

Scilab Textbook Companion for
A Text Book of Physical Chemistry
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Book Description

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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 1

Equilibrium between phases

Scilab code Exa 1.5.1 Ex 1

```
1 clear ;
2 clc ;
3 T1 = 234.5 ;// Temperature in K
4 P = 1 ; // Pressure in atm
5 rho1 = 14.19 // Density of solid Hg in g/(cm^3)
6 rho2 = 13.70 // Density of liquid Hg in g/(cm^3)
7 V = 200.59 // volume of liquid and solid in g/mol
8 delV = ((V/rho2)-(V/rho1))*(10^-3) // in dm^3/mol
9 delTdelP = 0.0051 // K/atm
10 R1 = 8.314 // in J
11 R2 = 0.082 // in (dm)^3/atm
12 delH = ((delV*T1)/(delTdelP))*(R1/R2)*10^-3; //molar
    heat of fusion in kJ/mol
13 printf('delH = %.3f (KJ)/mol',delH)
14 T2 = 273 // in K
15 delP = (delH*(R2/R1)*(T2-T1))/(delV*T1)*10^3; //
    pressure required to raise melting point to T2 in
    atm
16 printf('\ndelP = %d atm ',delP)
17
18
```

19 //Example in page 10

Scilab code Exa 1.5.2 Ex 2

```
1 clear;
2 clc;
3 T1=373.15; //in K
4 P=1; //atm
5 Vv=1674; //in cm^3/gm
6 delPdelT=27.12; //in torr/K
7 R1=8.314; //in J
8 R2=0.082; //in atm/(dm)^3
9 delH=((delPdelT)/760)*T1*((Vv*10^-3)*18)*(R1/R2)
10 printf('delH =%d J/mol', delH)
11
12 ////Example in page 15
```

Scilab code Exa 1.5.3 Ex 3

```
1 clear;
2 clc;
3 T1=313.75; //in K
4 P1=59.1; //in torr
5 T2=353.15; //in K
6 P2=298.7; //in torr
7 R=2.303*8.314; //in J/(K*mol)
8 delH=R*log10(P2/P1)*((T2*T1)/(T2-T1))
9 printf('delH=%d J/mol', delH)
```

Scilab code Exa 1.5.4 Ex 4

```

1 clear
2 clc
3 T1=325.15; //in K
4 T2=338.15; //in K
5 P2=760; //in torr
6 DelHm_v=10.5; //
7 P1=P2/(10^((DelHm_v/2.303)*((T2/T1)-1))); //in torr
8 printf('P1=%0.1f torr',P1)
9 P=200; //in torr
10 T=T2/(1+((2.303/10.5)*log10(P2/P))); //in K
11 printf('\nT=%0.1d K',T)
12 I=log10(P2)-(((DelHm_v*T2)/2.303)*(-1/T2)); //
13 printf('\nI=%0.3f',I)
14
15 //There are some errors in the solution given in
    textbook
16 //page 16

```

Scilab code Exa 1.5.5 Ex 5

```

1 clear;
2 clc;
3 P=760; //in torr
4 dP=52; //in torr
5 dT=2; //in K
6 DelH_RTb=10.5; //Trouton rule
7 Tb=(DelH_RTb*P)/(dP/dT)
8 printf('Tb=%0.1f K',Tb)
9 R=8.314; //in J/Kmol
10 DelH_v=(DelH_RTb*R*Tb)
11 printf('\nDelH_v=%01d J/mol',DelH_v)
12
13 //There are some errors in the solution given in
    textbook
14 //page 17

```

Scilab code Exa 1.5.6 Ex 6

```
1 clear;
2 clc;
3 T1=373.15; //in K
4 P1=76.0; //in cmHg
5 P2=77.0; //in cmHg
6 DelHm_v=2255; //in J/gm
7 Vm_v=1664; //in cm^3/mol
8 Vm_l=1; //in cm^3/mol
9 R=8.314; //in J/Kmol
10 T2=(1/((1/T1)-((2.303*R/(DelHm_v*18))*log10(P2/P1))))
11 printf('T2=%0.1f K',T2)
12
13 //There are some errors in the solution given in
14 //page 18
```

Scilab code Exa 1.5.7 Ex 7

```
1 clear;
2 clc;
3 T1=373.15; //in K
4 P1=76.0; //in cmHg
5 T2=363.15 //in K
6 DelHm_v=2268; //in J/gm
7 R=8.314; //in J/Kmol
8 P2=P1*(10^((DelHm_v*18/(2.303*R))*(1/T1-1/T2)))
9 printf('P2=%0.1f cmHg',P2)
10
11 //page 19
```

Scilab code Exa 1.5.8 Ex 8

```
1 clear
2 clc
3 T1=456.15; //boiling temperature of iodine in K
4 T2=389.65; //vapour pressure temperature of iodine in
   K
5 P1=760; //pressure in torr
6 P2=100; //vapour pressure in torr
7 DelHm_f=15.65; //heat of fusion in kJ/mol
8 R=8.314; //in J/K
9 DelHm_v=(2.303*R*log10(P1/P2))/((1/T2)-(1/T1)); //
   heat of vapourization in J/mol
10 DelHm_s=(DelHm_f*1000)+DelHm_v; //heat of sublimation
   in J/mol
11 T=311.85; //temperature at solid vapour equilibrium
   in K
12 P=1; //pressure at solid vapour equilibrium in torr
13 K1=(DelHm_s)/(2.30*R); //
14 K2=(DelHm_v)/(2.30*R); //
15 T0=(K1-K2)/((K1*(1/T))-(K2*(1/T2))-log10(P2)); //
   triple point temperature in K
16 printf('\nT0=%0.1 f K',T0)
17 P0=10^(K1*((1/T)-(1/T0))); //triple point pressure in
   torr
18 printf('\nP0=%0.2 f torr',P0)
19
20 //There are some errors in the solution given in
   textbook
21 //page 19
```

Scilab code Exa 1.5.9 Ex 9

```

1 clear
2 clc
3 K1=5.36; //
4 K2=4.95; //
5 T1=-2875; //in K
6 T2=-2740; //in K
7 R=8.314; //in J/Kmol
8 T=(T2-T1)/(K1-K2); //triple point temperature in K
9 printf('T=%0.1f K',T)
10 P=(10^((T1/T)+K1)); //triple point pressure in atm
11 printf('\nP=%0.7f K atm',P)
12 DelHm_s=2.303*R*(-T1); //molar enthalpy of
    sublimation in J/mol
13 DelHm_v=2.303*R*(-T2); //molar enthalpy of
    vapourization in J/mol
14 DelHm_f=DelHm_s-DelHm_v; //molar enthalpy of fusion
    in J/mol
15 printf('\nDelHm_f=%0.1d J/mol',DelHm_f)
16 DelSm_f=DelHm_f/T; //molar entropy of fusion in J/
    Kmol
17 printf('\nDelSm_f=%0.2f J/Kmol',DelSm_f)
18
19 //There are some errors in the solution given in
    textbook
20 //page 20

```

Scilab code Exa 1.5.10 Ex 10

```

1 clear
2 clc
3 T=273; //in K
4 R=8.314; //in J/Kmol
5 DelHm=T*2.303*R; //in J/mol
6 printf('DelHm=%0.1d J/mol',DelHm)
7

```

Scilab code Exa 1.5.11 Ex 11

```
1 clear;
2 clc;
3 DelG=2866; //in J/mol
4 rhoG=2.25; //in gm/cm^3
5 rhoD=3.52; //in gm/cm^3
6 MC=12; //mass of carbon
7 P1=1; //in atm
8 P2=(-DelG/(MC/rhoD-MC/rhoG))+P1
9 printf('P2=%0.1f Jcm^-3',P2)
10 R1=0.082; //in dm^3atm
11 R2=8.314; //in J
12 P21=P2*(R1*1000/R2)
13 printf('\nP21=%0.1d atm',P21)
14
15 //There are some errors in the solution given in
    textbook
16 //page 23
```

Scilab code Exa 1.5.12 Ex 12

```
1 clear
2 clc
3 rho1=1.21; //in gm^cm-3
4 rho2=1.10; //in gm^cm-3
5 P2=3260; //in atm
6 T2=298.15; //in K
7 P1=2450; //in atm
8 T1=242.15; //in K
9 MI=18; //molar mass of ice in gm/mol
```

```

10 R1=8.314; //in J
11 R2=0.082; //in atm dm^3
12 DelH_Tr=((P2-P1)*(R1/R2)*(MI/rho2-MI/rho1)*T1)/(T2-
    T1)
13 printf('DelH_Tr=%0.3 f J/mol',DelH_Tr)
14
15 //There is an error in the answer given in the
    textbook
16 //In text book he took (T1-T2)=6,but actually (T1-T2
    )=56
17 //page 23

```

Scilab code Exa 1.5.13 Ex 13

```

1 clear
2 clc
3 DelVm_tr=0.0126; //in cm^3/gm
4 P=1; //in atm
5 Ti=368.65; //in K
6 DelTDelP=0.035; //in K/atm
7 R1=8.314; //in J
8 R2=0.082; //in dm^3atm
9 DelHm_tr=Ti*(DelVm_tr*32/1000)*1/(DelTDelP)*(R1/R2)
10 printf('DelHm_tr=%0.1 f J/mol',DelHm_tr)

```

Scilab code Exa 1.7.1 Ex 14

```

1 clear
2 clc
3 T=263.15; //in K
4 P2=1.95; //in torr
5 rho=0.920; //in gm/cm^3
6 P=1; //in atm

```



```
7 R=0.082; //in dm^3atm/(molK)
8 P1=P2*exp((18/(rho*1000))*(P-(P2/760))/(R*T))
9 printf('P1=%0.3f torr ',P1)
10
11 //page 29
```

Scilab code Exa 1.7.2 Ex 15

```
1 clear
2 clc
3 P0=100; //in atm
4 P=1; //in atm
5 P2=31.82; //in torr
6 rho=0.996; //in gm/cm^3
7 R=0.082; //in dm^3atm/(molK)
8 T=303.15; //in K
9 P1=P2*(10^(((18/(rho*1000))*(P0-P))/(2.303*R*T)))
10 printf('P1=%0.1f torr ',P1)
11
12 //page 29
```

Chapter 2

Colligative properties

Scilab code Exa 2.2.1 Ex 1

```
1 clear
2 clc
3 M1=20; //mass of acetic acid in gm
4 M2=80; //mass of water in gm
5 mM=60; //molar mass of acetic acid in gm
6 Vm1=M1/60; //in mol
7 Vm2=M2/18; //in mol
8 rho=1.026; //in gm/cm^3
9 X=Vm1/(Vm1+Vm2); //mole fraction of acetic acid
10 printf('X=%0.3f',X)
11 B=Vm1/(M2/1000); //molality of acetic acid
12 printf('\nB=%0.3f mol/kg',B)
13 V=(M1+M2)/rho
14 C=(Vm1)/(V/1000); //molarity of acetic acid
15 printf('\nC=%0.3f moldm^-3',C)
16
17 //There are some errors in the solution given in
    textbook
18 //In textbook the value of X is given in fraction
19 //page 36
```

Scilab code Exa 2.2.2 Ex 2

```
1 clear
2 clc
3 C=5; //molarity in mol
4 mM=100; //molar mass in gm
5 rho=1.289; //in gm/cm^3
6 M1=C*mM; //mass of solute
7 M2=(rho*1000)-M1; //mass of solvent
8 V=(M2)/18; //volume of water solvent in mol
9 X=(C)/(V+C); //mole fraction of solute
10 printf('X=%0.4 f',X)
11 B=(C)/(M2/1000)
12 printf('\nB=%0.3 f mol/kg',B)
13
14 //page 23
```

