

LECTURE # 9

Note Title

10/8/2007

Solid State Physics

Chapter 4

Ashcroft-Mermin

Poly-crystalline

1915 BRAGG

distance between atoms $\sim 1 \text{ \AA} \sim 10^{-8} \text{ cm}$

need probe system with $\lambda \sim 1 \text{ \AA}$

Photons

$$h\nu_{\text{photon}} [\text{eV}] = \frac{12400}{\lambda [\text{\AA}]} \sim 10^4 \text{ eV}$$

10 keV

⇓
X RAY

MATHEMATICAL DESCRIPTION OF LATTICE

1) BRAVAIS LATTICE

(A) SAME "VIEW"

(B) $\vec{R} = m_1 \vec{a}_1 + m_2 \vec{a}_2 + m_3 \vec{a}_3$

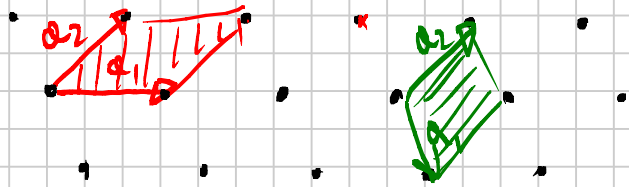
$\vec{a}_1, \vec{a}_2, \vec{a}_3$

PRIMITIVE
VECTORS

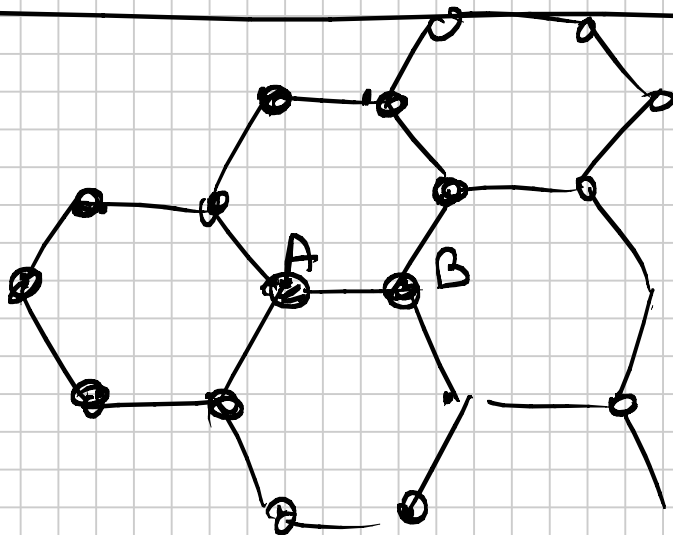
$m_1, m_2, m_3 \in \mathbb{Z}$

3D
3 vectors

2D
2 vectors



choice of primitive
vectors is not unique



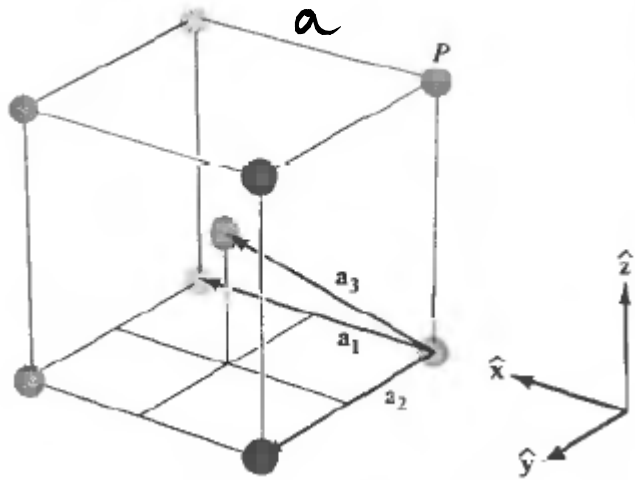
HONEYCOMB
LATTICE

NOT A BRAVAIS
LATTICE

3D BRAVAIS LATTICES (CUBIC not very common)

