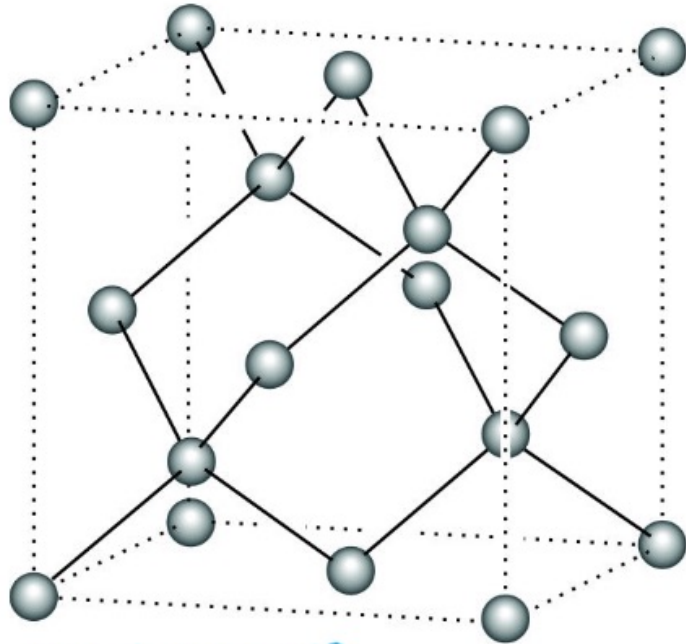


# CRYSTAL STRUCTURE



DIAMOND

## LATTICE

$$r = \mu_1 \vec{a}_1 + \mu_2 \vec{a}_2 + \mu_3 \vec{a}_3$$

$\mu_1, \mu_2, \mu_3 = \text{integers}$

$a_1, a_2, a_3 = \text{vectors}$

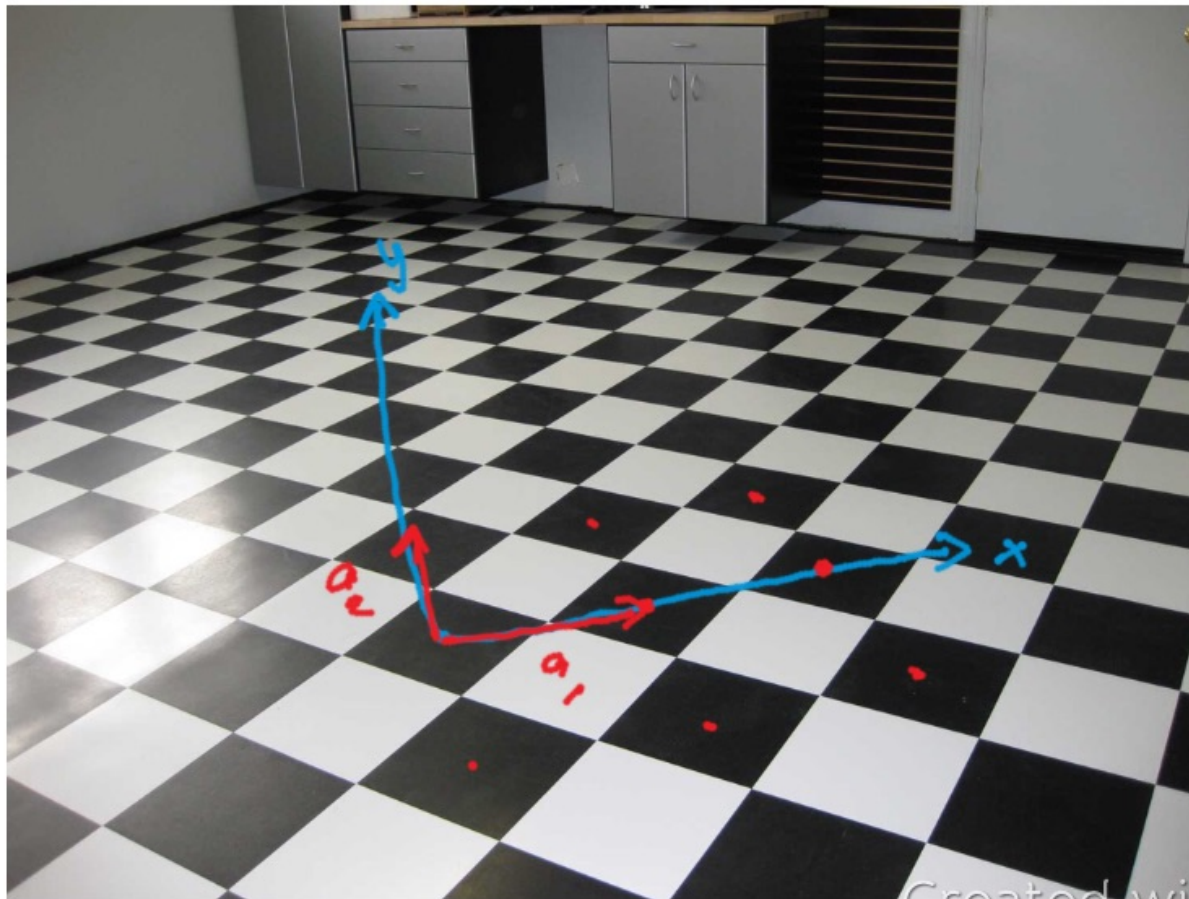
## BASIS

DESCRIBING  
"CHEMISTRY"

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# CHECKERBOARD - KITCHEN FLOOR:



## LATTICE

$$r = \mu_1 a_1 + \mu_2 a_2$$

$$a_1 = (1, 0)$$

$$a_2 = (0, 1)$$

## BASIS

◆ @ (0, 0)

◇ @  $(\frac{1}{2}, \frac{1}{2})$

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# CHECKERBOARD - KITCHEN FLOOR.