Scilab code Exa 2.4.1 Ex 3

```
1 clear
2 clc
3 T=303; //in K
4 m2=10; //mass of solute in gm
5 m1=80; //mass of solute acetone in gm
6 P1=271; //in torr
7 P2=283; //in torr
8 M1=58; //in gm/mol
9 M2=((m2*M1)/(((P2-P1)/P2)*m1))-((M1*m2)/m1)
10 printf('M2=%0.1 f gm/mol',M2)
11
12 //page 39
```

Scilab code Exa 2.4.2 Ex 4

```
1 clear
2 clc
3 P1=74.01; //in torr
4 P2=74.66; //in torr
5 m2=2; //in gm
6 m1=100; //in gm
7 M1=78; //in gm
8 M2=((m2*M1)/(((P2-P1)/P2)*m1))-((M1*m2)/m1)
9 printf('M2=%0.1f gm/mol',M2)
10 nCH=94.4/5.6; //mass ratio of C and H
11 N=nCH*(1/12); //atomic ratio
12 printf('\nN=%0.1f',N)
13
14 //atomic ratio is 7:5 (here N is showed decimals)
15
16 EM=(12*7)+(1*5); //empirical mass
17 K=M2/EM; //No. of units C7H5
18 printf('\nK=%0.1f',K)
19
20 //Approximately equal to 2, Molecular Formula C14H10
21 //There are some errors in the solution given in
    textbook
22 //page 40
```

Scilab code Exa 2.4.3 Ex 5

```
1 clear
2 clc
3 m1=100; //amount of water in gm
4 M1=18; //in gm
```

```

5 m2=1; //amount of urea in gm
6 M2=60; //in gm
7 m3=2; //amount of sucrose in gm
8 M3=342; //in gm
9 X=(m1/M1)/((m1/M1)+(m2/M2)+(m3/M3)); //mole fraction
    of solvent
10 P2=23.756; //in torr
11 T=298; //in K
12 P1=P2*X; //vapour pressure of solution intorr
13 printf('P1=%0.2f torr',P1)
14
15 //There are some errors in the solution given in
    textbook
16 //page 39

```

Scilab code Exa 2.7.3 Ex 6

```

1 clear
2 clc
3 m1=0.5126; //in dissolved mass in gm
4 mM1=128.2; //molar mass of naphthalene in gm
5 m0=50; //mass of solvent in gm
6 B=(m1/mM1)/(m0/1000); //Molality of solution in mol/
    kg
7 printf('B=%0.5f mol/kg',B)
8 delTb=0.402; //change in temperature of naphthalene
    in K
9 Kb=delTb/B;
10 delTbs=0.647 //change in temp for unknown solution in
    K
11 m2=0.6216; //mass of unknown solute
12 M=(Kb*m2*1000)/(delTbs*m0); //molar mass of unknown
    solute
13 printf('\nM=%0.2f gm/mol',M)
14

```

Scilab code Exa 2.7.4 Ex 7

```
1 clear
2 clc
3 P0=100; //vapour pressure in torr
4 P2=760; //in torr
5 T2=353.15; //in K
6 T1=300.15; // in K
7 DelSm_v=87.03; //entropy in J/Kmol
8 R=8.314; //in J/Kmol
9 P1=P2/(10^((DelSm_v*(T2-T1))/(2.303*R*T1)))
10 printf('P1=%0.1f torr',P1)
11 X=(P1-P0)/P1; //Mole fraction of solute
12 printf('\nX=%0.4f',X)
13 T0=(1/T2)+((R*log(1-X))/(DelSm_v*T2))
14 Tb=1/T0; //Boiling point of solution
15 printf('\nTb=%0.1f K',Tb)
16
17 //page 52
```

Scilab code Exa 2.7.5 Ex 8

```
1 clear
2 clc
3 M1=76; //molar mass of CS2 in gm
4 w2=3.795; //weight of S in 100gm of CS2 in gm
5 w1=100; //weight of CS2
6 R=8.314; //in J/Kmol
7 Tb=319.81; //boiling point of CS2 in K
8 Tbp=319.45; //boiling point of pure CS2 in K
9 DelHm_v=351.87; //enthalpy of vaporization in J/gm
```

```

10 M2=(w2*M1*R*(Tb^2))/(w1*(Tb-Tbp)*DelHm_v*76)
11 printf('M2=%0.1f gm/mol',M2)
12 N=M2/32;//no. of s atoms
13 printf('\nN=%0.1f',N)
14
15 //Molecular formula S8
16 //There are some errors in the solution given in
    textbook
17 //page 53

```

Scilab code Exa 2.8.2 Ex 9

```

1 clear
2 clc
3 M1=152.2;//molar mass of carbon in gm
4 T1=451.55;//melting point temp in K
5 T2=433.85;//melting point temp in K(for unknown
    compound)
6 w2=0.0386;//mass of unknown compound in gm
7 w1=0.522;//mass of camphor in solution in gm
8 R=8.314;//J/Kmol
9 DelHm_f=6.844;//in KJ
10 Kf=((R*T1^2)/(DelHm_f*10^3))*(M1/1000)
11 printf('Kf=%0.1f',Kf)
12 DelT_f=(T1-T2);
13 B=(DelT_f/Kf);//molality of the solution in mol/kg
14 printf('\nB=%0.2f mol/kg',B)
15 M2=(w2/B)*(1000/w1);
16 printf('\nM2=%0.1f gm/mol',M2)
17 Z=92.3/7.7;//mass ratio of wC and wH
18 N0=1/12;//atomic ratio of H and C
19 K=(Z*N0);
20 printf('\nK=%0.1f',K)
21
22 //Clearly we get K=1.0 implies empirical formula is

```

```

      CH
23 Me=13; //empirical mass in gm
24 N=(M2/Me); //no. of units of CH
25 printf( '\nN=%0.1 f ',N)
26
27 //Taking approximately equal to 12 Molecular formula
    is C12H12
28 //There are some errors in the solution given in
    textbook
29 //page 58

```

Scilab code Exa 2.8.3 Ex 10

```

1 clear
2 clc
3 n1=0.1; //amount of naphthalene in mol
4 n2=0.9; //amount of benzene in mol
5 Tf=278.5; //freezing temperature of C6H6 in K
6 Tb=353; //boiling temperature of C6H6 in K
7 P1=670; //vapour pressure in torr
8 P2=760; //in torr
9 R=8.314; //in J/Kmol
10 M1=78; //atomic mass of C6H6
11 DelHm_f=10.67; //in KJ
12 X1=(P2-P1)/P2; //
13 nT=(n1/X1); //
14 nb=(nT-n1); //
15 Kfb=((R*Tf^2)/(DelHm_f*1000))*(M1/1000); //
16 printf( 'Kfb=%0.3 f Kg/mol ',Kfb)
17 B=(n1/(nb*M1)*1000); //molality of the solution
18 printf( '\nB=%0.3 f mol/kg ',B)
19 DelTf=(Kfb*B); //in K
20 T=(Tf-DelTf); //in K
21 printf( '\nT=%0.2 f K ',T)
22

```



```
23 //There are some errors in the solution given in
    textbook
24 //page 59
```

Scilab code Exa 2.9.1 Ex 11

```
1 clear
2 clc
3 x1=[0.0200,0.0150,0.0100,0.0075,0.0050,0.00025]
4 y2=[0.104,0.101,0.099,0.098]
5 x2=[0.0200,0.0150,0.0100,0.0050]
6 y1=[0.585,0.440,0.300,0.230,0.18,0.140]
7 plot(x1,y1,'go-',x2,y2,'ro-')
8 [m1,c1]=reglin(x1,y1)
9 [m2,c2]=reglin(x2,y2)
10 R=82.0; //in cm^2atm/Kmol
11 T=298; //in K
12 M=R*T/c2; //molar mass of polyisobutylene in gm/mol
13 printf('M=%0.1d gm/mol',M)
14
15 //There is some error in the solution given in
    textbook
16 //There are some errors in the solution given in
    textbook
17 //page 68
```

Scilab code Exa 2.9.2 Ex 12

```
1 clear
2 clc
3 h=3.9; //height in mm
4 rho=1.0; //density of solution
5 g=980.7; //acceleration due to gravity in cm/s^2
```

```

6 P=((h/10)*rho*g); //osmotic pressure in gm/(cms^2)
7 V=1000; //volume in cm^3
8 T=25; //temperatur in C
9 w2=1; //weight of serum albumin
10 R=8.314; //in J/Kmol
11 M2=(w2*(R*10^7)*(T+273))/(P*V); //molar mass of serum
    albumin
12 printf('M2=%0.3f *10^4 g/mol',M2/10^4)
13
14 //The above result is in CGS units
15
16 //The following results are in SI units
17 p=(P/10); //osmotic pressure in N/m^2
18 m2=(M2/10^3); //molar mass of serum albumin
19 printf('\nm2=%0.2f Kg/mol',m2)
20
21 //page 67

```

Scilab code Exa 2.9.3 Ex 13

```

1 clear
2 clc
3 w1=1; //amount of glucose C6H12O6 in gm
4 w2=1; //amount of sucrose C12H12O22 in gm
5 n=(w1/180)+(w2/342); //amount of solute
6 R=8.314; //in J/Kmol
7 T=25; //in C
8 V=1000; //volume of water in gm
9 P=(n*R*(T+273))/(V*10^-6); //osmotic pressure of
    solution
10 printf('P=%0.3f *10^4 N/m^2',P/10^4)
11 w=(w1+w2); //weight of solute
12 M=(w*R*(T+273))/(P*(V*10^-3)); //molar mass of solute
13 printf('\nM=%0.4f kg/mol',M)
14 Mn=((w1*10^-3)+(w2*10^-3))/(n); //average molar mass

```

```
    in Kg/mol
15 printf( '\nMn=%0.4 f kg/mol ', Mn)
16
17 //page 67
```

Scilab code Exa 2.10.1 Ex 14

```
1 clear
2 clc
3 P=2.47; //osmotic pressure in atm
4 DelHm_v=539*18; //in cal/mol
5 R=0.082; //in litreatm
6 Vm=18.1; //molar volume of water
7 T=303; //in K
8 Tb=373; //boiling point temperature in K
9 DelTb=(P*Vm*10^-3*(Tb^2))/(DelHm_v*(R/1.987)*T)
10 printf( 'DelTb=%0.4 f K', DelTb)
11
12 //The above calculations are done in CGS units
13
14 //To convert them into SI units the following
    changes are done
15 R=8.314; //in J/Kmol
16 P=2.47*101325; //in N/m^2
17 Vm=18.1*10^-6; //in m^3/mol
18
19 //Both answers come out be same
20 //page 70
```

Scilab code Exa 2.10.2 Ex 15

```
1 clear
2 clc
```

```

3 rho=1.59; //density of CCl4 in kg/dm^3
4 M1=154; //molar mass of CCl4 in kg/mol
5 DelTb=0.60; //boiling point of CCl4 in K
6 Kb=5.03; //in Kkg/mol
7 m=DelTb/Kb;
8 m2=3; //amount added to CCl4 in gm
9 m1=100; //amount of CCl4 in gm
10 M2=(m2*10^-3)/(m1*10^-3*m); //molar mass of substance
11 printf('M2=%0.3 f kg/mol',M2)
12 Kf=31.8; //freezing point depression in Kkg/mol
13 DelTf=Kf*m
14 printf('\nDelTf=%0.3 f K',DelTf)
15 P=(m2*10^-3/M2)/(((m2*10^-3)/M2)+((m1*10^-3)/(M1
    *10^-3))); //relative vapour pressure DelP/P1
16 printf('\nP=%0.5 f',P)
17 V1m=m1*10^-3/(rho); //volume in dm^3
18 R=0.082; //in dm^3atm/Kmol
19 T=298; //Temperature in K
20 P0=((m2/250)*R*T)/V1m; //osmotic pressure in atm
21 printf('\nP0=%0.3 f atm',P0)
22
23 //There are some errors in the solution given in
    textbook
24 //page 71

