sqm/Vsec
14 n_h=c/(e*u_h);
15 disp(n_h,"Impurity concentration in rod (in per
    cubic m) = ");

```

---

**Scilab code Exa 15.8** Calculate Conduction electron and hole density

```

1 //Calculate Conduction electron and hole density
2 //Ex:15.8
3 clc;
4 clear;
5 close;
6 ni=1.5*10^16; //intrinsic carrier concentration per
    cu. m
7 n=10^19; //no. of conduction electrons in per cu. m
8 p=ni^2/n; //in per cu.m
9 disp(p,"Conduction electron and hole density (per
    cubic m) = ");

```

---

**Scilab code Exa 15.9** Calculate Hole concentration

```

1 //Calculate Hole concentration
2 //Ex:15.9
3 clc;
4 clear;
5 close;

```

```

6 nd=10^17; //in per cu cm
7 ni=1.5*10^10; //in cu cm
8 ne=nd; //nd>>ni
9 nh=ni^2/ne;
10 disp(nh,"Hole concentration (in per cubic cm) = ");
11 t=300; //in K
12 e=0.0259*log(ne/ni); //in eV
13 disp(e,"Location of Fermi Level (in eV) = ");

```

---

**Scilab code Exa 15.10 Find thickness of insulation**

```

1 //Find thickness of insulation
2 //Ex:15.10
3 clc;
4 clear;
5 close;
6 d=40000; //dielectric strength in V/mm
7 v=33*10^3; //in volts
8 t=v/d; //in mm
9 disp(t,"thickness of insulation (in mm) = ");

```

---

**Scilab code Exa 15.16 Find Band gap energy**

```

1 //Find Band gap energy
2 //Ex:15.16
3 clc;
4 clear;
5 close;
6 c=2.99*10^8; //speed of light in m/s

```

```
7 h=6.62*10^-24; //planck 's constant
8 w=1.771*10^-6; //wavelength in J
9 eg=h*c/w; //in J
10 disp(eg,"Band gap energy (in J) = ");
```

---

## Chapter 16

# SUPERCONDUCTIVITY AND SUPERCONDUCTORS

Scilab code Exa 16.1 find susceptibilty and relative permeability of a superconductor

```
1 //find susceptibilty and relative permeability of a
  superconductor
2 //Ex:16.1
3 clc;
4 clear;
5 close;
6 b=0;
7 //m=-h
8 //m=x*h
9
10 //==>> -h=x*h
11 x=-1; //from above realtions
12 disp(x, "Susceptibility of superconductor = ");
13 ur=x+1; //relative permeability
14 disp(ur, "Relative permeability of superconductor = "
  );
```

---

### Scilab code Exa 16.2 Determine critical field

```
1 //Determine critical field
2 //Ex:16.2
3 clc;
4 clear;
5 close;
6 ho=0.0803; //in A/m
7 t1=3; //in K
8 t2=10; //in k
9 tc=7.17; //in K
10 hc1=ho*(1-(t1/tc)^2);
11 disp(hc1," Critical field at 3K (in A/m) = ");
12 hc2=ho*(1-(t2/tc)^2);
13 disp(hc2," Critical field at 10K (in A/m) = ");
```

---

### Scilab code Exa 16.3 Find critical current

```
1 //Find critical current
2 //Ex:16.3
3 clc;
4 clear;
5 close;
6 r=1*10^-3; //in m
7 hc=7.9*10^3; //in A/m
8 ic=2*3.14*r*hc; //in m
9 disp(ic," Critical current in superconducting state (
    in A) = ");
```



---

Scilab code Exa 16.4 Find electron and current density

