

BINDURA UNIVERSITY OF SCIENCE EDUCATION

ENGINEERING AND PHYSICS DEPARTMENT

PH 208: SOLID STATE PHYSICS

DURATION: THREE HOURS

JAN 2025

Answer ALL parts of Section A and any THREE questions from Section B. Section

A carries 40 marks and Section B carries 60 marks.

SECTION A

- 1 (a) (i) Inspite of their long range order in the arrangement of particles, why are the crystals usually not perfect? [2]
- (ii) Which of the following statements are true about semiconductors? [4]
- (a) Silicon doped with electron rich impurity is a p-type semiconductor
(b) Silicon doped with an electron rich impurity is an n-type semiconductor
(c) Delocalised electrons increase the conductivity of doped silicon
(d) An electron vacancy increases the conductivity of n-type semiconductor
- (b) (i) Under which situations can an amorphous substance change to crystalline form? [3]
- (ii) Match the items given in Column I with the items given in Column II. [5]

Column I	Column II
A. Mg in solid state	1. P-type semiconductor
B. MgCl ₂ in molten state	2. N-type semiconductor
C. Silicon with phosphorus	3. Electrolytic conductors
D. Germanium with Boron	4. Electronic com

- (c) Find the kinetic energy of an electron with a wavelength of 0.5 nm [4]
- (d) Determine the energy and radius of the electron orbit of s hydrogen atom for n=1. [10]
- (e) Using Figure 1.1, determine the dipole moment associated with the ionic model of the water molecule. NB: the length of the O-H bond is 0.097 nm and the angle between the bonds is 104.5°. [6]

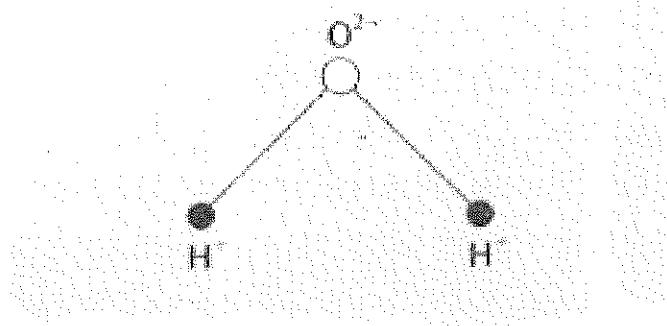


Figure 1.1 Ionic model

- (f) Estimate the stress required to open a surface crack in a glass rod given that the crack has a depth of $2.0 \mu\text{m}$ and a radius of curvature at the tip of 1.0 nm . [4]
- (g) Estimate the typical conductivity of a metal at 295 K assuming that the mean free path is about 1 nm and that the number of valency electrons is about 10^{29} m^{-3} . [3]

Section B

- 2 (a) [8]

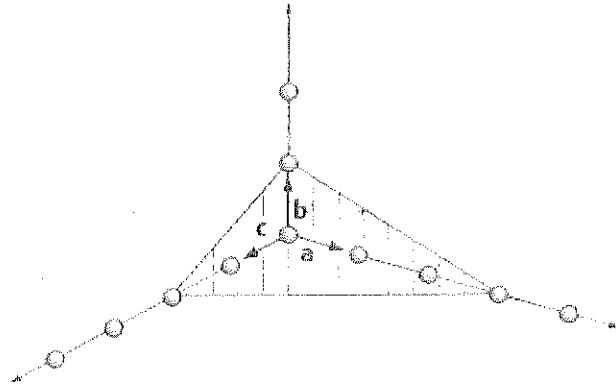


Figure 2

Determine the Miller indices for the plane shown in Fig. 2

- (b) Determine the actual volume occupied by the spheres in the simple cubic structure as a percentage of the total volume. [12]
- 3 (a) The energy-wavevector ($E - k$) dispersion relation for a particle in two dimensions is $E = Ck$, where C is a constant. If its density of states $D(E)$ is proportional to E^p . What is the value of p ? [4]