```

Scilab code Exa 2.11.1 Ex 16

```

1 clear
2 clc
3 w2=0.122; //amount of benzoic acid in kg
4 w1=1; //amount of benzene in kg
5 Tb1=353; //boiling point of benzene in K
6 Tb2=354.5; //boiling point at which actually boiling
    of benzene starts in K
7 DelH_v=394.57; //in J/gm

```

```

8 M1=w2/0.122;//amount of benzoic acid in mol
9 R=8.314;//in J/Kmol
10 M2=((M1*78*10^-3)*R*Tb1^2*w2)/(w1*(Tb2-Tb1)*(DelH_v
    *78));//apparant molar mass of benzoic acid in kg
    /mol
11 printf('M2=%0.4f kg/mol',M2)
12 alpha=2*(1-(w2/M2));//degree of dimerisation in mol
13 printf('\nalpha=%0.4f mol',alpha)
14
15 //There are some errors in the solution given in
    textbook
16 //page 75

```

Scilab code Exa 2.11.2 Ex 17

```

1 clear
2 clc
3 w2=0.011;//amount of barium nitrate in kg
4 M2=0.2613;//molar mass of barium nitrate inkg/mol
5 w1=0.1;//amount of water in kg
6 Kb=5.2;//for 100gm of water in K
7 m=(w2/M2)/w1;//molality of solution in mol/kg
8 DelTb_0=Kb/10*m;//in K
9 T=100.46;//boiling point of water
10 i=(T-100)/DelTb_0;//van't hoff factor
11 v=3;
12 alpha=(i-1)/(v-1);//degree of ionization
13 printf('alpha=%0.2f',alpha)
14
15 //page 75

```

Scilab code Exa 2.11.3 Ex 18

```

1 clear
2 clc
3 M1=324.6; //molar mass of Hg(NO3)2 in gm
4 m1=3.24; //amount of Hg(NO3)2 in gm dissolved in
   water
5 w=1; //amount of water in kg
6 M0=(m1/M1)*(1/w); //molality of the solution in K
7 Kf=1.86; //in Kkg/mol
8 DelTf_0=(Kf*M0); // here DelTf_0 is negative
9 DelTf=0.0558; //freezing point of the solution here
   DelTf is negative
10 i=(DelTf/DelTf_0); //van't hoff factor
11 v=3;
12 alpha=(i-1)/(v-1); //degree of dissociation
13 printf('alpha=%1d',alpha)
14 M2=271.5; //molar mass of HgCl2 in gm
15 m2=10.84; //amount of HgCl2 dissolved in water in gm
16 M=(m2/M2)*(1/w); //molality of HgCl2 solution in mol/
   kg
17 DelTf1_0=Kf*M; //for HgCl2 solution
18 printf('\nDelTf1_0=%0.3 f K',DelTf1_0)
19
20 //page 76

```

Scilab code Exa 2.11.4 Ex 19

```

1 clear
2 clc
3 M=78*10^-3; //molar mass of C6H6 in Kg/mol
4 R=8.314; //gas constant in J/Kmol
5 Tf2=278.4; //melting point of pure C6H6 in K
6 DelHm_v=10.042*10^3; //heat of fusion in J/mol
7 Kf=((M*R*Tf2^2)/DelHm_v); //inKkg/mol
8 Tf1=277.4; //melting point of C6H6 in Kg/mol
9 M1=(Tf2-Tf1)/Kf; //molality in mol/kg

```

```

10 X1=0.02; //molefraction of CH3COOH
11 M2=X1/M; //molality in mol/kg
12 Md=(M2-M1); //molality of dimer in mol/kg
13 Mm=M1-Md; //molality of monomer in mol/kg
14 Keq=(Md)/(Mm^2); //equilibrium constant fo
    dimerization of CH3COOH
15 printf( 'Keq=%0.2 f ', Keq)
16
17 //There are some errors in the solution given in
    textbook
18 //page 76

```

Scilab code Exa 2.11.5 Ex 20

```

1 clear
2 clc
3 DelTf1=0.704; //freezing point of aqueous KCN in K
4 Kf=1.86; //in kg/mol
5 M1=(DelTf1)/Kf; //molality of the solution containing
    KCN
6 DelTf2=0.530; //freezing point on addition of Hg(CN)2
7 M2=(DelTf2)/Kf; //molality on addition og Hg(CN)2
8 Kplus=0.1892; //amount of K+ in 1000gm of solvent
9 HgCN2=0.095; //amount of Hg(CN)2 added to form
    complex
10 M=(Kplus+Kplus+HgCN2-M2); //
11 N=(M/HgCN2); //no. of CN- units combined
12 printf( 'N=%0.1 f ', N)
13
14 //Formula is Hg(CN)2^-4
15 //page 77

```

Scilab code Exa 2.11.6 Ex 21

```

1 clear
2 clc
3 M1=148.31; //molar mass of Mg(NO2)2 in gm
4 m1=6.69; //amount of Mg(NO2)2 dissolved in water
5 m2=100; //amount of water
6 P1=747; //pressure in torr
7 T2=373; //temperature in K
8 P=760; //pressure at normal temperature
9 X=(m1/M1)/((m1/M1)+(m2/18)); //mole fraction of
    solute in solution \
10 DelP=X*P; //
11 i=(P-P1)/DelP; //van't hoff factor
12 v=3; //
13 alpha=(i-1)/(v-1); //degree of dissociation of salt
    in solution
14 printf('alpha=%0.3f',alpha)
15
16 //page 78

```

Scilab code Exa 2.11.7 Ex 22

```

1 clear
2 clc
3 C_Mg=0.5; //concentration of Mg2+ ion
4 C_SO4=0.7; //concentration of SO4_2- ion
5 C_Al=0.1; //concentration of Al3+ ion
6 C_Cl=0.3; //cocncentration of Cl- ion
7 C_NH4=0.4; //concentration of NH4+ ion
8 Z1=2; //valence of Mg2+ ion
9 Z2=2; //valence of SO4_2- ion
10 Z3=3; //valence of Al3+ ion
11 Z4=1; //valence of Cl- ion
12 Z5=1; //valence of NH4+ ion
13 mu=1/2*(C_Mg*(Z1^2)+C_SO4*(Z2^2)+C_Al*(Z3^2)+C_Cl*(
    Z4^2)+C_NH4*(Z5^2)); //ionic strength

```



```
14 printf('mu=%.1 f',mu)
15
16 //page 78
```

Scilab code Exa 2.11.8 Ex 23

```
1 clear
2 clc
3 Kf=1.86; //in Kkg/mol
4 m=0.2; //amount of aqueous solution of KCL freezes in
      mol/kg
5 DelTf_0=Kf*m; //in K
6 DelTf_1=0.680; //in K
7 i1=DelTf_1/DelTf_0; //van't hoff factor
8 printf('i1=%.2 f',i1)
9 v=2;
10 alpha=(i1-1)/(v-1); //degree of dissociation
11 printf('\nalpha=%.2 f',alpha)
12 z=1; //valency
13 mu=(1/2)*((m*z^1)+(m*z^1))
14 printf('\nmu=%.1 f',mu)
15 i2=v*(1-((0.375+z-z)*(sqrt(mu))))
16 printf('\ni2=%.4 f',i2)
17 Kb=0.52; //in Kkg/mol
18 DelTb=i1*Kb*m
19 printf('\nDelTb=%.3 f K',DelTb)
20 R=8.314; //in J/Kmol
21 T=273; //in K
22 P=i1*(m*10^3)*R*T*(1/101325); //osmotic pressure in
      atm
23 printf('\nP=%.1 f atm',P)
24
25 //page 79
```

Scilab code Exa 2.12.1 Ex 24

```
1 clear
2 clc
3 DelH2m_f=10; //molar heat of fusion in kJ/mol
4 T1=298.15; //temperature in K
5 T2=353.35; //freezing temperature in K
6 R=8.314; //in J/Kmol
7 X=(10^-(((DelH2m_f*10^3)/R)*((1/T1)-(1/T2)))); //
   solubility of naphthalene
8 printf('X=%0.4f',X)
9
10 //solution is wrong
11 //There are some errors in the solution given in
   textbook
12 //page 81
```

Scilab code Exa 2.12.2 Ex 25

```
1 clear
2 clc
3 n2=6.2; //in mol
4 n1=1000; //in mol
5 X=(n2)/((n1/18)+n2); //solubility of sucrose
6 T1=298; //in K
7 T2=473; //freezing point temperature in K
8 R=8.314; //in J/Kmol
9 DelH2m_f=-((R*2.303*log10(X))/((1/T1)-(1/T2))); //molar
   heat of fusion
10 printf('DelH2m_f=%0.1d J/mol',DelH2m_f)
11
```

- 12 //There are some errors in the solution given in
textbook
 - 13 //page 82
-

Chapter 3

Phase Rule

Scilab code Exa 3.3.1 Ex 1

```
1 clear
2 clc
3 //In KCL NaCl H2O system
4 r=3;//no.of reactions
5 C=8;//no. of constituents
6 Z=2;//no.of restricting equations
7 C1=C-r-Z;//no. of components
8 printf('C1=%0.1d',C1)
9
10 //If salts present in equal amounts
11 C1=C-r-(Z+1);//no. of components
12 printf('\nC1=%0.1d',C1)
13
14 //If KCL NaCl as strong electrolytes
15 r=1;//no.of reactions
16 C=6;//no. of constituents
17 Z=2;//no.of restricting equations
18 C3=C-r-Z;//no. of components
19 printf('\nC3=%0.1d',C3)
20
21 //If salts present in equal amounts
```

```

22 C4=C-r-(Z+1); //no. of components
23 printf( '\nC4=%0.1d', C4)
24
25 //In KCL NaCl H2O system
26 r=5; //no. of reactions
27 C=11; //no. of constituents
28 Z=2; //no. of restricting equations
29 C5=C-r-Z; //no. of components
30 printf( '\nC5=%0.1d', C5)
31
32 //If salts present in equal amounts
33 C6=C-r-(Z+1); //no. of components
34 printf( '\nC6=%0.1d', C6)
35
36 //If KCL NaCl NaBr and KBr as strong electrolytes
37 r=1; //no. of reactions
38 C=7; //no. of constituents
39 Z=2; //no. of restricting equations
40 C7=C-r-Z; //no. of components
41 printf( '\nC7=%0.1d', C7)
42
43 //If salts present in equal amounts
44 C8=C-r-(Z+1); //no. of components
45 printf( '\nC8=%0.1d', C8)
46
47 //page 103

```

Scilab code Exa 3.3.2 Ex 2

```

1 clear
2 clc
3 //For system when P_NH3=P_HCl
4 r=1; //no. of equations
5 C=3; //no. of constituents
6 Z1=1; //no. of restricting equations

```

```

7 C1=C-r-Z1;//no. of components
8 printf('C1=%0.1d',C1)
9
10 //For system when P_NH3 not equal P_HCl
11 Z2=0;//no. of restricting equations
12 C1=C-r-Z2
13 printf('\nC1=%0.1d',C1)
14
15 //page 103

```

Scilab code Exa 3.3.3 Ex 3

```

1 clear
2 clc
3 C=9;//no. of constituents
4 r=5;//no. of equilibrium reactions
5 Z=1;//no. of restricting conditions
6 C1=C-r-Z;//no. of components
7 printf('C1=%0.1d',C1)
8
9 //page 103