BODY CENTERED CUBIC (BCC)

a LATTICE CONSTANT



$$\mathbf{a}_1 = a(100)$$

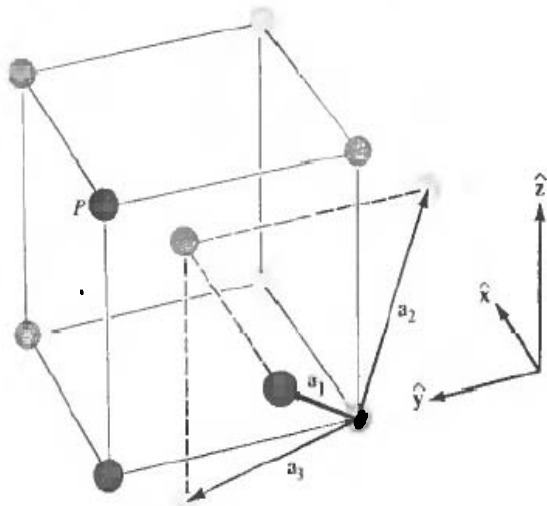
$$\mathbf{a}_2 = a(010)$$

$$\mathbf{a}_3 = \frac{a}{2}(111)$$

Fe
Li
Na
K
W

BCC

$\bar{1} = -1$

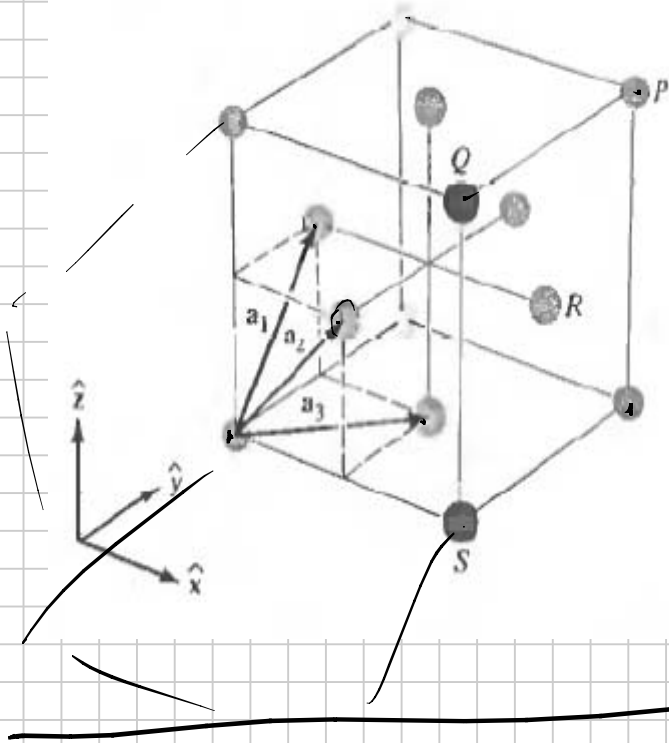


$$\mathbf{a}_1 = \frac{a}{2}(11\bar{1})$$

$$\mathbf{a}_2 = \frac{a}{2}(1\bar{1}1)$$

$$\mathbf{a}_3 = \frac{a}{2}(\bar{1}11)$$

FACE CENTERED CUBIC (FCC)