LATTICE


$$a_1 = (1, 0)$$

$$a_2 = (0, 1)$$

$$r = \mu_1 a_1 + \mu_2 a_2$$

BASIS

 @  $(0, 0)$   
 $(\frac{1}{2}, \frac{1}{2})$

 @  $(\frac{1}{2}, 0)$   
 $(0, \frac{1}{2})$

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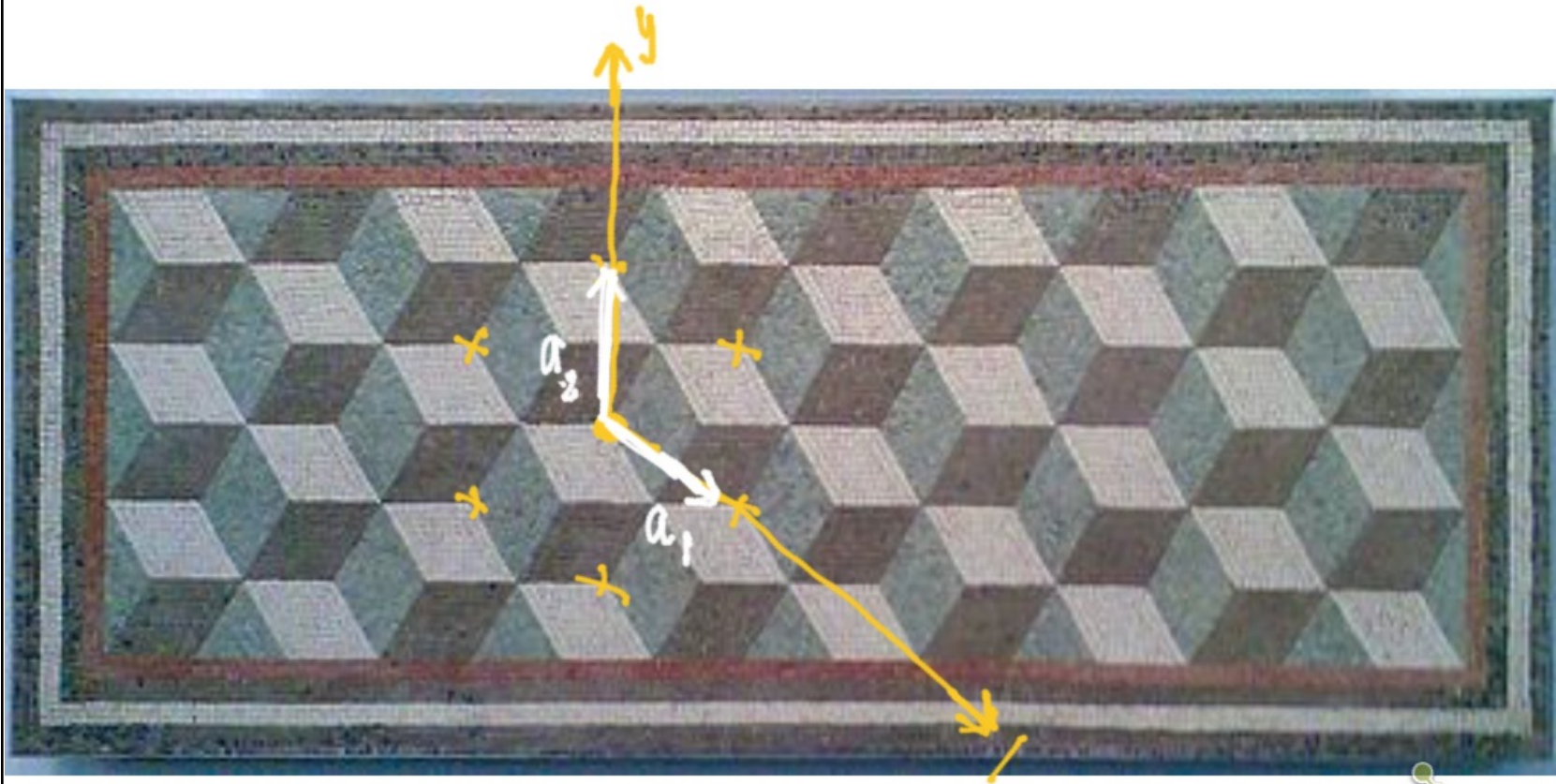