```

Scilab code Exa 3.3.4 Ex 4

```

1 clear
2 clc
3 //Arbitrary amounts of A1 and A2 only
4 C=4;//no. of constituents
5 r=1;//no. of reactions
6 Z1=1;//no. of restrictions
7 C1=C-r-Z1;//no. of components
8 printf('C1=%0.1d',C1)
9

```

```
10 //Arbitrary amounts of A1,A2,A3,A4
11 Z2=0
12 C1=C-r-Z2;//no. of components
13 printf( '\nC1=%0.1d',C1)
14
15 //Different moles of A1 and A2 only
16 Z3=2
17 C1=C-r-Z3;//no. of components
18 printf( '\nC1=%0.1d',C1)
19
20 //page 103
```

Chapter 4

Solutions

Scilab code Exa 4.6.1 Ex 1

```
1 clear
2 clc
3 T1=273; //in K
4 T2=283; //in K
5 R=8.314; //in J/Kmol
6 alpha1=0.04889; //absorption coefficients in /atm
7 alpha2=0.03802; //absorption coefficients in /atm
8 DelH=(2.303*R*log10(alpha2/alpha1))/((1/T1)-(1/T2));
   //enthalpy of solution
9 printf('DelH=%0.1d J/mol', DelH)
10
11 //Answer comes negative, error in the textbook
12 //page 118
```

Scilab code Exa 4.6.2 Ex 2

```
1 clear
2 clc
```



```

3 V1=500; //volume of H2O in cm^3
4 V2=15.03; //volume of CH4 in cm^3
5 V=V2/V1; //volume dissolved in 1 cm^3 water
6 P=1; //pressure in atm
7 T=273; //Temperature in K
8 R=82.06; //In cm^3atm/Kmol
9 X=(P*V)/(R*T); //amount of gas dissolved in mol
10 M=(X*16); //mass of gas dissolved in gm
11 K=M/P; //
12 m1=0.001; //amount of CH4 in mol
13 m2=300; //amount of H2O in cm^3
14 M1=(m1*16)/m2; //mass of gas dissolved in 1 cm^3
15 P0=M1/K; //pressure if Henry's law holds in atm
16 printf('P0=%0.3 f atm',P0)
17
18 //There are some errors in the solution given in
    textbook
19 //page 58

```

Scilab code Exa 4.6.3 Ex 3

```

1 clear
2 clc
3 Kh=150; //Henry's law constant in torr
4 X1=0.12; //mole fraction of acetone
5 P=(Kh*X1); //vapour pressure of acetone in torr
6 printf('P=%0.1d torr',P)
7 Kh1=175; //Henry's law constant for chloroform in
    torr
8 X2=(P/Kh1);
9 printf('\nX2=%0.3 f',X2)
10
11 //page 119

```

Scilab code Exa 4.6.4 Ex 4

```
1 clear
2 clc
3 X1=4/100; // amount of NH3 solution
4 X2=(1-X1); //amount of water]
5 P=17; //vapour pressure of pure water
6 PT=50; //total pressure in torr
7 P2=(P*X2); //vapour pressure of water in torr
8 P1=(PT-P2); //vapour pressure of NH3 in torr
9 Kh=P1/X1; //Henry's constant for NH3 in torr
10 X=5/100; //mol % of solution
11 P10=Kh*X; //pressure of NH3 at 5% mol in torr
12 printf('P10=%0.1f torr',P10)
13 P20=P*(1-X); //pressure of water at 5% mol in torr
14 printf('\nP20=%0.1f torr',P20)
15 PT0=(P10+P20); //total pressure for 5% of mol
    solution in torr
16 printf('\nPT0=%0.1f torr',PT0)
17
18 //There are some errors in the solution given in
    textbook
19 //page 119
```

Scilab code Exa 4.6.5 Ex 5

```
1 clear
2 clc
3 P1=2/100*101325; //partial pressure of O2 in Pa
4 P2=8/100*101325; //partial pressure of N2 in Pa
5 Kh1=2.53*10^9; //Henry's law constant for O2 in Pa
6 Kh2=5.47*10^9; //Henry's law constant for N2 in Pa
```

```

7 X1=(P1/Kh1); //mole fraction of O2
8 X2=(P2/Kh2); //mole fraction of N2
9 K=(X1/X2);
10 P=1; //in atm
11 M1=(K/(P+K))*100; //mol % of O2
12 printf('M1=%0.2 f ',M1)
13 M2=100-M1; //mol % of N2
14 printf('\nM2=%0.2 f ',M2)
15 X=X1+X2; //total mole fraction
16 N=X*(1000/18); //in mol
17 Kf=1.86; //Kkg/mol
18 DelTf=(Kf*N); //freezing point of saturated water in
    K
19 printf('\nDelTf=%0.5 f ',DelTf)
20
21 //Freezing point will be negative of DelTf
22 //There are some errors in the solution given in
    textbook
23 //page 120

```

Scilab code Exa 4.7.3 Ex 6

```

1 clear
2 clc
3 Xt=1/2; //mole fraction of toluene
4 Xb=1/2; //mole fraction of benzene
5 Pt=4.274; //Partial pressure of toluene in kNm^2
6 Pb=13.734; //Partial pressure of benzene in kNm^2
7 P=(Xt*Pt)+(Xb*Pb); //total pressure in kNm^2
8 printf('P=%0.4 f kNm^2 ',P)
9 Yt=(Xt*Pt)/P; //composition of toluene
10 printf('\nYt=%0.4 f ',Yt)
11 Yb=(1-Yt); //composition of benzene
12 printf('\nYb=%0.4 f ',Yb)
13 P0=(Pb*Pt)/(Pt+((Pb-Pt)*Xt)); //pressure at which

```

```

    last trace liquid disappear
14 printf( '\nP0=%0.3 f kNm^2 ',P0)
15 Xt1=(Xt*P0)/Pt;//composition of last trace of
    toluene
16 printf( '\nXt1=%0.4 f kNm^2 ',Xt1)
17 Xb1=(1-Xt1);//composition of last trace of benzene
18 printf( '\nXb1=%0.4 f kNm^2 ',Xb1)
19 P=sqrt(Pt*Pb);//pressure when 1 mol of mixture is
    vaporized in kN/m^2
20 printf( '\nP=%0.3 f kN/m^2 ',P)
21 Yb1=1-((P-Pt)/(Pb-Pt));//composition of benzene when
    1 mol of mixture is vaporized
22 printf( '\nYb1=%0.3 f ',Yb1)
23 Yt1=(1-Yb1);//composition of toluene when 1 mol of
    mixture is vaporized
24 printf( '\nYt1=%0.3 f ',Yt1)
25
26 //There are some errors in the solution given in
    textbook
27 //page 143

```

Scilab code Exa 4.7.4 Ex 7

```

1 clear
2 clc
3 XA=0.70;
4 YA=0.35;
5 P=600;//in torr
6 PA=(YA*P)/XA;//vapour pressure of pure A
7 printf( 'PA=%0.1d torr ',PA)
8 PB=((1-YA)*P)/(1-XA);//vapour pressure of pure B
9 printf( '\nPb=%0.1 f torr ',PB)
10
11 //page 145

```

Scilab code Exa 4.7.5 Ex 8

```
1 clear
2 clc
3 PA=54.4; //vapour pressure of n-hexane in kN/m^2
4 PB=18.8; //vapour pressure of n-heptane in kN/m^2
5 YA=0.85; //molar fraction of n-hexane
6 XA=(YA*PB)/(PA-((PA-PB)*YA)); //mole fraction of n-
    hexane in equilibrium with vapour
7 printf('XA=%0.3f',XA)
8 R=8.314; //J/Kmol
9 DelS_mix=(-((XA*2.303*log10(XA))+((1-XA)*2.303*log10
    (1-XA))))*R
10 printf('\nDelS_mix=%0.3f J/K',DelS_mix)
11
12 //page 146
```

Scilab code Exa 4.7.6 Ex 9

```
1 clear
2 clc
3 P1=36.7; //vapour pressure of pure toluene in torr
4 P2=118.2; //vapour pressure of pure benzene in torr
5 nt=50; //% amount of toluene in gm
6 nb=50; //% amount of benzene in gm
7 Nt=92; //molar mass of toluene in gm/mol
8 Nb=78; //molar mass of benzene in gm/mol
9 Xt=(nt/Nt)/((nt/Nt)+(nb/Nb)); //mole fraction of
    toluene
10 Pt=Xt*P1; //partial pressure of toluene on torr
11 Pb=(1-Xt)*P2; //partial pressure of benzene on torr
12 P=Pt+Pb; //total pressure of toluene on torr
```

```

13 printf('P=%0.2f torr',P)
14 Yt=Pt/P;//mole fraction of toluene in vapour phase
15 printf('\nYt=%0.3f ',Yt)
16 Yb=(1-Yt);//mole fraction of benzene in vapour phase
17 printf('\nYb=%0.3f ',Yb)
18 P0=50;//in torr
19 Xt1=(P0-P2)/(P1-P2);//mole fraction of toluene at P
    =50 torr
20 printf('\nXt1=%0.4f ',Xt1)
21 Xb1=(1-Xt1);//mole fraction of benzene at P=50 torr
22 printf('\nXb1=%0.4f ',Xb1)
23
24 //page 146

```

Scilab code Exa 4.7.7 Ex 10

```

1 clear
2 clc
3 PA=22.93;//vapour pressure of pure ethyl bromide in
    kNm-2
4 PB=16.93;//vapour pressure of pure propylene bromide
    in kNm-2
5 nA=3;//in mol
6 nB=2;//in mol
7 P=20.4;//in kNm-2
8 XA=(P-PB)/(PA-PB);//mole fraction of ethyl bromide
9 printf('XA=%0.3f ',XA)
10 XB=(1-XA);//mole fraction of propylene bromide
11 printf('\nXB=%0.3f ',XB)
12 YA=(XA*PA)/P;
13 printf('\nYA=%0.4f ',YA)
14 NA=(nA-(XA*(nA+nB)))/(1-(XA/YA));//amount of
    vaporized ethyl bromide at P
15 printf('\nNA=%0.4f ',NA)
16 NB=(NA/YA)-NA;//amount of vaporized propylene

```

```

    bromide at P
17 printf( '\nNB=%0.4 f ', NB)
18 //There are some errors in the solution given in
    textbook
19 //page 147

```

Scilab code Exa 4.7.8 Ex 11

```

1 clear
2 clc
3 YA=0.6497; //
4 XA=0.578; //
5 nA=3; //
6 nB=2; //
7 N1=(YA-(nA/(nA+nB)))/((nA/(nA+nB))-XA); //amount of
    liquid phase
8 N2=(1/(1+N1))*(nA+nB); //amount of vapour phase
9 NA=YA*((nA+nB)/(1+N1)); //mole fraction of ethyl
    bromide at P
10 printf( '\nNA=%0.4 f mol ', NA)
11 NB=(1-YA)*((nA+nB)/(1+N1)); //
12 printf( '\nNB=%0.4 f mol ', NB)
13
14 //page 148

```

Scilab code Exa 4.7.9 Ex 12

```

1 clear
2 clc
3 PA=300; //in torr
4 PB=800; //in torr
5 YA=0.25;

```

```

6  XA=(YA*PB)/(PA-((PA-PB)*YA)); //mole fraction of
    component A
7  printf('XA=%0.4 f ',XA)
8  XB=(1-XA)
9  P=(PA*XA)+(PB*XB); //total pressure P in torr
10 printf('\nP=%0.1 f torr ',P)
11 P0=760; //in torr
12 XA1=(P0-PB)/(PA-PB); //mole fraction at normal
    boiling point
13 printf('\nXA1=%0.2 f ',XA1)
14 XB1=(1-XA1); //
15 printf('\nXB1=%0.2 f ',XB1)
16 P1=(PA*YA)+(PB*(1-YA)); //
17 printf('\nP1=%0.1 d torr ',P1)
18 YA1=(YA*PA)/P1; //
19 printf('\nYA1=%0.3 f ',YA1)
20 YB1=(1-YA1); //
21 printf('\nYB1=%0.3 f ',YB1)
22
23 //page 149