```
1 //Find electron and current density
2 //Ex:16.4
3 clc;
4 clear;
5 close;
6 p=11.4*10^3;//in kg/m^3
7 aw=207.2;//in kg/kg-mol
8 v=1200;//in m/s
9 na=60.23*10^26;//avagadro's no
10 e=1.6*10^-19;//charge in C
11 m=9.1*10^-31;//mass of electron in kg
12 mo=4*3.14*10^-7;//in H/m
13 ne=2*p*na/aw;//in per m^3
14 disp(ne,"Electron density (in per m^3) = ");
15 ied=ne*e*v;//in A/m^2
16 disp(ied,"Current density (in A/m^2) = ");
17 dp=(m/(mo*ne*(e^2)))^(1/2);
18 dp1=dp*10^10;
19 disp(dp1,"Depth of penetration (in angstorm) = ");
```

---

Scilab code Exa 16.9 determine critical current density

```
1 //determine critical current density
2 //Ex:16.9
3 clc;
```

```

4 clear;
5 close;
6 ho=65*10^3; //in A/m
7 tc=7.18; //in K
8 t=4.2; //in K
9 r=0.5*10^-3; //in m
10 hc=ho*(1-(t/tc)^2); //in A/m
11 ic=2*3.14*r*hc; //in A
12 a=3.14*r^2; //area in m^2
13 j=ic/a; //in A/m^2
14 disp(j,"current density (in A/m^2) = ");

```

---

Scilab code Exa 16.10 Determine transition temperature and critical field

```

1 //Determine transition temperature and critical
  field
2 //Ex:16.10
3 clc;
4 clear;
5 close;
6 hc1=21; //in A/m
7 hc2=10; //in A/m
8 tc=7; //in K
9 t=14; //in K
10 h=hc1/hc2;
11 //Determining critical temperature
12 tc1=sqrt(3626/11); //by quadratic eqn in the example
13 ho=hc1/(1-(tc^2/tc1^2));
14 disp(ho,"Critical field at 0 K (in A/m) = ");
15 t=4.2; //in k
16 hc=ho*(1-(t/tc1)^2);
17 disp(hc,"Critical field At 4.2 k (in A/m) = ");

```

---



# Chapter 17

## CERAMICS AND PLASTICS

Scilab code Exa 17.1 Calculate molecular weight

```
1 //Calculate molecular weight
2 //Ex:17.1
3 clc;
4 clear;
5 close;
6 mc=12;//mol wt of carbon
7 mh=1;;//mol wt of hydrogen
8 m=8*(mc+mh);//mol wt of C8H8
9 DOP=10000;//degree of polarization , given
10 mp=DOP*m;
11 disp(mp,"Molecular weight of Styrene polymer = ");
```

---

Scilab code Exa 17.2 Determine molecular weight of teflon

```
1 //Determine molecular weight of teflon
2 //Ex:17.2
3 clc;
```

```
4 clear;
5 close;
6 DOP=10000;
7 mc=12; //mol wt of carbon
8 mf=19; //mol wt of fluorine
9 m=(2*mc)+(4*mf); //mol wt of teflon monomer
10 mp=DOP*m;
11 disp(mp,"Molecular weight of Teflon polymer = ");
12 mh=1; //mol wt of hydrogen
13 m1=(2*mc)+(4*mh); //mol wt of polyethylene
14 //for same DOP
15 x=m/m1; //ratio of molecular weights
16 disp(x,"Ratio of molecular weights of Teflon and
    Polyethylene = ");
```

---

# Chapter 19

## COMPOSITE MATERIALS

Scilab code Exa 19.1 Find flexural rigidity of sandwich construction

```
1 //Find flexural rigidity of sandwich construction
2 //Ex:19.1
3 clc;
4 clear;
5 close;
6 t_s=3; //in mm
7 t_c=24; //in mm
8 b=100; //in mm
9 d=(t_s+t_c)/2; //in mm
10 is=((b*t_s^3)/12)+(b*t_s*d^2); //in mm^4
11 ic=b*t_c^3/12; //in mm^4
12 m_p=7000; //moduli of polyester skin in N/mm^2
13 m_f=20; //moduli of foam core in N/mm^2
14 d_fr=(2*m_p*is)+(m_f*ic); //in N/mm^2
15 disp(d_fr,"Flexural rigidity (in N/sqm) = ");
```

---

Scilab code Exa 19.2 Determine volume ratio of Al and B in aluminium boron composi