- (b) The energy dispersion for electrons in one dimensional lattice with lattice parameter a is given by $E(k) = \varepsilon_0 - \frac{1}{2}W \cos ka$, where W and ε_0 are constants. What is the effective mass of the electron near the bottom of the band? [4]
- (c) Discuss the advantages and disadvantages of using electron microscopes in comparison to X-ray diffraction for determining the structure of a crystal. [6]
- (d) A lift cable of diameter 2 cm has a yield stress of $2 \times 10^8 \text{ Nm}^{-2}$. Assuming that the empty lift has a mass of 150 kg and that the average mass of a person is 70 kg, determine the maximum number of people that can be carried in the lift given that the safe operating limit is 20% of the elastic limit. [6]
- 4 (a) Explain why the conductivity of germanium crystals increases on doping with gallium? [10]
- (b) How does the doping increase the conductivity of semiconductors? [10]
- 5 (a) (i) n-type semiconductors.
Use Bohr's model to determine the radius of the donor electron orbit for a phosphorus atom in a silicon crystal.
(ii) Estimate the number of silicon atoms in a sphere of this radius. (The nearest neighbour separation in silicon is 0.234 nm.) [10]
- (b) By making use of the Bohr's model of the hydrogen atom estimate the energy of the donor electron in Silicon. [4]
- (c) Given that the grains in a polycrystalline metal are typically $50 \mu\text{m}$ across and that metal ions have radius of 0.5 nm, estimate the average number of ions in a grain and the proportion of these ions which are adjacent to a grain boundary. [6]
- 6 (a) By considering the group velocity v_g of an electron in a crystal, show that the effective mass of the particle is given by $m_e^* = \hbar^2 \left(\frac{d^2 E}{dk^2} \right)^{-1}$ [14]
- (b) Explain why silicon is opaque, whereas diamond is transparent.
(Hint: consider the minimum frequency of light that is absorbed in each case.) [6]

Formula Sheet

A.1 Physical Constants

The following table contains the values in SI units for some physical constants that are particularly important in solid-state physics.

