

(b) **Cohesion energy.** From the very existence of solids, two general conclusions can be drawn.

- (i) There must be an *attractive force* between atoms or molecules in a solid which keeps them together.
- (ii) There must also be a *repulsive force* acting between the atoms. This is why large external pressure is required to compress a solid to any appreciable extent.

When two identical or non-identical atoms, which form a molecule, are brought from infinity to a close proximity, they first attract each other and then, if they are brought closer than a certain distance, repel each other. In other words, a kind of spring effect takes place. As the potential V is related to the force by the relation $F = -\frac{dV}{dr}$, the resulting potential is a function of distance of separation r . The variation of corresponding potential energy V in eV with distance r in Å is shown in Fig. 3.1

The attractive force gives rise to a negative potential and the repulsive force gives rise to a positive potential shown by dotted curves. The resultant potential energy is shown by full line curve. It has a minimum at a distance r_0 and represents the equilibrium position, where the attractive and the repulsive forces balance each other and potential energy has a maximum negative value. The potential energy $V(r_0)$ at a distance r_0 is thus responsible for keeping the atoms bound in a solid and gives the cohesive or binding energy of the solid.

The potential energy $V(r_0)$ is a negative quantity. The positive quantity $-V(r_0) = D$ is called the dissociation energy of the molecule.

The cohesive energy of a solid is defined as the energy which will be given out in the process of the formation of a crystal by bringing neutral atoms from infinity to the position of equilibrium separation.

If the potential energy due to attraction V_a is supposed to vary as the m th power of the distance and that due to repulsion V_r , as n th power of r , then

$$V_a \propto \frac{1}{r^m} = -\frac{A}{r^m} \text{ and } V_r \propto \frac{1}{r^n} = \frac{B}{r^n}$$

∴ Resultant potential or cohesive energy

$$V = V_a + V_r = -\frac{A}{r^m} + \frac{B}{r^n}$$

where A and B are constants of proportionality. The cohesive energy is generally expressed in eV per atom.

The force between the two atoms is given by

$$F = -\frac{dV}{dr} = \frac{mA}{r^{m+1}} - \frac{nB}{r^{n+1}}$$

At $r = r_0$ the attractive and repulsive forces balance each other and $F(r_0) = 0$. In other words at $r = r_0$

$$\frac{mA}{r_0^{m+1}} - \frac{nB}{r_0^{n+1}} = 0$$

$$\frac{m}{n} = \frac{B}{A} r_0^{m-n}$$

$$r_0^{m-n} = \frac{A}{B} \frac{m}{n} \quad \dots (i)$$

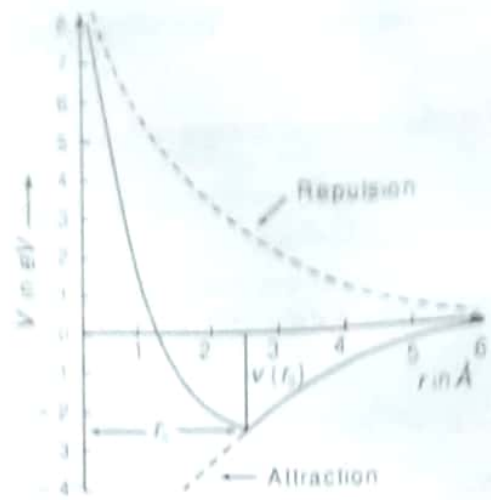


Fig. 3.1

Hence, the potential energy in the equilibrium state

$$V(r_0) = \frac{-A}{r_0^m} + \frac{B}{r_0^n} = \frac{-A}{r_0^m} \left[1 - \frac{B}{A} r_0^{m-n} \right] = \frac{-A}{r_0^m} \left[1 - \frac{m}{n} \right]$$

Although the attractive and repulsive forces are in equilibrium at $r = r_0$, the potential energy is not zero as $m \neq n$. If $n \gg m$ the total energy is essentially the energy of attraction given by $-\frac{A}{r_0^m}$. For the potential energy to have a minimum at $r = r_0$ we have $\left. \frac{d^2V}{dr^2} \right|_{r=r_0} > 0$ i.e., it must be positive.

Now,

$$\frac{d^2V}{dr^2} = \frac{-m(m+1)A}{r^{m+2}} + \frac{n(n+1)B}{r^{n+2}} > 0$$

Substituting the value of $r = r_0$ from (i), we have

$$-m(m+1)A \times \frac{B}{A} \frac{n}{m} \left(\frac{1}{r_0^{n+2}} \right) + \frac{n(n+1)B}{r_0^{n+2}} > 0$$

or

$$-m-1+n+1 > 0$$

or

$$n-m > 0 \text{ or } n > m$$

In other words, a minimum in the energy curve is possible only if $n > m$. Thus the formation of a chemical bond requires that the repulsive forces be of shorter range than the attractive forces.

The energy $V(r_0)$ at the equilibrium distance r_0 is called the *binding energy*, the energy of *cohesion* or *dissociation energy* of the molecule. This much energy is required to separate the atoms of a diatomic molecule to an infinite distance apart.

The cohesive energy may also be defined as the energy released when two atoms are brought close to each other at the equilibrium distance r_0 .

Larger the energy released, more stable the bond formed and hence more stable is the crystal structure.

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