```

Scilab code Exa 4.7.10 Ex 13

```

1  clear
2  clc
3  PA=300; //in torr
4  PB=800; //in torr
5  XA=0.60; //
6  XB=1-XA; //
7  P=(PA*XA)+(PB*XB); //pressure at which first bubble
    of vapour is formed
8  printf('P=%0.1 d torr ',P)
9  YA=(XA*PA)/P; //mole fraction of components in first
    bubble of vapour
10 printf('\nYA=%0.2 f ',YA)

```



```

11 YB=(1-YA); //ole fraction of components in first
    bubble of vapour
12 printf( '\nYB=%0.2 f ', YB)
13 XA1=(XA*PB)/(PA+((PB-PA)*XA)); //mole fraction of
    last drop of liquid
14 printf( '\nXA1=%0.2 f ', XA1)
15 XB1=(1-XA1); //mole fraction of last drop of liquid
16 printf( '\nXB1=%0.2 f ', XB1)
17 P=(PA*XA1)+(PB*XB1); //pressure when the last droplet
    of liquid remains
18 printf( '\nP=%0.1d torr ', P)
19
20 //page 151

```

Scilab code Exa 4.7.11 Ex 14

```

1 clear
2 clc
3 Tb=353.25; //temperature of benzene in K
4 Tt=383.75; //temperature of toluene in K
5 T=368.15; //temperature in K
6 DelS_vR=-10.6; //
7 Xb=((exp(DelS_vR))-(exp((DelS_vR)*(Tt/T))))/((exp((
    DelS_vR)*(Tb/T)))-(exp((DelS_vR)*(Tt/T)))); //mole
    fraction of benzene
8 printf( 'Xb=%0.4 f ', Xb)
9 Xt=(1-Xb); //mole fraction of benzene
10 printf( '\nXt=%0.4 f ', Xt)
11 Yb=Xb*(exp((-DelS_vR)*(1-(Tb/T)))); //
12 printf( '\nYb=%0.4 f ', Yb)
13 Yt=1-Yb; //
14 printf( '\nYt=%0.4 f ', Yt)
15
16 //There are minor errors in solution in textbook
17 //page 151

```

Scilab code Exa 4.7.12 Ex 15

```
1 clear
2 clc
3 T1=5100; //in K
4 T2=4530; //in K
5 A=16.24; //
6 B=13.38; //
7 PA=760; //in torr
8 PB=PA
9 TA=-(T1/(log10(PA)-A)); //in K
10 printf('TA=%0.1f K',TA)
11 TB=-(T2/(log10(PB)-B)); //in K
12 printf('\nTB=%0.1f K',TB)
13 l=round(TA)+3;
14 u=round(TB)-6;
15 T=1:5:u;
16 for i=1:length(T)
17 P_A=10^(-T1/T(i)+A);
18 P_B=10^(-T2/T(i)+B);
19 x_A(i)=(PA-P_B)/(P_A-P_B);
20 y_A(i)=x_A(i)*P_A/PB
21 end
22 plot(x_A,T,y_A,T); xlabel('Mole fraction xA');ylabel
    ('T/K');
23
24 //There is no numerical solution to the given
    question only a graph is plotted
25 //page 152
```

Scilab code Exa 4.7.13 Ex 16

```

1 clear
2 clc
3 T=391; //temperature in K
4 Yb=0.045; //
5 Ya=0.955; //
6 T=410; //in K
7 X=50/100; //
8 XA=0.09; //composition of liquid at l2
9 XB=0.91; //composition of liquid at l2
10 YA=0.74; //composition of liquid at v2
11 YB=0.26; //composition of liquid at v2
12 N=(YA-X)/(X-XA); //
13 M1=(X-XA)/((YA-X)+(X-XA))*100; //mol % of vapour
14 printf('M1=%0.2 f ',M1)
15 M2=100-M1; //mol % of vapour
16 printf('\nM2=%0.2 f ',M2)
17 XA1=0.035; //composition of liquid at l3
18 XB1=0.965; //composition of liquid at l3
19 Yaf=0.743; //
20 X1=(Ya+Yaf)/2; //
21 printf('\nX1=%0.2 f ',X1)
22
23 //page 154

```

Scilab code Exa 4.7.14 Ex 17

```

1 clear
2 clc
3 n1=270; //amount of sugar in gm
4 N1=358; //molar mass of sugar in gm/mol
5 n2=1; //amount of water in kg
6 N2=18; //molar mass of water in gm/mol
7 M1=n1/N1; //amount of sugar in mol
8 M2=(n2*1000)/N2; //amount of water in mol
9 Mt=M1+M2; //total amount in mol

```

```

10 Xs=M1/Mt; //mole fraction of sugar
11 Xw=M2/Mt; //mole fraction of water
12 R=8.314; //in J/Kmol
13 T=298; //in K
14 DelG_m=(Xs*R*T*log(Xs))+(Xw*R*T*log(Xw))
15 printf('DelG_m=%0.3 f J/mol', DelG_m)
16 DelGm=Mt*DelG_m; //
17 printf('\nDelGm=%0.2 f J', DelGm)
18 DelS_m=-(DelG_m/T); //
19 printf('\nDelS_m=%0.3 f J/Kmol', DelS_m)
20 DelSm=-(DelGm/T); //
21 printf('\nDelSm=%0.3 f J/Kmol', DelSm)
22
23 //There are some errors in the solution given in
    textbook
24 //page 154
25
26 //there are some minor errors in solutions in
    textbook

```

Scilab code Exa 4.7.15 Ex 18

```

1 clear
2 clc
3 R=8.314; //in J/Kmol
4 T=300; //in K
5 Nt=10; //in mol
6 m1=1; //in mol
7 m2=9; //in mol
8 M=10; //in mol
9 DelGm1=Nt*R*T*((m1/M*log(m1/M))+(m2/M*log(m2/M)))
    *(10^-3); //
10 printf('DelGm1=%0.3 f kJ', DelGm1)
11 DelSm1=-((DelGm1/T)*1000); //
12 printf('\nDelSm1=%0.2 f J/K', DelSm1)

```

```

13 Nt1=20; //in mol
14 m3=19; //in mol
15 M1=20; //in mol
16 DelGm3=Nt1*R*T*((m1/M1*log(m1/M1))+(m3/M1*log(m3/M1)
    ))*(10^-3); //
17 printf('\nDelGm3=%0.3 f kJ', DelGm3)
18 DelSm3=-((DelGm3/T)*1000); //
19 printf('\nDelSm3=%0.2 f J/K', DelSm3)
20 DelGm2=DelGm3-DelGm1
21 printf('\nDelGm2=%0.3 f kJ', DelGm2)
22 DelSm2=DelSm3-DelSm1
23 printf('\nDelSm2=%0.2 f J/K', DelSm2)
24
25 //There are some errors in the solution given in
    textbook
26 //page 155

```

Scilab code Exa 4.10.1 Ex 19

```

1 clear
2 clc
3 P=760; //total vapour pressure in torr
4 MA=112.5; //molar mass of chlorobenzene in gm
5 MB=18; //molar mass of water in gm
6 P1=538.9; //vapour pressure of water at 90.6 C
7 PA=(P-P1); //vapour pressure of pure chlorobenzene in
    torr
8 W1=(PA*MA)/(P1*MB); //
9 W2=1/W1; //
10 W=W2+1; //
11 M=100; //in gm
12 WA=M/W; //
13 printf('WA=%0.1 d gm', WA)
14
15 //There are some errors in the solution given in

```

textbook
16 //page 191

Scilab code Exa 4.10.2 Ex 20

```
1 clear
2 clc
3 Pt=747.3; //toatal pressure in torr
4 PB=638.6; //vapour pressure of water
5 PA=Pt-PB; //vapour pressure of liquid
6 WA=1.27; //in gm
7 WB=1; //in gm
8 MB=18; //molar mass of water in gm/mol
9 MA=(WA/WB)*((PB*MB)/PA); //molar mass of liquid in gm
  /mol
10 printf('MA=%0.1 f gm/mol',MA)
11
12 //page 192
```

Scilab code Exa 4.11.1 Ex 21

```
1 clear
2 clc
3 Xw1=0.01; //in gm/dm^3
4 Xw2=0.12; //in gm/dm^3
5 Xw3=0.24; //in gm/dm^3
6 Xb1=1.848*10^-5; //in gm/dm^3
7 Xb2=2.661*10^-3; //in gm/dm^3
8 Xb3=1.089*10^-2; //in gm/dm^3
9 //Taking Xw1,Xw2,Xb1,Xb1 to calculate n
10 n=((log10(Xb1))-(log10(Xb2)))/((log10(Xw1))-(log10(
  Xw2))); //degree of complexity
11 printf('n=%0.1 f',n)
```

```
12
13 //Similarly can be done using lines(4,5,7,8) and
    also for lines (3,5,6,8)
14 //For all we get n=2
15 //page 200
```

Scilab code Exa 4.11.2 Ex 22

```
1 clear
2 clc
3 KD=9; //distribution coefficient
4 M1=0.10825; //amount of p-nitroaniline in gm
5 N=0.00693; //amount of p-nitroaniline chlorine
    dissolved in mol
6 N0=0.04342; //molarity of dil HCl
7 m1=138; //molar mass of p-nitroaniline in gm/mol
8 N1=60; //amount of benzene added in cm3
9 N2=25; //amount of benzene withdrawn in cm3
10 M2=(M1/m1); //amount of free base in 25cm3 of
    benzene
11 X=(M2*(N1/N2)); //amount in mol
12 M=(X/(N1/1000)); //in mol/dm3
13 M0=(M/KD); //molar concentration of free bas e in
    aqeous solution
14 C=(N-(X+M0)); //concentration of unhydrolyzed cation
15 C1=(X+M0); //amount of free base in benzene and water
16 Ct=(N0+C1); //total amount of acid
17 Kh=(M0*Ct)/C; //hydrolysis constant
18 printf('kh=%0.4f mol/dm3',Kh)
19
20 //There are some errors in the solution given in
    textbook
21 //page 201
```

Scilab code Exa 4.11.3 Ex 23

```
1 clear
2 clc
3 KD=25.8; //
4 M1=0.385; // concentration of NH3 in aqueous CuSO4
   solution in mol/dm-3
5 M2=0.0112; // concentration of NH3 in chloroform in
   mol/dm-3
6 m=0.025; // concentration of CuSO4 in mol/dm-3
7 M0=(M2*KD); // concentration of NH3 in aqueous layer
   in mol/dm-3
8 M=M1-M0; // concentration of combined NH3 in mol/dm-3
9 X=(M/m); //
10 printf('X=%0.2 f',X)
11
12 // X is approximately equal to 4
13 //page 202
```

Scilab code Exa 4.11.4 Ex 24

```
1 clear
2 clc
3 Ac=10; //
4 Ab=1; //
5 Kd=Ab/Ac; //
6 wn=0.01; //in gm
7 w=1.00; //in gm
8 Vb=100; //in cm3
9 Vc=10; //in cm3
10 n=log10(wn/w)/(log10((Kd*Vb)/((Kd*Vb)+Vc))); //
11 printf('n=%0.1 f',n)
```



```

12 V=n*10; //in am^3
13 printf( '\nV=%0.1d',V)
14
15 //approximately equal to 7,n=7 is taken in the text
    book
16 ////There are some errors in the solution given in
    textbook
17 //page 202