# M.C. ESCHER'S TESSELLATION



Created with Doceri



# ROMAN MOSAIC:



LATTICE

$$r = u_1 a_1 + u_2 a_2$$

BASIS =

Created with Doceri



# M.C. ESCHER'S TESSELLATION



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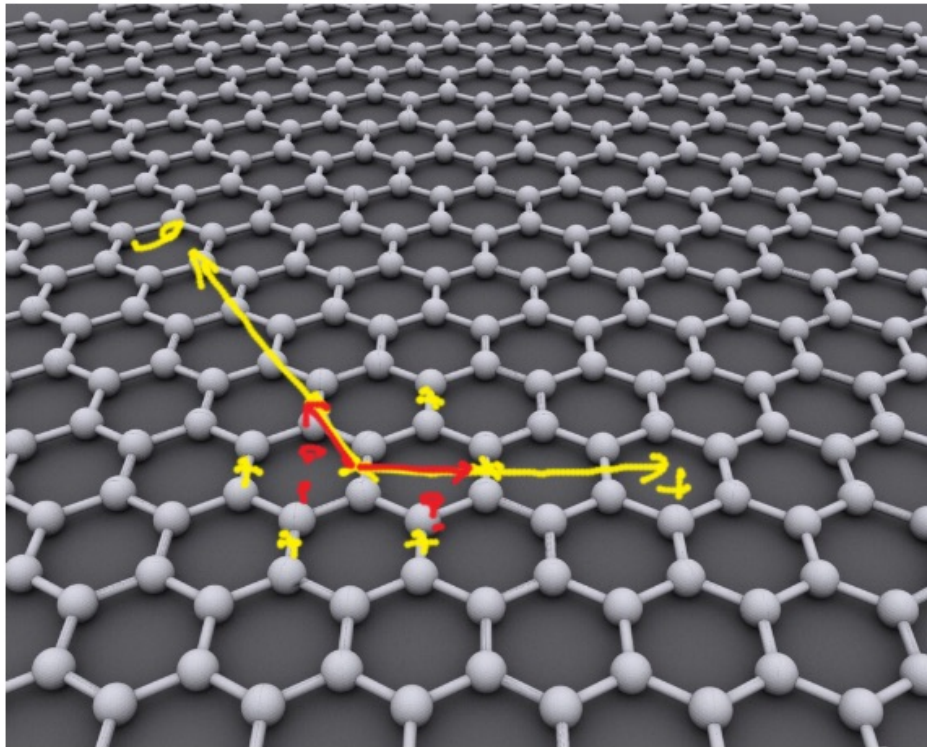