```

1 //Determine volume ratio of Al and B in aluminium
  boron composite
2 //Ex:19.2
3 clc;
4 clear;
5 close;
6 ec=210;//in GPa
7 ea=71;//in GPa
8 eb=440;//in GPa
9 va=(ec-eb)/(ea-eb);
10 disp(va,"Va = ");
11 vb=1-va;
12 disp(vb,"Vb = ");
13 c=vb/va;
14 disp(c,"Volume ratio = ");

```

---

**Scilab code Exa 19.3** Calculate fraction of load carried by fibres

```

1 //Calculate fraction of load carried by fibres
2 //Ex:19.3
3 clc;
4 clear;
5 close;
6 ef=430;//in GPa
7 e=3.6;//in GPa
8 m=ef/e;
9 vf=0.15;//by volume
10 vm=1-vf;
11 x=vm/vf;
12 pf=m;
13 pc=m+x;
14 y=pf/pc;
15 disp(y,"fraction of load carried by fibres (15 % by

```

```

        volume) = ");
16 vf1=0.65
17 vm1=1-vf1;
18 z=vm1/vf1;
19 pc1=m+z;
20 zz=pf/pc1;
21 disp(zz,"fraction of load carried by fibres (65 % by
        volume) = ")

```

---

Scilab code Exa 19.4 Find longitudinal strength longitudinal modulus transverse modulus Poisson ratio Shear modulus

```

1 //Find longitudinal strength longitudinal modulus
  transverse modulus Poisson ratio Shear modulus
2 //Ex:19.4
3 clc;
4 clear;
5 close;
6 vf=0.65;
7 vm=1-vf;
8 kts=2.8; //in GPa
9 ets=0.0025; //in GPa
10 ac=(kts*vf)+(ets*vm); //in GPa
11 disp(ac,"Longitudinal Strength (in GPa) = ");
12 ktm=130; //in GPa
13 etm=3.5; //in GPa
14 ec=(ktm*vf)+(etm*vm);
15 disp(ec,"Longitudinal Modulus (in GPa) = ");
16 e_c=1/((vf/ktm)+(vm/etm));
17 disp(e_c,"Transverse Modulus (in GPa) = ");
18 kp=0.34; //in GPa
19 ep=0.36; //in GPa
20 vlt=(vf*kp)+(vm*ep);
21 disp(vlt,"Poissons Ratio = ");

```



```
22 glt=1/((vf/2.2)+(vm/1.2)); //in GPa
23 disp(glt," Shear Modulous (in GPa) = ");
```

---

## Chapter 20

# PERFORMANCE OF MATERIALS IN SERVICE

Scilab code Exa 20.1 Find fracture strength and ratio

```
1 //Find fracture strength and ratio
2 //Ex:20.1
3 clc;
4 clear;
5 close;
6 l=1.5*10^-6;//crack length in m
7 e=70*10^9;//Young's modulus in N/m^2
8 y_e=1.05;//specific surface energy in j/m^2
9 a_f=sqrt((2*y_e*e)/(3.14*1));
10 a_f1=a_f*10^-6;//in MPa
11 disp(a_f1,"Fracture strength (in MPa) = ");
12 r=a_f/e;//ratio
13 disp(r,"Ratio of fracture strength to Youngs
    modulus = " );
```

---

Scilab code Exa 20.2 Investigate an oxidation film will form over Nickel or not

```
1 //Investigate an oxidation film will form over
   Nickel or not
2 //Ex:20.2
3 clc;
4 clear;
5 close;
6 m_m=58.71; //molecular weight of ni
7 m_c=74.71; //molecular weight of nio
8 p_m=8900; //density of ni in kg/m^3
9 p_c=7080; //desity of nio in kg/m^3
10 x=m_m/p_m; //molar volume of ni in m^3/mol
11 disp(x,"Mc/Pc (in m^3/mol)");
12 y=m_c/p_c; //molar volume of nio in m^3/mol
13 disp(y,"Mm/Pm (in m^3/mol)");
14 printf("          Mc/Pc > Mm/Pm Hence protective
   layer of NiO will form over Ni          ");
```

---

Scilab code Exa 20.3 Find how much loss will occur in 300 hours