Name	Symbol	Value
speed of light in vacuum	c	$299\,792\,458 \text{ m s}^{-1}$
magnetic constant (permeability of free space)	μ_0	$4\pi \times 10^{-7} \text{ N A}^{-2}$
electric constant (permittivity of free space)	$\epsilon_0 = 1/\mu_0 c^2$	$8.854\,188 \times 10^{-12} \text{ F m}^{-1}$
elementary charge	e	$1.602\,176 \times 10^{-19} \text{ C}$
Planck constant in eV	h	$6.626\,069 \times 10^{-34} \text{ J s}$
reduced Planck constant in eV	$h/\{e\}$	$4.135\,667 \times 10^{-15} \text{ eV s}$
fine-structure constant inverse of α	$b = h/2\pi$	$1.054\,572 \times 10^{-31} \text{ J s}$
magnetic flux quantum	$h/\{e\}$	$6.582\,119 \times 10^{-16} \text{ eV s}$
conductance quantum inverse of G_0	$\alpha = e^2/4\pi\epsilon_0\hbar c$	$7.297\,353 \times 10^{-3}$
Josephson constant	α^{-1}	$137.035\,999$
von Klitzing constant	$\Phi_0 = h/2e$	$2.007\,834 \times 10^{-15} \text{ Wb}$
Boltzmann constant	$G_0 = 2e^2/h$	$7.748\,092 \times 10^{-5} \text{ S}$
Avogadro constant	$G_0^{-1} = h/2e^2$	$12\,906.404 \Omega$
molal gas constant	$R = N_A k_B$	$483\,597.9 \times 10^3 \text{ Hz V}^{-1}$
Bohr magneton	$\mu_B = eh/2m_e$	$25\,812.807 \Omega$
nuclear magneton	$\mu_N = eh/2m_p$	$1.380\,650 \times 10^{-23} \text{ J K}^{-1}$
Bohr radius	$a_0 = 4\pi\epsilon_0\hbar^2/m_e e^2$	$6.022\,142 \times 10^{23} \text{ mol}^{-1}$
electron mass	m_e	$8.31\,472 \text{ J mol}^{-1} \text{ K}^{-1}$
electron magnetic moment	μ_e	$9.274\,009 \times 10^{-24} \text{ J T}^{-1}$
electron g -factor	$g_e = 2\mu_e/\mu_B$	$5.050\,783 \times 10^{-27} \text{ J T}^{-1}$
electron gyromagnetic ratio	$\gamma_e = 2 \mu_e /h$	$0.529\,177 \times 10^{-10} \text{ m}$
neutron mass	m_n	$9.109\,382 \times 10^{-31} \text{ kg}$
neutron magnetic moment	μ_n	$-9.284\,764 \times 10^{-25} \text{ J T}^{-1}$
neutron g -factor	$g_n = 2\mu_n/\mu_N$	$-1.001\,160 \mu_B$
proton mass	m_p	$-2.002\,319$
proton-electron mass ratio	m_p/m_e	$1.760\,860 \times 10^{11} \text{ s}^{-1} \text{ T}^{-1}$
proton magnetic moment	μ_p	$28324.9532 \text{ MHz T}^{-1}$
proton g -factor	$g_p = 2\mu_p/\mu_N$	$1.674\,927 \times 10^{-27} \text{ kg}$
proton gyromagnetic ratio	$\gamma_p = 2\mu_p/h$	$-0.936\,236 \times 10^{-26} \text{ J T}^{-1}$
Rydberg constant	$\gamma_p/2\pi$	$-1.943\,043 \mu_N$
Rydberg energy in electronvolts	$R_\infty = \alpha^2 m_e c/2\hbar$	$-3.826\,085$
Hartree energy in electronvolts	$Ry = R_\infty \hbar c$	$1.672\,622 \times 10^{-27} \text{ kg}$
	$E_h = e^2/4\pi\epsilon_0\hbar c$	$1836.152\,607$
		$5.446\,170\,10^{-1}$
		$1.410\,607 \times 10^{-26} \text{ J T}^{-1}$
		$2.792\,847 \mu_N$
		$5.585\,695$
		$2.675\,222 \times 10^8 \text{ s}^{-1} \text{ T}^{-1}$
		$42.577\,481 \text{ MHz T}^{-1}$
		$10.973\,731.569 \text{ m}^{-1}$
		$2.179\,872 \times 10^{-18} \text{ J}$
		$13.605\,692 \text{ eV}$
		$4.359\,714 \times 10^{-18} \text{ J}$
		$27.211\,384 \text{ eV}$

Quantum theory of solid states

$$\begin{aligned}
 \psi(x) &= u(x)e^{ikx} \\
 \frac{\partial^2 u_1(x)}{\partial x^2} + 2jk \frac{\partial u_1(x)}{\partial x} - (k^2 - \alpha^2) u_1(x) &= 0 \\
 \alpha^2 &= \frac{2mE}{\hbar^2} \\
 \frac{2m}{\hbar^2}(E - V_0) &= \alpha^2 - \frac{2mV_0}{\hbar^2} = \beta^2 \\
 \frac{\partial^2 u_2(x)}{\partial x^2} + 2jk \frac{\partial u_2(x)}{\partial x} - (k^2 - \alpha^2 + \frac{2mV_0}{\hbar^2}) u_2(x) &= 0 \\
 P' \frac{\sin \alpha a}{\alpha a} + \cos \alpha a &= \cos ka \\
 P' &= \frac{mV_0ba}{\hbar^2} \\
 \frac{1}{\hbar} \frac{\partial E}{\partial k} &= \frac{p}{m} = v \\
 \frac{1}{m} &= \frac{1}{\hbar^2} \frac{\partial^2 E}{\partial k^2} \\
 g(E) &= \frac{4\pi(2m)^{\frac{3}{2}}}{\hbar^3} \sqrt{E} \\
 \frac{N(E)}{g(E)} &= f_F(E) = \frac{1}{1 + \exp\left\{\frac{(E - E_F)}{k_B T}\right\}}
 \end{aligned}$$