$$\vec{a}_1 = \frac{a}{2} (011)$$

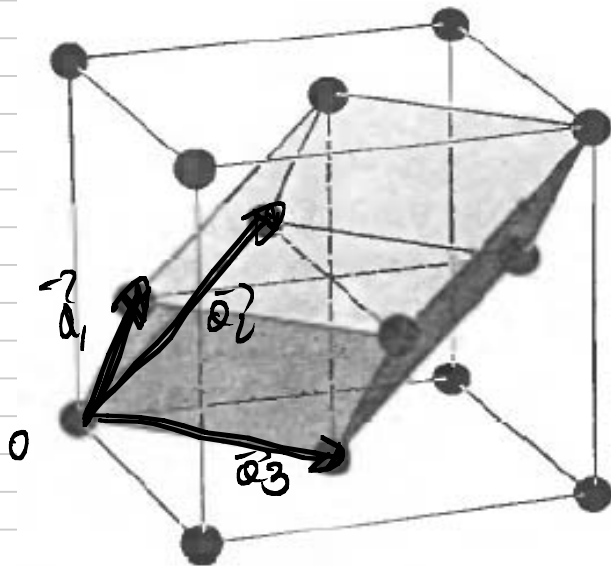
$$\vec{a}_2 = \frac{a}{2} (101)$$

$$\vec{a}_3 = \frac{a}{2} (110)$$

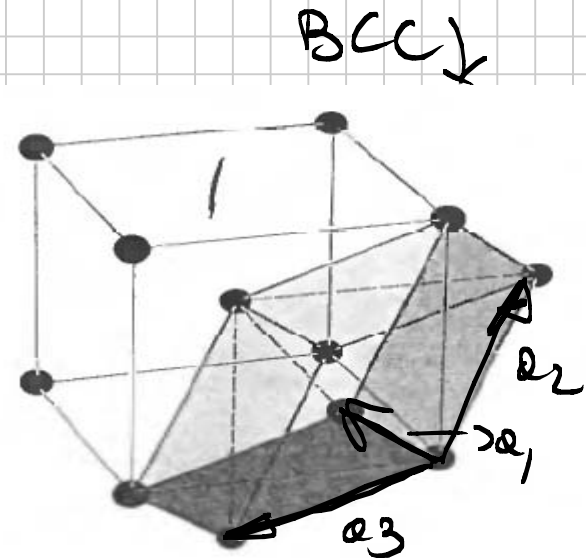
Ag
Au
Cu

PRIMITIVE UNIT CELL

= SOLID (AREA IN 2D) GENERATED
BY PRIMITIVE VECTORS



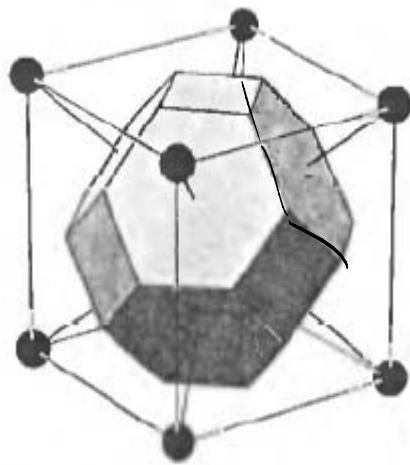
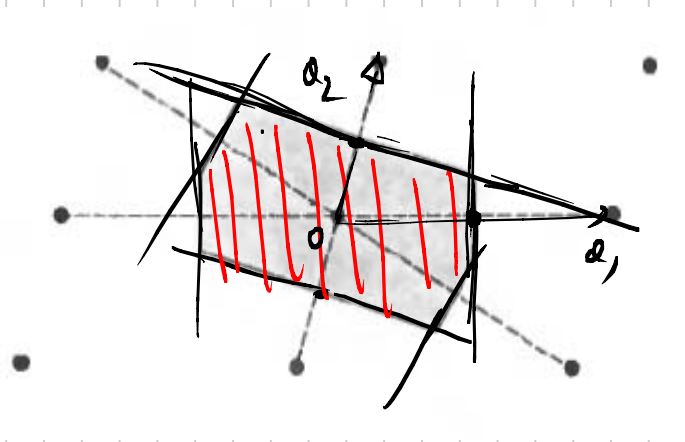
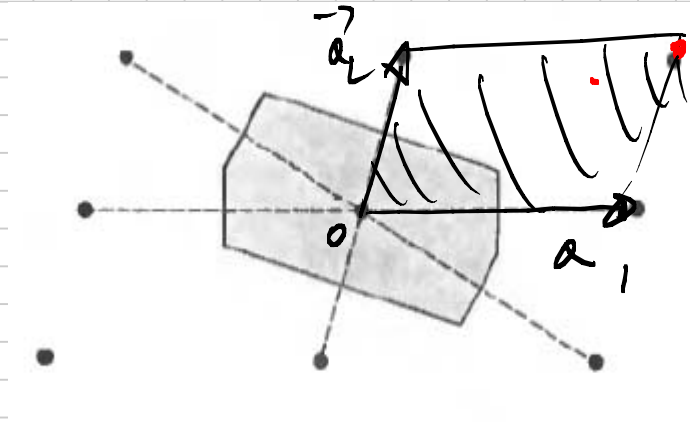
FCC