# HONEYCOMB:



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# GRAPHENE:

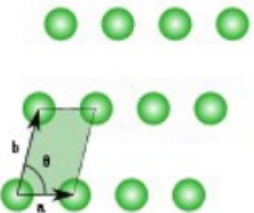
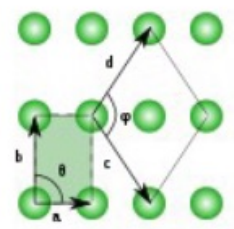
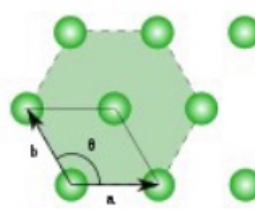
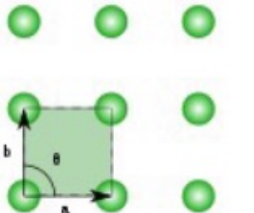


Two C-atoms  
in the BASIS.

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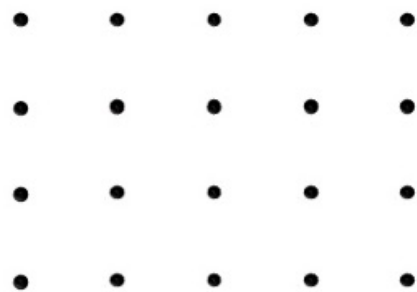
# POSSIBLE TWO-DIMENSIONAL LATTICES

 <p>1</p>	 <p>2</p>	 <p>4</p>	 <p>5</p>
<p><math> a  \neq  b , \theta = 90^\circ</math></p> <p><b>m</b></p>	<p><math> a  \neq  b , \theta = 90^\circ</math>  <math> c  =  d , \phi = 90^\circ</math></p> <p><b>o</b></p>	<p><math> a  =  b , \theta = 120^\circ</math></p> <p><b>h</b></p>	<p><math> a  =  b , \theta = 90^\circ</math></p> <p><b>t</b></p>

# BRAVAIS LATTICES (2D)



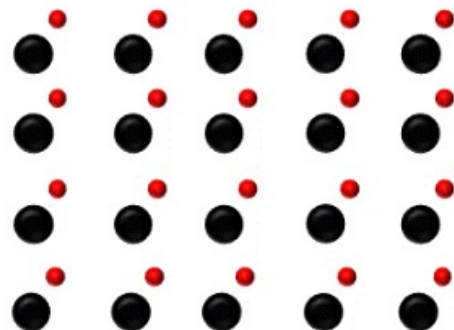
# LATTICE + BASIS = CRYSTAL STRUCTURE



Space Lattice



Basis (atoms)