Semiconductor

$$\begin{aligned}
 N_c &= 2 \left(\frac{2\pi m_c^* k_B T}{\hbar^2} \right)^{\frac{3}{2}} \\
 N_v &= 2 \left(\frac{2\pi m_v^* k_B T}{\hbar^2} \right)^{\frac{3}{2}} \\
 n(E) &= g_c(E)f_F(E) \\
 E_{F_i} - E_{\text{midgap}} &= \frac{3}{4} k_B T \ln \left(\frac{m_v^*}{m_n^*} \right) \\
 n_i &= N_c \exp \left\{ \frac{-(E_c - E_{F_i})}{k_B T} \right\} \\
 n_0 &= N_c \exp \left\{ \frac{-(E_c - E_F)}{k_B T} \right\} \\
 p_0 &= N_v \exp \left\{ \frac{-(E_F - E_v)}{k_B T} \right\} \\
 n_i^2 &= N_c N_v \exp \left\{ \frac{-(E_c - E_v)}{k_B T} \right\} \\
 n_0 &= n_i \exp \left\{ \frac{(E_p - E_{F_i})}{k_B T} \right\} \\
 p_0 &= n_i \exp \left\{ \frac{-(E_F - E_{F_i})}{k_B T} \right\} \\
 n_d &= \frac{N_d}{1 + \frac{1}{2} \exp \frac{E_d - E_F}{k_B T}} = N_d - N_d^+ \\
 p_a &= \frac{N_a}{1 + \frac{1}{2} \exp \frac{E_F - E_a}{k_B T}} = N_a - N_a^+ \\
 n_0 &= \frac{N_d \cdot N_a}{2} + \sqrt{\left(\frac{N_d \cdot N_a}{2}\right)^2 + n_i^2} \\
 p_0 &= \frac{N_a - N_d}{2} + \sqrt{\left(\frac{N_a - N_d}{2}\right)^2 + n_i^2} \\
 E_c - E_F &= k_B T \ln \left(\frac{N_c}{n_i} \right) \\
 E_F - E_{F_i} &= k_B T \ln \left(\frac{n_i}{n_i} \right) \\
 E_F - E_v &= k_B T \ln \left(\frac{N_v}{p_0} \right) \\
 E_{F_i} - E_F &= k_B T \ln \left(\frac{p_0}{n_i} \right) \\
 n_0 p_0 &= n_i^2
 \end{aligned}$$

Properties of Si and GaAs

Property	Si	GaAs
Atom density (cm^{-3})	5.02×10^{22}	4.43×10^{22}
Atomic weight	28.09	144.64
Crystal structure	Diamond	Zincblende
Density (g/cm^3)	2.329	5.317
Lattice constant (\AA)	5.43102	5.6533
Dielectric constant	11.9	12.9
Electron affinity χ (V)	4.05	4.07
Energy gap (eV)	1.12 (indirect)	1.42 (direct)
Effective density of states in conduction band, N_c (cm^{-3})	2.8×10^{19}	4.7×10^{17}
Effective density of states in valence band, N_v (cm^{-3})	2.65×10^{19}	7.0×10^{18}
Intrinsic carrier concentration n_i (cm^{-3})	9.65×10^9	2.1×10^6
Effective mass (m^*/m_0)	Electrons	$m_e^* = 0.98$ $m_h^* = 0.19$
	Holes	$m_e^* = 0.16$ $m_h^* = 0.49$
Drift mobilities ($\text{cm}^2/\text{V}\cdot\text{s}$)	Electrons μ_n Holes μ_p	1,450 500 1×10^3 $2.5-8 \times 10^3$ $\approx 10^{-3}$
Saturation velocity (cm/s)		3.42×10^7
Breakdown field (V/cm)		$8,000$ 400 7×10^6 $3-9 \times 10^4$
Minority-carrier lifetime (s)		$\approx 10^{-3}$ $\approx 10^{-3}$
Index of refraction		3.3
Optical-phonon energy (eV)		0.063
Melting point ($^\circ\text{C}$)		1414
Linear coefficient of thermal expansion $\Delta L/L\Delta T$ ($^\circ\text{C}^{-1}$)		2.59×10^{-6}
Thermal conductivity (W/cm-K)		1.56
Thermal diffusivity (cm^2/s)		0.9
Specific heat (J/g- $^\circ\text{C}$)		0.713
Heat capacity (J/mol- $^\circ\text{C}$)		20.07
Young's modulus (GPa)		130
		85.5

Note: All properties at room temperature.