```

Scilab code Exa 4.11.5 Ex 25

```

1 clear
2 clc
3 KD=4.7; //distribution coefficient
4 W1=20; //amount of ether added in cm^3
5 W2=50; //amount of solution in cm^3
6 M=0.20; //amount of aspirin in gm
7 w2=(M/(1+(W1*KD)/W2)); //mass of aspirin in ether
    phase in gm
8 printf( 'w2=%0.4 f gm',w2)
9 w1=M-w2; //mass of aspirin in aqueous phase in gm
10 printf( '\nw1=%0.4 f gm',w1)
11 n=2
12 W=10; //
13 wn=((1/KD)*W2)/(((1/KD)*W2)+W))^n*(M); //amount of
    aspirin unextracted in gm
14 printf( '\nwn=%0.4 f gm',wn)
15 w=(M-wn); //amount of aspirin extracted in gm
16 printf( '\nw=%0.4 f gm',w)
17
18 //page 203

```

Chapter 5

Phase Diagrams of one component systems

Scilab code Exa 5.2.1 Ex 1

```
1 clear
2 clc
3 DelHm_f=6008.5; //in J/mol
4 m=18; //molar mass of water in gm/mol
5 rho_i=0.917; //density of ice in gm/cm^3
6 rho_l=0.99987; //density of liquid in gm/m^3
7 DelV=((m/rho_l)-(m/rho_i));
8 printf('DelV=%0.3f*10^-6 m^3/mol',DelV/10^-6)
9 T=273.15; //in K
10 P=760; //in mmHg
11 Pt=4.6; //triple point pressure in mmHg
12 DelPDelT=((DelHm_f)/(T*DelV*10^-6));
13 printf('\nDelPdelT=%0.3f 10^6 J/Km^3',DelPDelT/10^6)
14 DelP=((P-Pt)/P)*101.325*10^3; //in N/m^3
15 DelT=(DelP/DelPDelT);
16 printf('\nDelT=%0.4f K',DelT)
17
18 //There are some errors in the solution given in
    textbook
```

19 //page 222

Scilab code Exa 5.2.2 Ex 2

```
1 clear
2 clc
3 R=8.314; //in J/Kmol
4 T=273.15; //in K
5 m=18; //molar mass of water in gm /mol
6 rho_l=0.99987; //density of water ingm/cm^3
7 P2=101.325*10^3; //atmospheric pressure in N/m^2
8 Pt=4.6; //triple point pressure in mmHg
9 P1=(Pt/760)*P2; //
10 P=Pt*10^((((m*10^-3)/(rho_l*10^3))*(P2-P1))/(2.303*R
    *T)); //vapour pressure of liquid water in mmHg
11 printf('P=%0.3 f mmHg',P)
12
13 //page 223
```

Chapter 6

Phase Diagrams of two component systems

Scilab code Exa 6.6.1 Ex 1

```
1 clear
2 clc
3 DelHmA_f=28.87; //enthalpy of fusion of NaCl in KJ/
  mol
4 DelHmB_f=24.06; //enthalpy of fusion of Na2SO4 in KJ/
  mol
5 R=8.314; //in J/Kmol
6 TA=1074; //melting point temperature of NaCl
7 XB=48.2/100; //composition of Na2SO4
8 XA=(1-XB); //composition of NaCl
9 TB=1/((1/TA)-(2.303*R*log10(XA)/(DelHmA_f*10^3))
  +(2.303*R*log10(XB)/(DelHmB_f*10^3))); //melting
  point of Na2SO4 in K
10 printf('TB=%0.1d K',TB)
11 T=(1)/((-2.303*R*log10(XA)/(DelHmA_f*10^3)+(1/TA))
  ); //temperature of the sysytem in K
12 printf('\nT=%0.1f K',T)
13
14 //There are some errors in the solution given in
```

textbook
15 //page 313

Scilab code Exa 6.6.2 Ex 2

```
1 clear
2 clc
3 M=20; //in Kg
4 BC=35; //in mm
5 BA=31; //in mm
6 M1=(BA/(BA+BC))*M; //mass of Sb in Kg
7 printf('M1=%0.2 f Kg',M1)
8 L=(BA+20); //in mm
9 M2=(L/(L+BC))*20; //mass of Sb in Kg
10 printf('\nM2=%0.2 f Kg',M2)
11
12 //page 314
```

Scilab code Exa 6.6.3 Ex 3

```
1 clear
2 clc
3 X1=80; //mol % of n-heptane
4 X2=90; //mol % of n-heptane
5 X3=95; //mol % of n-heptane
6 X=24; //mol % of n-heptane at -114.4 C
7 N1=(100-X1)/(X1-X); //at 80% of n-heptane
8 N2=(X1-X)/(100-X); //at 80% of n-heptane
9 N=((N2*100)/X1)*100; //%of n-heptane recovered w.r.t
    to original n-heptane
10 printf('N=%0.1 f',N)
11 N3=(100-X2)/(X2-X); //at 90% of n-heptane
12 N4=(X2-X)/(100-X); //at 90% of n-heptane
```

```

13 N=((N4*100)/X2)*100; // % of n-heptane recovered w.r.t
    to original n-heptane
14 printf( '\nN=%0.1 f ', N)
15 N5=(100-X3)/(X3-X); // at 95% of n-heptane
16 N6=(X3-X)/(100-X); // at 95% of n-heptane
17 N=((N6*100)/X3)*100; // % of n-heptane recovered w.r.t
    to original n-heptane
18 printf( '\nN=%0.1 f ', N)
19
20 //page 315

```

Scilab code Exa 6.6.5 Ex 4

```

1 clear
2 clc
3 X1=1.04; // solubility of KBr in gm/gm
4 w=1; // amount of H2O in gm
5 X2=0.64; // solubility of KBr after cooling in gm/gm
6 M1=(w/X1)*(X1-X2); // mass of water to be added in gm
7 M2=(X2/w)*M1; // mass of KBr in the solution in gm
8 M=(X1-X2)-M2; // mass of KBr separated in gm
9 N1=M*(100/X1); // percent yield of pure KBr
10 printf( 'N1=%0.2 f ', N1)
11 M3=(w/X1)*X2; // mass of water remained in above
    evaporation process in gm
12 M4=(X2/w)*(M3); // mass of water remained after
    cooling in above evaporation process in gm
13 M=(X2)*M4; // mass of KBr separated in second crop in
    gm
14 Mt=(X1-X2)+M; // total mass of kBr separated in two
    crops in gm
15 N2=Mt*(100/X1); // percent of KBr recovered
16 printf( '\nN2=%0.2 f ', N2)
17
18 // There are some errors in the solution given in

```

Scilab code Exa 6.6.6 Ex 5

```
1 clear
2 clc
3 M1=40; //total amount of mixture of calcium and
    aluminium in gm
4 w1=54; //amount of aluminium in CaAl2 in gm
5 w2=81; //amount of aluminium in CaAl3 in gm
6 W1=70; //total amount of aluminium in gm
7 X1=(W1-((w2/M1)*M1))/((w1/M1)-(w2/M1)); //amount of
    calcium in gm
8 printf('X1=%0.1 f gm',X1)
9 N1=(w1/M1)*X1; //amount of calcium in mixture of
    CaAl2 in gm
10 printf('\nN1=%0.1 f gm',N1)
11 N2=(w2/M1)*(M1-X1); //amount of aluminium in mixture
    of CaAl3 in gm
12 printf('\nN2=%0.1 f gm',N2)
13 M2=20; //total amount of mixture of calcium and
    aluminium in gm
14 W2=90; //total amount of aluminium in gm
15 w3=86; //amount of aluminium after melting
16 M3=14; //percent of calcium mass melted
17 X2=(W2-((w3/M3)*M2))/((w2/M1)-(w3/M3)); //amount of
    calcium in gm
18 printf('\nX2=%0.1 f gm',X2)
19 N3=(w2/M1)*X2; //amount of calcium in mixture of
    CaAL2 in gm
20 printf('\nN3=%0.1 f gm',N3)
21 N4=(w3/M3)*(M2-X2); //amount of aluminium in mixture
    of CaAL3 in gm
22 printf('\nN4=%0.1 f gm',N4)
```

23
24 //There are some errors in the solution given in
 textbook
25 //page 318
26
27 //There are some errors in the solution given in
 textbook

Chapter 8

Electrochemical cells

Scilab code Exa 8.9.2 Ex 1

```
1 clear
2 clc
3 EFe_Pt=0.771; //in V
4 EFe=-0.440; //in V
5 Ecell=(EFe_Pt-EFe); //in V
6 printf('Ecell=%0.3f V',Ecell)
7 E1=1.510; //in V
8 E2=1.223; //in V
9 Ecell=(E1-E2); //in V
10 printf('\nEcell=%0.3f V',Ecell)
11 E3=0.401; //in V
12 E4=-0.601; //in V
13 Ecell=(E3-E4); //in V
14 printf('\nEcell=%0.3f V',Ecell)
15 E5=0.337; //in V
16 E6=0.799; //in V
17 Ecell=(E5-E6); //in V
18 printf('\nEcell=%0.3f V',Ecell)
19 E7=1.44; //in V
20 E8=0.5355; //in V
21 Ecell=(E7-E8); //in V
```

```

22 printf( '\nEcell=%0.4 f V',Ecell)
23 E9=0.7991; //in V
24 E10=-0.126; //in V
25 Ecell=(E9-E10); //in V
26 printf( '\nEcell=%0.4 f V',Ecell)
27 E11=1.51; //in V
28 E12=-0.49; //in V
29 Ecell=(E11-E12); //in V
30 printf( '\nEcell=%0.3 f V',Ecell)
31 E13=1.33; //in V
32 E14=0.771; //in V
33 Ecell=(E13-E14); //in V
34 printf( '\nEcell=%0.3 f V',Ecell)
35 E15=0.771; //in V
36 E16=0.150; //in V
37 Ecell=(E15-E16); //in V
38 printf( '\nEcell=%0.3 f V',Ecell)
39 E17=0.771; //in V
40 E18=0.76; //in V
41 Ecell=(E17-E18); //in V
42 printf( '\nEcell=%0.3 f V',Ecell)
43 E19=0.771; //in V
44 E20=1.080; //in V
45 Ecell=(E19-E20); //in V
46 printf( '\nEcell=%0.3 f V',Ecell)
47
48 //page 448

```

Scilab code Exa 8.9.3 Ex 2

```

1 clear
2 clc
3 E1=1.3595; //in V
4 E2=0.337; //in V
5 Ecell=(E1-E2); //in V

```

```

6  printf( '\nEcell=%0.4 f V',Ecell)
7  E3=1.510; //in V
8  E4=0.337; //in V
9  Ecell=(E3-E4); //in V
10 printf( '\nEcell=%0.3 f V',Ecell)
11 E5=0.7791; //in V
12 E6=0.337; //in V
13 Ecell=(E5-E6); //in V
14 printf( '\nEcell=%0.4 f V',Ecell)
15 E7=0.771; //in V
16 E8=0.150; //in V
17 Ecell=(E7-E8); //in V
18 printf( '\nEcell=%0.3 f V',Ecell)
19 E9=0.771; //in V
20 E10=1.51; //in V
21 Ecell=(E9-E10); //in V
22 printf( '\nEcell=%0.3 f V',Ecell)
23 E11=0.771; //in V
24 E12=-0.126; //in V
25 Ecell=(E11-E12); //in V
26 printf( '\nEcell=%0.3 f V',Ecell)
27
28 //All the positive values of Ecell gives forward
    reaction ,negative values of Ecell gives backward
    reaction
29 //page 451

```

Scilab code Exa 8.9.4 Ex 3

```

1  clear
2  clc
3  E1=1.3595; //in V
4  E2=0.337; //in V
5  Ecell=(E1-E2); //in V
6  printf( '\nEcell=%0.4 f V',Ecell)