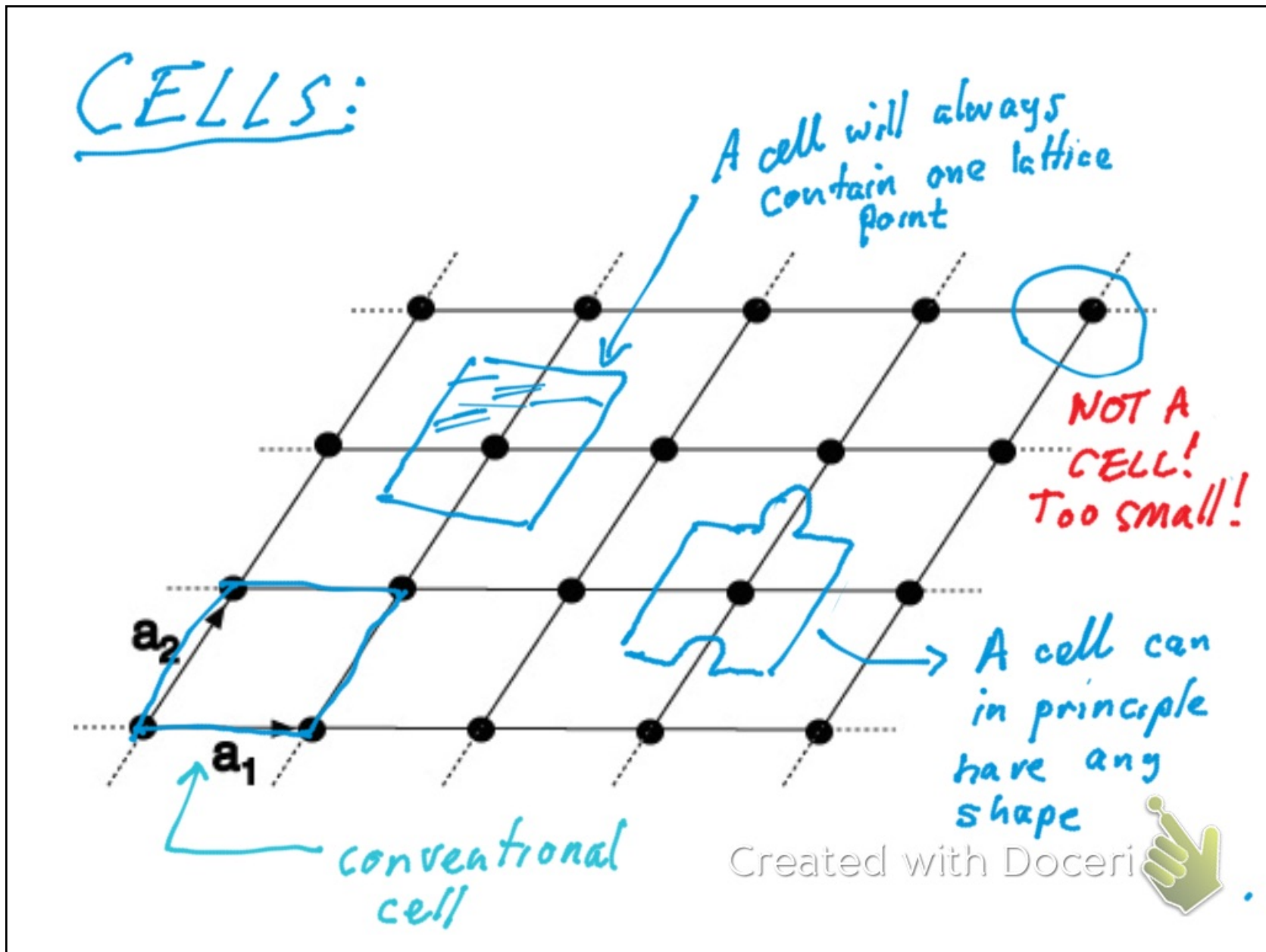


Crystal structure

The crystal structure is formed by adding basis (atoms) to every lattice points of the lattice. The number of atoms in the basis may be one or more than one.

