

Spacing of Planes: —

The distance between successive lattice planes of the same type is called the Spacing of Planes or Inter-Planar distance between the planes.

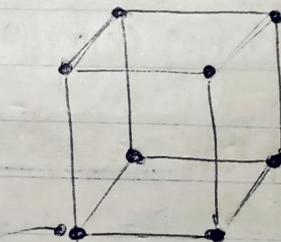
The crystal belonging to cubic system have three kinds of Bravais lattices, depending upon the shape of the unit cells.

These are:

(1) Simple Cubic lattices: — (Cubic - P - lattices)

When one unit is situated only at each corner of the cubic cell, it is called a simple cubic lattice, as shown in fig.

In this case



(d_{100}) is the length of the face or 1, so the cube is of unit length.

(d_{110}) is equal to half-diagonal of a face or $\frac{\sqrt{2}}{2}$ or $\frac{1}{\sqrt{2}}$

(d_{111}) is equal to one third of the diagonal of cube i.e. $\frac{\sqrt{3}}{3}$ or $\frac{1}{\sqrt{3}}$

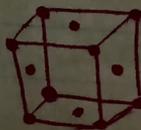
So, the ratio between these distances is given by,

$$d_{100} : d_{110} : d_{111} \quad \therefore \quad 1 : \frac{1}{\sqrt{2}} : \frac{1}{\sqrt{3}}$$

$$\text{or } \frac{1}{d_{100}} : \frac{1}{d_{110}} : \frac{1}{d_{111}} = 1 : \sqrt{2} : \sqrt{3}$$

(2) Face Centred Cubic lattice or Cubic F - lattices: —

When one unit is situated at each corner and one at the centre of each of the faces, then it is known as face centred cubic lattices.



In this case

d_{100} is same as the Simple Cubic lattice i.e. 1

d_{110} is also same as the Simple Cubic lattice i.e. $\frac{1}{\sqrt{2}}$

but d_{111} is double the value of d_{111} of the Simple Cubic lattice i.e. $\frac{2}{\sqrt{3}}$

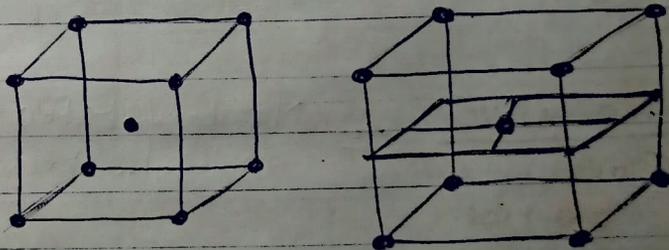
$$\therefore d_{100} : d_{110} : d_{111} = 1 : \frac{1}{\sqrt{2}} : \frac{2}{\sqrt{3}}$$

$$= 1 : 0.707 : 1.154$$

$$\text{or, } \frac{1}{d_{100}} : \frac{1}{d_{110}} : \frac{1}{d_{111}} = 1 : \sqrt{2} : \frac{\sqrt{3}}{2}$$

(iii) One body Centred Cubic lattices: —

In a body Centred Cubic lattice, One unit is situated at each corner and one at the centre of the cube, as shown in fig.



In this case.

d_{100} is same as in Simple Cubic lattice i.e. 1

d_{110} is double as the value of d_{110} in Simple Cubic lattice i.e. $\frac{2}{\sqrt{2}}$

d_{111} is same as in Simple Cubic lattice i.e. $\frac{1}{\sqrt{3}}$

$$\therefore d_{100} : d_{110} : d_{111} = 1 : \frac{2}{\sqrt{2}} : \frac{1}{\sqrt{3}}$$

$$= 1 : \sqrt{2} : \frac{1}{\sqrt{3}}$$

$$\text{or } \frac{1}{d_{100}} : \frac{1}{d_{110}} : \frac{1}{d_{111}} = 1 : \frac{1}{\sqrt{2}} : \sqrt{3}$$

Example: - Iron II Oxide, FeO , crystal has a cubic structure and each edge of the unit cell is 5.0 \AA .

Taking density of the oxide as 4 g/cc .
 Calculate - the number of Fe^{++} and O^{--} ions present in each unit cell.

$$\therefore \text{Edge of the unit cell} = 5 \text{ \AA} = 5 \times 10^{-8} \text{ cm}$$

$$\text{Volume of the unit cell} = (5 \times 10^{-8} \text{ cm})^3$$

$$= 125 \times 10^{-24} \text{ c.c} = 1.25 \times 10^{-22}$$

$$\therefore \text{density of FeO} = 4 \text{ g/cc}$$

$$d = \frac{M}{V} \quad \therefore \text{Mass of unit cell} = \text{Density} \times \text{Volume}$$

$$= 1.25 \times 10^{-22} \times 4$$

$$= 5.00 \times 10^{-22}$$

$$= 5 \times 10^{-22} \text{ g}$$

$$\therefore \text{Mass of FeO (one molecule)} = \frac{\text{Molecular wt.}}{\text{Avogadro Number}}$$

$$= \frac{72}{6.023 \times 10^{23}} = 1.195 \times 10^{-22} \text{ g}$$

Mass of unit cell
 Mass of one molecule
 of unit cell.

$$\therefore \text{Number of FeO molecules per unit cell} = \frac{5 \times 10^{-22}}{1.195 \times 10^{-22}}$$

$$= \frac{5}{1.196} = 4.19$$

$$= 4$$

Hence, there are Four (Fe^{++}) ions and Four (O^{--}) ions in Each unit cell.