```
1 //Find how much loss will occur in 300 hours
2 //Ex:20.3
3 clc;
4 clear;
5 close;
6 x1=0.1; //in mm
7 t1=25; //in hours
8 t2=300; //in hours
9 x2=x1*sqrt(t2/t1); //in mm
10 disp(x2,"Oxidation loss in 300 hours (in mm) = ");
```

---

#### Scilab code Exa 20.4 Determine PBR ratio

```
1 //Determine PBR ratio
2 //Ex:20.4
3 clc;
4 clear;
5 close;
6 p_mg=1.74; //density of magnesium in gm/cm^3
7 p_mgo=3.65; //density of magnesium oxide in gm/cm^3
8 m_mg=24; //mol wt of mg
9 m_mgo=40.3; //mol wt of mgo
10 PBR=(m_mgo/p_mgo)/(m_mg/p_mg);
11 disp(PBR,"PBR = ");
12 printf("Since PBR < 1. So porous film will form
        which will be non protective");
```

---

#### Scilab code Exa 20.5 Find quantity of magnesium needed

```
1 //Find quantity of magnesium needed
2 //Ex:20.5
3 clc;
4 clear;
5 close;
6 m=0.0243; //one mole of magnesium in kg
7 c=2*96490; //in C
8 j=20*10^-3; //in A/m^2
9 t=15*365*24*3600; //in sec
10 x=j*t; //in A s
```

```
11 w_mg=m*x/c; //in kg/sqm
12 disp(w_mg,"Amount of Magnesium needed (in Kg/m^2) =
    ");
```

---

# Chapter 21

## DIFFUSION IN SOLIDS

Scilab code Exa 21.1 Find Concentration Gradient and diffusivity

```
1 //Find Concentration Gradient and diffusivity
2 //Ex:21.1
3 clc;
4 clear;
5 close;
6 c_cu=2*10^13;//concentration of copper in /m^3
7 c_al=4*10^6;//concn of copper on other side of Al in
  /m^3
8 t=3*10^-3;//thickness in m
9 z=(c_cu-c_al)/t;//z=dm/dx, concentration graient
10 disp(z,"Concentration Gradient (in /m^4) = ");
11 jx=10^21;//outward flux of copperv atoms in /sq m/
  sec
12 d=-jx/z;//diffusivity in sq m/sec
13 disp(d,"Diffusivity (in sq m/sec) = ");
```

---

Scilab code Exa 21.2 Find Concentration Gradient and diffusivity and rate

```

1 //Find Concentration Gradient and diffusivity and
  rate
2 //Ex:21.2
3 clc;
4 clear;
5 close;
6 c_n=12;//nitrogen concentration in kg/m^3
7 t=6*10^-3;//thickness in m
8 z=(c_n-0)/t;//concentration gradient in kg/m^4
9 disp(z,"concentration gradient (in kg/m4) = ");
10 d0=5*10^-7;//in sqm/sec
11 q=75*10^3;//in j/mol
12 r=8.314;//in J/mol/K
13 t=400;//in K
14 dx=d0*exp(-q/(r*t));//diffusivity in sqm/sec
15 disp(dx,"Diffusivity (in sqm/sec) = ");
16 jx=-z*dx;//rate of flow of nitrogen in kg/sqm/sec
17 disp(jx,"Rate at which nitrogen escapes (in kg/sqm/
  sec) = ");

```

---

**Scilab code Exa 21.4 Calculate Activation Energy in Silver diffusion**

```

1 //Calculate Activation Energy in Silver diffusion
2 //Ex:21.4
3 clc;
4 clear;
5 close;
6 z=8;//ratio of diffusion in silicon at 1350 C and
  1100 C
7 x=log(z);
8 q=x/(1.35*10^-5);//activation energy for silver
  diffusion in J/mol
9 q1=q/1000;//in kJ/mol

```

```
10 disp(q1," Activation Energy in Silver diffusion (in  
kJ/mol) = ");
```

---