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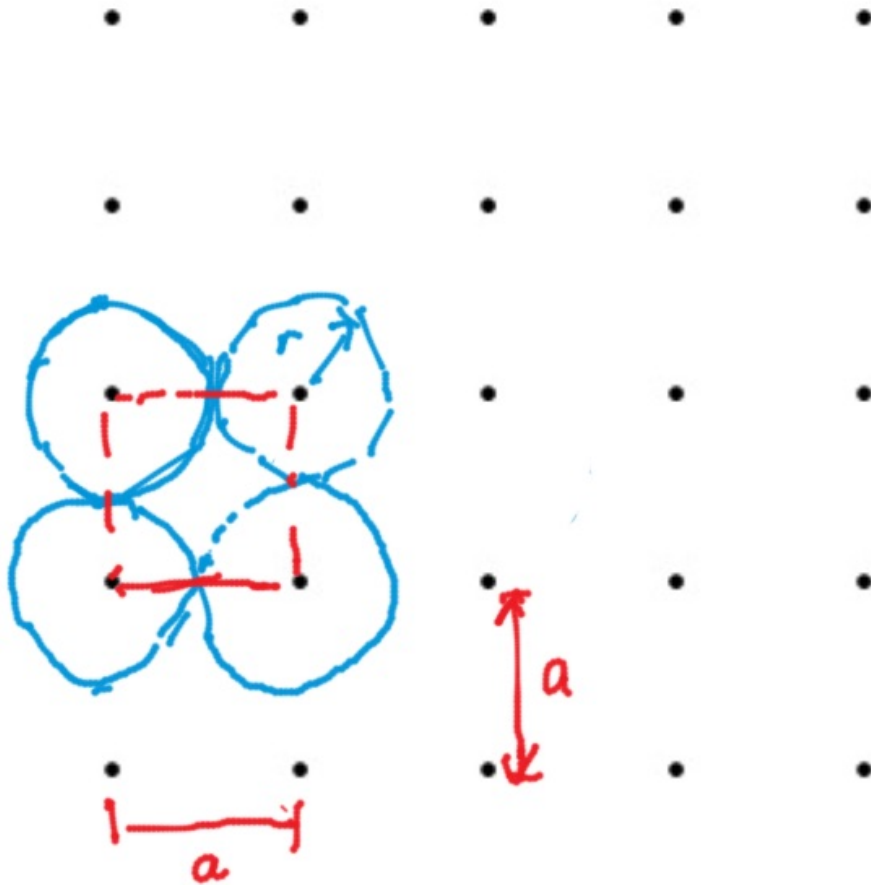
# WIGNER-SEITZ CELL:

"PIZZA DELIVERY STRUCTURE"

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# PACKING RATIOS

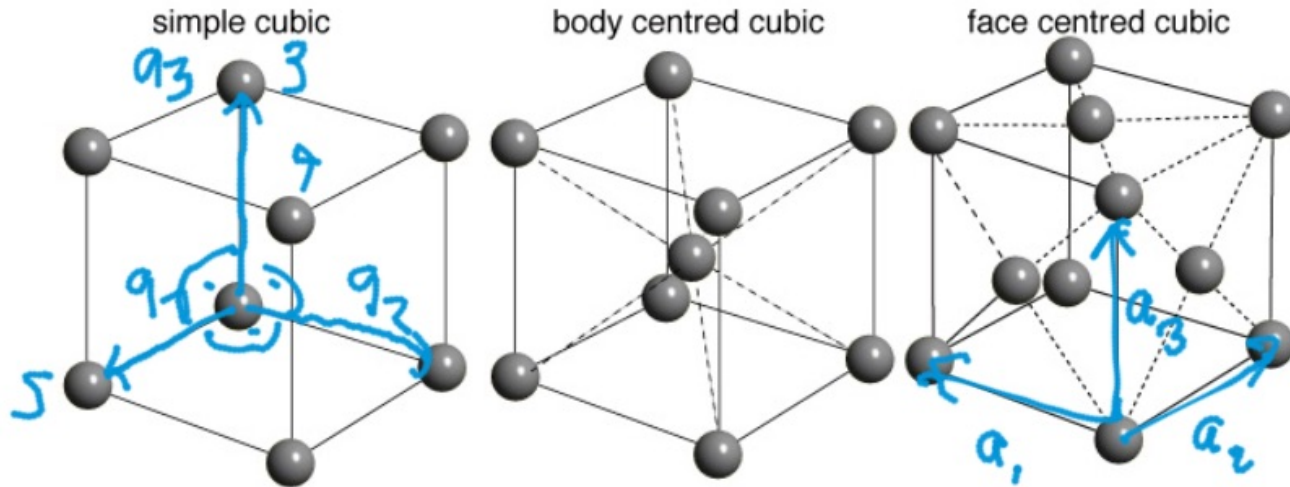


$$\begin{aligned}
 \rho\text{-ratio} &= \frac{\text{Area Circle}}{\text{Cell area}} \\
 &= \frac{\pi \cdot r^2}{a^2} \\
 &= \frac{\pi}{4}
 \end{aligned}$$