BCC

WIGNER-SEITZ CELLS

2D



WIGNER SEITZ
FOR BCC

TRUNCATED OCTAHEDRON

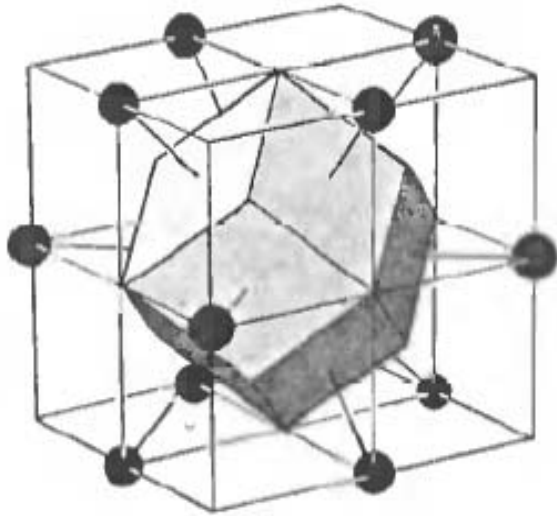
8 ATOMS ARE FIRST
NEIGHBORS \rightarrow CALLED

COORDINATION NUMBER

FCC

→ COORDINATION NUMBER

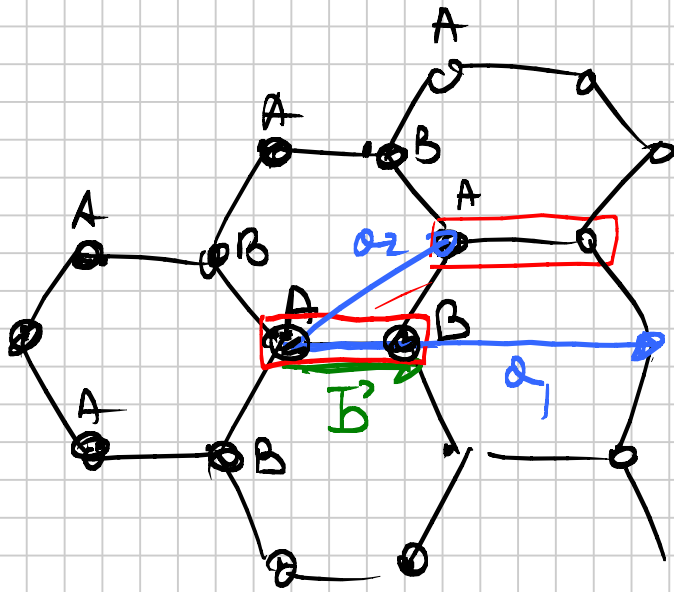
12



TRUNCATED
DODECAHEDRON

CRYSTAL STRUCTURES

BRAVAIS LATTICE
WITH BASIS



CONSIDER A AND B TOGETHER

\vec{a}_1 \vec{a}_2 PRIMITIVE VECTORS

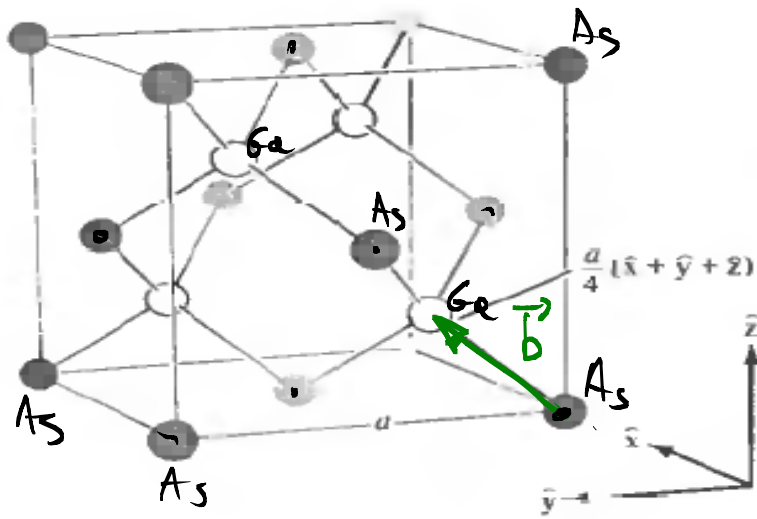
\vec{b} BASIS VECTOR

2D

3D LATTICE WITH