```

```

7 E3=1.510; //in V
8 E4=0.337; //in V
9 Ecell=(E3-E4); //in V
10 printf( '\nEcell=%0.3 f V',Ecell)
11 E5=0.7791; //in V
12 E6=0.337; //in V
13 Ecell=(E5-E6); //in V
14 printf( '\nEcell=%0.4 f V',Ecell)
15 E7=0.771; //in V
16 E8=0.150; //in V
17 Ecell=(E7-E8); //in V
18 printf( '\nEcell=%0.3 f V',Ecell)
19 E9=0.771; //in V
20 E10=1.51; //in V
21 Ecell=(E9-E10); //in V
22 printf( '\nEcell=%0.3 f V',Ecell)
23 E11=0.771; //in V
24 E12=-0.126; //in V
25 Ecell=(E11-E12); //in V
26 printf( '\nEcell=%0.3 f V',Ecell)
27
28 //page 451

```

Scilab code Exa 8.9.6 Ex 4

```

1 clear
2 clc
3 E1=1.07; //in V
4 E2=0.45; //in V
5 Ecell=(E1-E2); //in V
6 printf( '\nEcell=%0.2 f V',Ecell)
7 E3=0.71; //in V
8 E4=0.54; //in V
9 Ecell=(E3-E4); //in V
10 printf( '\nEcell=%0.2 f V',Ecell)

```

11
12 //page 454

Scilab code Exa 8.9.7 Ex 5

```
1 clear
2 clc
3 E_RHE=(0.5335-(-2.363)); //reduction reaction at RHE
  in V
4 RT_F=0.05915; //
5 E_LHE=((RT_F/2)*log10(0.1*0.2^2)); //reduction
  reaction at LHE in V
6 Ecell=E_RHE-E_LHE; //cell reaction in V
7 printf('Ecell=%0.4 f V',Ecell)
8 E_RHE=(0.0-0.0713); //reduction reaction at RHE in V
9 RT_F=0.05915; //
10 E_LHE=((RT_F)*log10((0.5^(1/2))/(0.02*0.02))); //
  reduction reaction at LHE in V
11 Ecell=E_RHE-E_LHE; //cell reaction in V
12 printf('\nEcell=%0.4 f V',Ecell)
13 E_RHE=(0.337-(-0.441)); //reduction reaction at RHE
  in V
14 RT_F=0.05915; //
15 E_LHE=((RT_F/2)*log10(0.05/0.01)); //reduction
  reaction at LHE in V
16 Ecell=E_RHE-E_LHE; //cell reaction in V
17 printf('\nEcell=%0.4 f V',Ecell)
18 E_RHE=(0.0-0.0); //reduction reaction at RHE in V
19 RT_F=0.05915; //
20 E_LHE=((RT_F/2)*log10(6.43/0.127)); //reduction
  reaction at LHE in V
21 Ecell=E_RHE-E_LHE; //cell reaction in V
22 printf('\nEcell=%0.4 f V',Ecell)
23 E_RHE=(-0.763-0.337); //reduction reaction at RHE in
  V
```

```

24 RT_F=0.05915; //
25 E_LHE=((RT_F/2)*log10((0.1^2)*0.732)); //reduction
    reaction at LHE in V
26 Ecell=E_RHE+E_LHE; //cell reaction in V
27 printf('\nEcell=%.3f V',Ecell)
28
29 //There are some errors in the solution given in
    textbook
30 //page 455

```

Scilab code Exa 8.10.1 Ex 6

```

1 clear
2 clc
3 RT_F=0.05915; //in V
4 Ecell=0.0295; //in V
5 A=0.1; //
6 B=0.01; //
7 n=(RT_F/Ecell)*(log10(A/B)); //
8 printf('n=%0.1f',n)
9
10 //page 459

```

Scilab code Exa 8.11.1 Ex 7

```

1 clear
2 clc
3 k = [24.4,48.8,73.2,85.4]
4 E = [0.101,0.116,0.129,0.139]
5 l = log10(k./(100-k))
6 plot(l,E,'mo-')
7 [m,c]=reglin(l,E)
8 V=0.0603; //in V

```

```
9 n=V/m; //
10 printf('n=%0.1 f',n)
11
12 //page 460
```

Scilab code Exa 8.12.1 Ex 8

```
1 clear
2 clc
3 E0=0.7991; //in V
4 RT_F=0.05915; //in V
5 K1=6.02*10^-8; //
6 K2=1.995*10^-19; //
7 E1=(E0-(RT_F*(-log10(K1)))); //
8 printf('E1=%0.4 f V',E1)
9 E2=(E0-(RT_F*(-log10(K2)))); //
10 printf('\nE2=%0.4 f V',E2)
11
12 //page 464
```

Scilab code Exa 8.12.2 Ex 9

```
1 clear
2 clc
3 T=298; //temperature in K
4 R=8.314; //J/K
5 F=96500; //in C
6 Kw=(10^-14); //
7 E=((2.303*R*T)/F)*log10(Kw); //reduction potential in
   V
8 printf('E=%0.3 f V',E)
9
10 //page 464
```

Scilab code Exa 8.14.1 Ex 10

```
1 clear
2 clc
3 E3=0.54; //in V
4 E4=0.45; //in V
5 n3=4; //
6 n4=1; //
7 n1=5; //
8 E1=((-n3*(E3))-n4*(E4))/(-n1); // in V
9 printf('E1=%0.2 f V',E1)
10 n2=6; //
11 n5=1; //
12 E5=1.07; //in V
13 E2=((-n3*(E3))-n4*(E4))-n5*(E5))/(-n2); // in V
14 printf('\nE2=%0.2 f V',E2)
15
16 //page 468
```

Scilab code Exa 8.16.1 Ex 11

```
1 clear
2 clc
3 E_RHE=(0.1385); //reduction reaction at RHE in V
4 RT_F=0.05915; //
5 E_LHE=((RT_F*2)*log10(0.2)); //reduction reaction at
   LHE in V
6 Ecell=E_RHE-E_LHE; //cell reaction in V
7 printf('Ecell=%0.4 f V',Ecell)
8
9 //page 473
```

Scilab code Exa 8.16.2 Ex 12

```
1 clear
2 clc
3 Ecell=0.2860; //in V
4 E_RHE=(-0.1522-(-0.403)); //in V
5 RT_F=0.05915; //
6 a=10^((-2/(3*RT_F))*(Ecell-E_RHE))
7 printf('a=%0.4 f',a)
8 a1=a^3; //
9 printf('\na1=%0.5 f',a1)
10
11 //page 474
```

Scilab code Exa 8.16.3 Ex 13

```
1 clear
2 clc
3 m=0.01021; //in mol/kg
4 m1=m*(2*m)^2
5 Ecell=1.1566; //in V
6 E_RHE=(0.222-(-0.762)); //in V
7 RT_F=0.05915; //
8 K=10^((-2/(3*RT_F))*((Ecell-E_RHE)+((RT_F/2)*log10(
   m1)))); //ion activity coefficient
9 printf('K=%0.4 f',K)
10
11 //There are some errors in the solution given in
   textbook
12 //page 474
```

Scilab code Exa 8.18.1 Ex 14

```
1 clear
2 clc
3 n1=2; //
4 F=96500; //in C
5 E=0.0455 //in V
6 DelG=-(n1*F*E); //free energy change in J
7 printf('DelG=%0.1d J',DelG)
8 T=298; //in K
9 dEdT_p=(3.38*10^-4)
10 DelH=-(n1*F*(E-(T*dEdT_p))); //enthalpy change in J
11 printf('\nDelH=%0.1d J',DelH)
12 DelS=(n1*F*dEdT_p); //entropy change in J/K
13 printf('\nDelS=%0.2 f J/K',DelS)
14
15 //There are some errors in the solution given in
    textbook
16 //page 480
```

Scilab code Exa 8.18.2 Ex 15

```
1 clear
2 clc
3 n1=2; //
4 F=96500; //in C
5 E=0.1634 //in V
6 DelG=-(n1*F*E); //free energy change in J
7 printf('DelG=%0.1 f J',DelG)
8 T=298; //in K
9 dEdT_p=(0.000837); //in V/K
10 DelH=-(n1*F*(E-(T*dEdT_p))); //enthalpy change in J
```

```

11 printf( '\nDelH=%.1d J', DelH)
12 DelS=(n1*F*dEdT_p); //entropy change in J/K
13 printf( '\nDelS=%.2 f J/K', DelS)
14
15 //There are some errors in the solution given in
    textbook
16 //page 481

```

Scilab code Exa 8.18.3 Ex 16

```

1 clear
2 clc
3 x=[293,298,303]
4 y=[0.0663,0.06839,0.07048]
5 plot(x,y,'mo-')
6 [m,c]=reglin(x,y)
7 n=2; //
8 F=96500; //in C
9 T=298; //in K
10 E=0.06839; //in V
11 DelG=-n*F*E; //in J
12 printf( 'DelG=%.1 f J', DelG)
13 DelH=-n*F*(E-(T*m)); //in J
14 printf( '\nDelH=%.1 f J', DelH)
15 DelS=n*F*m; //in J/K
16 printf( '\nDelS=%.1 f J/K', DelS)
17
18 //There are some errors in the solution given in
    textbook
19 //page 482

```

Scilab code Exa 8.18.4 Ex 17

```

1 clear
2 clc
3 n1=2; //
4 F=96500; //in C
5 E=0.490 //in V
6 DelG=-(n1*F*E); //free energy change in J
7 printf('DelG=%0.1 f J',DelG)
8 T=298; //in K
9 dEdT_p=-(1.86*10^-4); //in V/K
10 DelH=-(n1*F*(E-(T*dEdT_p))); //enthalpy change in J
11 printf('\nDelH=%0.1 f J',DelH)
12 DelS=(n1*F*dEdT_p); //entropy change in J/K
13 printf('\nDelS=%0.2 f J/K',DelS)
14
15 //page 483

```

Scilab code Exa 8.18.5 Ex 18

```

1 clear
2 clc
3 n=2; //
4 F=96500; //
5 DelH=-217780; //in J
6 T=273; //in K
7 E=1.015; //in V
8 dEdT_p=(1/T)*(E+(DelH/(n*F))); //
9 printf('dEdT_p=%0.3 f*10^-4 V/K',dEdT_p/10^-4)
10
11 //There are some errors in the solution given in
    textbook
12 //page 483

```

Scilab code Exa 8.18.6 Ex 19

```

1 clear
2 clc
3 DelG1=-237.23; //in kJ
4 DelG2=79.71; //in kJ
5 n=2; //
6 DelG=(DelG1+(n*DelG2)); //in kJ
7 F=96500; //in C
8 T=298; //in K
9 E=-((DelG*10^3)/(n*F)); //in V
10 printf('E=%0.3 f V',E)
11 DelH1=-285.85; //in kJ
12 DelH2=56.9; //in kJ
13 DelH=(DelH1+(n*DelH2)); //in kJ
14 dEdT_p=((DelH-DelG)*10^3)/(n*F*T); //in V/K
15 printf('\ndEdT_p=%0.5 f V/K',dEdT_p)
16
17 //error in solution
18 ////There are some errors in the solution given in
    textbook
19 //page 484

```

Scilab code Exa 8.18.7 Ex 20

```

1 clear
2 clc
3 E1=0.771; //in V
4 E2=0.150; //in v
5 E=(E1-E2); //in V
6 n=2; //
7 F=96500; //in C/mol
8 R=8.314; //in J/Kmol
9 T=298; //in K
10 Keq=10^((n*F*E)/(2.303*R*T)); //equilibrium constant
11 printf('Keq=%0.1 f *10^21 ',Keq/10^21)
12