Appendix F Properties of Important Semiconductors

Semiconductor	Crystal Structure	Lattice Const. at 300 K (Å)	Band Gap (eV)	Band gap (eV)	Mobility at 300 K (cm ² /V·s)	Effective Mass (m ₁ , m ₂)	$\mu_1 \mu_2$
C Carbon (diamond)	D	3.66683	5.42	5.38	1.890	1.270	0.23
Ge Germanium	D	5.64612	0.66	0.74	1.340	1.940	0.645, 0.258
Si Silicon	D	5.45102	1.12	1.17	1.450	5.00	0.98, 0.19
IV-VI-III-VI-SIC SiC/SiC/SiC	W $\sigma = 3.086 \times 10^{17}$	2.926	3.03	4.00	80	0.60	1.99
III-V AlN AlN	Z	5.6605	2.36	2.23	1.180	—	—
AlP Aluminum arsenide	Z	5.4635	2.42	2.51	60	450	0.212
AlPb Aluminum phosphide	Z	6.1385	1.58	1.68	200	420	0.12
AlSb Aluminum antimonide	Z	3.6157	6.4	—	260	590	0.26
Boron nitride	W	$\sigma = 2.653 \times 10^7$	5.8	—	—	—	—
BP Boron phosphide	Z	4.5383	2.0	—	40	200	0.24
GaAs Gallium arsenide	Z	5.6653	1.42	1.52	8.000	400	0.003
GaN Gallium nitride	W $\sigma = 3.189 \times 5.182$	3.44	3.59	400	10	0.27	0.8
GaP Gallium phosphide	Z	5.4512	2.26	2.34	110	73	0.82
GaSb Gallium antimonide	Z	6.0959	0.72	0.81	5.000	850	0.042
InAs Indium arsenide	Z	6.0584	0.76	0.42	23,000	450	0.023
InP Indium phosphide	Z	5.8636	1.35	1.42	0	4,600	150
InSb Indium antimonide	Z	6.4794	0.17	0.23	0	80,000	1,250
CaS Calcium sulfide	Z	5.825	2.3	—	—	—	—
CaSe Calcium selenide	W $\sigma = 4.136 \times 6.714$	2.49	—	D	150	40	0.14
CaTe Calcium telluride	Z	6.0850	1.70	1.85	D	300	0.13
ZnO Zinc oxide	Z	6.182	1.58	—	D	1,050	150
ZnS Zinc sulfide	Z	4.530	3.35	3.42	D	250	150
W _x T _y PS Lead sulfide	Z	5.410	1.66	1.84	D	600	0.19
W _x T _y Se Lead selenide	W $\sigma = 1.825 \times 6.736$	2.78	—	D	750	650	0.25
PtTe Lead telluride	R	5.9362	0.41	0.26	1	6,000	4,000
ZnCl ₂ Zinc chloride	Z	6.4620	0.31	0.19	1	6,000	0.17

D = Diamond, W = Wurtzite, Z = Zincblende, R = Rock salt. I, D = Indirect-direct bandgap. CaS, CaSe, CaTe, CaS_xTe_y have effective masses.

Group	1	2	3	4	5	6	7	8	9	10	11	12	13	14	15	16	17	18
Period																		
1	H																	
2	Li	B																
3	Na	Mg																
4	K	Ca	Sc	Ti	V	Cr	Mn	Fe	Co	Ni	Cu	Zn	Ga	Ge	As	S	Cl	Ar
5	Rb	Sr	Y	Zr	Nb	Mo	Tc	Ru	Rh	Pd	Au	Cd	In	Sn	Sb	Te	I	Xe
6	CS	Ba	*	Lu	Hf	Ta	V	Re	Os	Ir	Pt	Au	Hg	Tl	Pb	Bi	Po	Rn
7	Fr	Pb	**	Lr	Rf	Db	Sg	Bh	Ms	Mt	Ds	Fm	Uut	Uup	Uuj	Uus	Uuo	
*Lanthanoids																		
**Actinoids																		

End of exam