BASIS : DIAMOND

C, Si, Ge



FCC

$$\vec{b}_1 = \frac{a}{4}(111)$$

$$\vec{b}_2 = \frac{a}{2}(110)$$

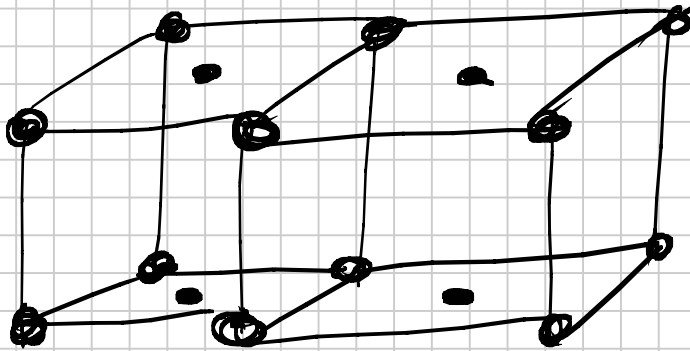
$$\vec{b}_3 = \frac{a}{2}(101)$$

$$\vec{b}_4 = \frac{a}{2}(011)$$

ATOM AT \vec{b} IS DIFFERENT : ZINC BLEND LATTICE

GeAs, CdTe

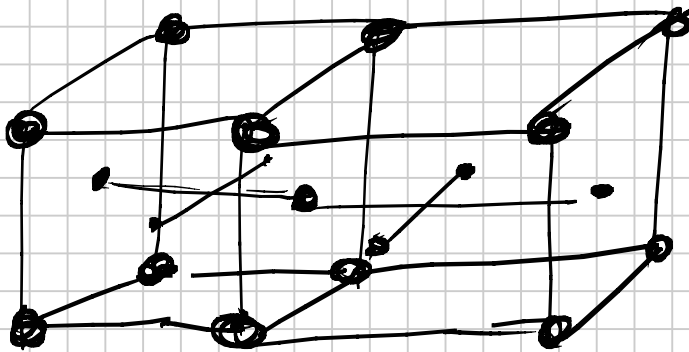
① BASE CENTERED CUBIC



BRAVAS LATTICE?

\vec{b} ?

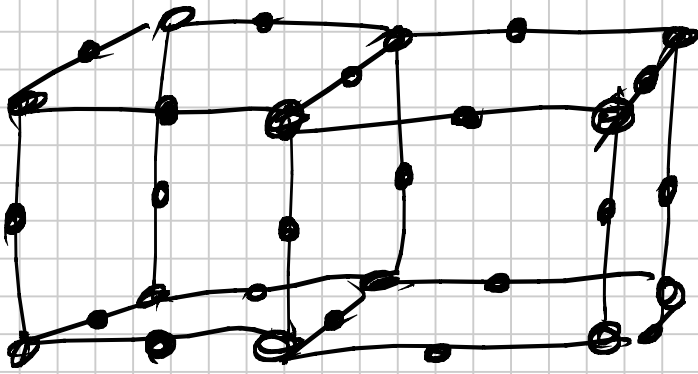
②



BRAVAS LATTICE?

\vec{b} ?

3



BRAVILS LATTICE?

$\vec{0}$?