```

13 //page 485

Scilab code Exa 8.18.8 Ex 21

```
1 clear
2 clc
3 E1=1.51; //in V
4 E2=-0.49; //in v
5 E=(E1-E2); //in V
6 n=2; //
7 F=96500; //in C/mol
8 R=8.314; //in J/Kmol
9 T=298; //in K
10 Keq=10^((n*F*E)/(2.303*R*T)); //equilibrium constant
11 printf('Keq=%0. d10^331 ', Keq/10^331)
12 //There are some errors in the solution given in
    textbook
13 //page 486
```

Scilab code Exa 8.18.9 Ex 22

```
1 clear
2 clc
3 E1=-0.40; //in V
4 E2=-0.61; //in v
5 E=(E1-E2); //in V
6 n=2; //
7 F=96500; //in C/mol
8 R=8.314; //in J/Kmol
9 T=298; //in K
10 Keq=10^((n*F*E)/(2.303*R*T)); //equilibrium constant
11 printf('Keq=%0.2 f*10^7 ', Keq/10^7)
12
```

13 //page 486

Scilab code Exa 8.18.10 Ex 23

```
1 clear
2 clc
3 E1=-0.224; //in V
4 E2=0.337; //in v
5 E=(E1-E2); //in V
6 n=2; //
7 F=96500; //in C/mol
8 R=8.314; //in J/Kmol
9 T=298; //in K
10 Keq=10^((n*F*E)/(2.303*R*T)); //equilibrium constant
11 printf('Keq=%0.1 f*10^-19', Keq/10^-19)
12
13 //page 486
```

Scilab code Exa 8.18.11 Ex 24

```
1 clear
2 clc
3 E1=-0.151; //in V
4 E2=0.799; //in v
5 E=(E1-E2); //in V
6 RT_F=0.05913; //in V
7 Ksp=10^(E/RT_F); //solubility product
8 printf('Ksp=%0.2 f*10^-17', Ksp/10^-17)
9
10 //page 487
```

Scilab code Exa 8.18.12 Ex 25

```
1 clear
2 clc
3 E1=0.222; //in V
4 E2=0.095; //in v
5 E=(E1-E2); //in V
6 n=1; //
7 F=96500; //in C/mol
8 R=8.314; //in J/Kmol
9 T=298; //in K
10 Keq=10^((n*F*E)/(2.303*R*T)); //equilibrium constant
11 printf('Keq=%0.1 f', Keq)
12 X=(Keq*0.1)/(1+Keq); //in moldm^3
13 printf('\nX=%0.6 f moldm^3', X)
14 Y=0.1-X; //in moldm^3
15 printf('\nY=%0.6 f moldm^3', Y)
16
17 //error in the solution
18 //There are some errors in the solution given in
   textbook
19 //page 487
```

Scilab code Exa 8.18.13 Ex 26

```
1 clear
2 clc
3 E1=0.337; //in V
4 E2=-0.763; //in v
5 E=(E1-E2); //in V
6 n=2; //
7 F=96500; //in C/mol
8 R=8.314; //in J/Kmol
9 T=298; //in K
10 Keq=10^((n*F*E)/(2.303*R*T)); //equilibrium constant
```



```
11 printf('Keq=%0.1f*10^37',Keq/10^37)
12
13 //page 488
```

Scilab code Exa 8.18.14 Ex 27

```
1 clear
2 clc
3 m=[0.01,0.02,0.05,0.10,0.20];//
4 n=0.01;//
5 mu=m+n;//
6 E=[1.0495,1.0315,1.0073,0.9885,0.9694]
7 E2=0.2225;//in V
8 R=0.05913;//in V
9 O=log10(m/n);//
10 K=(E-E2)/R +0;//
11 plot(mu,K,'mo-');//
12 [m,c]=reglin(mu,K)
13 Ksp=10^-c;//
14 printf('Ksp=%0.2f*10^-14',Ksp/10^-14)
15
16 //There are some errors in the solution given in
    textbook
17 //page 491
```

Scilab code Exa 8.18.15 Ex 28

```
1 clear
2 clc
3 RT_F=0.05913;//in V
4 pH=5;//
5 E1=0.280;//in V
6 E2=0.6996;//in V
```

```

7 E=(E1-E2)+(RT_F*pH); //in V
8 printf('E=%0.4 f V',E)
9 E=0; //
10 pH=(E-(E1-E2))/RT_F; //
11 printf('\npH=%0.1 f',pH)
12 pH=7.5
13 E=(E1-E2)+(RT_F*pH); //in V
14 printf('\nE=%0.4 f V',E)
15
16 //page 489

```

Scilab code Exa 8.18.16 Ex 29

```

1 clear
2 clc
3 RT_F=0.05913; //in V
4 pH=7; //
5 E=0.062; //in V
6 E1=(E-(RT_F*pH)); //in V
7 E2=0.145; //in V
8 pH1=(E2-E1)/RT_F; //
9 printf('pH1=%0.1 f',pH1)
10 E=-0.062; //in V
11 E1=(E-(RT_F*pH)); //in V
12 pH2=(E2-E1)/RT_F; //
13 printf('\npH2=%0.1 f',pH2)
14
15 //page 499

```

Scilab code Exa 8.20.1 Ex 30

```

1 clear
2 clc

```

```

3 RT_F=0.05913; //in V
4 m_LHC=0.01; //
5 gamma_LHC=0.383; //
6 m_RHC=1.0; //
7 gamma_RHC=0.042; //
8 Ecell=- (RT_F*log10((m_LHC*gamma_LHC)/(m_RHC*
    gamma_RHC))); //
9 printf('Ecell=%0.4f V',Ecell)
10
11 //page 525

```

Scilab code Exa 8.20.3 Ex 31

```

1 clear
2 clc
3 RT_F=0.05913; //in V
4 m_LHC=0.01; //
5 gamma_LHC=0.708; //
6 m_RHC=0.10; //
7 gamma_RHC=0.502; //
8 Ecell=((-3/2)*(RT_F*log10((m_LHC*gamma_LHC)/(m_RHC*
    gamma_RHC)))); //
9 printf('Ecell=%0.4f V',Ecell)
10
11 //page 527

```

Scilab code Exa 8.21.2 Ex 32

```

1 clear
2 clc
3 RT_F=0.05913; //in V
4 m_LHC=0.02; //
5 gamma_LHC=0.320; //

```

```

6 m_RHC=0.2; //
7 gamma_RHC=0.110; //
8 E1=0.370; //in V
9 Ecell_1=(-E1)*(RT_F*log10((m_LHC*gamma_LHC)/(m_RHC*
    gamma_RHC))); //in V
10 printf('Ecell_1=%0.4f V',Ecell_1)
11 Ecell_2=(Ecell_1)/(2*E1); //in V
12 printf('\nEcell_2=%0.5f V',Ecell_2)
13
14 //page 536

```

Scilab code Exa 8.23.1 Ex 33

```

1 clear
2 clc
3 E1=-0.277; //in V
4 E2=-0.744; //in V
5 Ecell_1=(E2-E1); //in V
6 printf('Ecell_1=%0.3f V',Ecell_1)
7 Ecell_2=(E1-E2); //in V
8 printf('\nEcell_2=%0.3f V',Ecell_2)
9
10 //page 539

```

Scilab code Exa 8.23.2 Ex 34

```

1 clear
2 clc
3 E1=-0.74; //in V
4 E2=-0.40; //in V
5 E3=-0.91; //in V
6 n1=3; //
7 n2=1; //

```

```

8 n3=6; //
9 n=2; //
10 E=((n1*E1)-(n2*E2))/n; //in V
11 printf('E=%0.2 f V',E)
12 E=E2-E1; //in V
13 DelG1=n1*E; //
14 printf('\nDelG1=%0.2 f V',DelG1)
15 E=E1-E3; //in V
16 DelG2=n3*E; //
17 printf('\nDelG2=%0.2 f V',DelG2)
18 E=E2-E3; //in V
19 DelG3=n*E; //
20 printf('\nDelG3=%0.2 f V',DelG3)
21 RT_F=0.05913; //in V
22 Keq=10^(DelG1/RT_F); //
23 printf('\nKeq=%0.2 f *10^17 ',Keq/10^17)
24
25 //page 540

```

Scilab code Exa 8.23.3 Ex 35

```

1 clear
2 clc
3 Ecell=0; //in V
4 E1=0; //in V
5 E2=-0.40; //in V
6 E=(E1-E2); //in V
7 T=298; //in K
8 RT_F=0.05913; //in V
9 Kw=10^-14; //in moldm^3
10 Ksp=Kw^2*(10^((-2/RT_F)*(-E))); //in (moldm^3)^2
11 printf('Ksp=%0.2 f *10^-15 (moldm^3)^2 ',Ksp/10^-15 )
12
13 dEdT_p=0.002; //in V/K
14 n=2

```

```

15 F=96500; //inC
16 DelG=n*F*E1; //change in gibbs energy
17 printf( '\nDelG=%0.1 d ', DelG)
18
19 DelS=n*F*dEdT_p; //change in entropy in J/K
20 printf( '\nDelS=%0.1 d J/K', DelS)
21
22 DelH=DelG+(T*DelS)*10^-3; //change in enthalpy in kJ
23 printf( '\nDelH=%0.3 f kJ', DelH)
24
25 //page 543

```

Scilab code Exa 8.23.4 Ex 36

```

1 clear
2 clc
3 K1=1; //
4 K2=1.66*10^6; //in dm^3/mol
5 Keq=(K1/K2); //equilibrium constant in mol/dm^3
6 RT_F=0.05913; //in V
7 n=1; //
8 Ecell=RT_F/n*(log10(Keq))
9 printf( ' Ecell=%0.4 f V', Ecell)
10 E_h=0.337; //
11 n2=2
12 Ecell_2=n2*E_h
13 printf( '\nEcell_2=%0.3 f V', Ecell_2)
14
15 //page 544

```

Scilab code Exa 8.23.5 Ex 37

```

1 clear

```

```

2  clc
3  RT_F=0.05913; //in V
4  Ecell=0.1185 //in V
5  K1=0.379*10^-3 //
6  K2=37.9*10^-3 //
7  m=-(RT_F/Ecell)*log10(K1/K2); //
8  printf('m=%0.1 f',m)
9  K3=0.1; //
10 K4=1; //
11 Ecell_1=0.1263; //in V
12 n=(-(Ecell_1*m)/RT_F)/log10(K3/K4); //
13 printf('\nn=%0.1 d',n)
14
15 //page 545

```

Scilab code Exa 8.23.6 Ex 38

```

1  clear
2  clc
3  E1=0.6994; //in V
4  RT_F=0.05913; //in V
5  E=0.7314; //in V
6  kRHE_kLHE=10^((E-E1)/RT_F); //
7  printf('kRHE_kLHE=%0.3 f',kRHE_kLHE)
8  K=0.1; //in mol/dm^3
9  K2=2.1*10^-4; //in mol/dm^3
10 K1=(K2/(kRHE_kLHE^2)); //in mol/dm^3
11 printf('\nK1=%0.2 f*10^-5 mol/dm^3',K1/10^-5)
12 pH1=-log10(sqrt(K2*K))
13 printf('\npH1=%0.3 f',pH1)
14 pH2=-log10(sqrt(K1*K))
15 printf('\npH2=%0.3 f',pH2)
16
17 //page 545

```

Scilab code Exa 8.23.7 Ex 39

```
1 clear
2 clc
3 Ecell=-0.188; //in V
4 RT_F=0.05913; //in V
5 H=10^(Ecell/RT_F); //in mol/dm^3
6 M=1/32; //
7 alpha=(H/M); //degree of freedom
8 printf('alpha=%0.3f*10^-2',alpha/10^-2)
9 Kh=(M*alpha^2)/(1-alpha); //
10 printf('\nKh=%0.2f*10^-5 mol/dm^3',Kh/10^-5)
11
12 //page 547
```