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# CUBIC LATTICES:

BASIS-BCC  
 $(0, 0, 0)$   
 $(\frac{1}{2}, \frac{1}{2}, \frac{1}{2})$



$|a_1| = |a_2| = |a_3|$

## FCC-BASIS

$(0, 0, 0)$   
 $(\frac{1}{2}, \frac{1}{2}, 0)$   
 $(\frac{1}{2}, 0, \frac{1}{2})$   
 $(0, \frac{1}{2}, \frac{1}{2})$

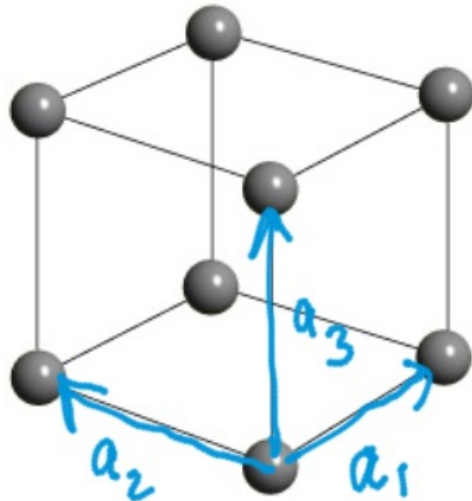
	SC	bcc	fcc
lattice points:	1	2	4
nearest neighbours:	6	8	12

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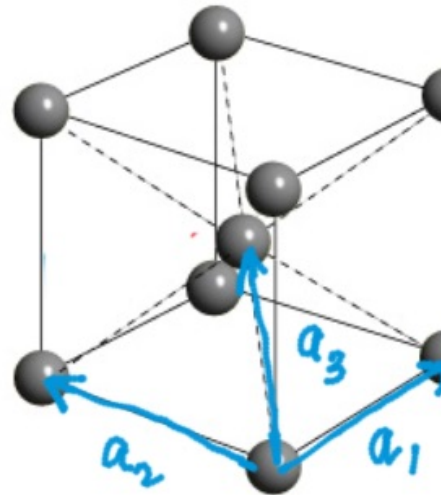


# CUBIC LATTICES:

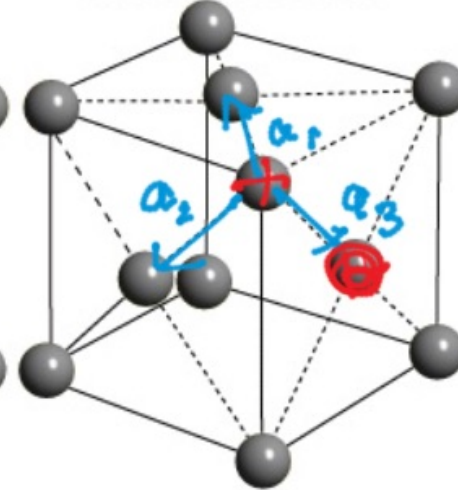
simple cubic



body centred cubic



face centred cubic



## LATTICE VECTORS

$$a_1 = (1, 0, 0)$$

$$a_2 = (0, 1, 0)$$

$$a_3 = (0, 0, 1)$$

BASIS:

$$(0, 0, 0)$$

$$a_1 = (1, 0, 0)$$

$$a_2 = (0, 1, 0)$$

$$a_3 = \left(\frac{1}{2}, \frac{1}{2}, \frac{1}{2}\right)$$

BASIS

$$(0, 0, 0)$$

$$a_1 = \left(\frac{1}{2}, \frac{1}{2}, 0\right)$$

$$a_2 = \left(0, \frac{1}{2}, \frac{1}{2}\right)$$

$$a_3 = \left(\frac{1}{2}, 0, \frac{1}{2}\right)$$

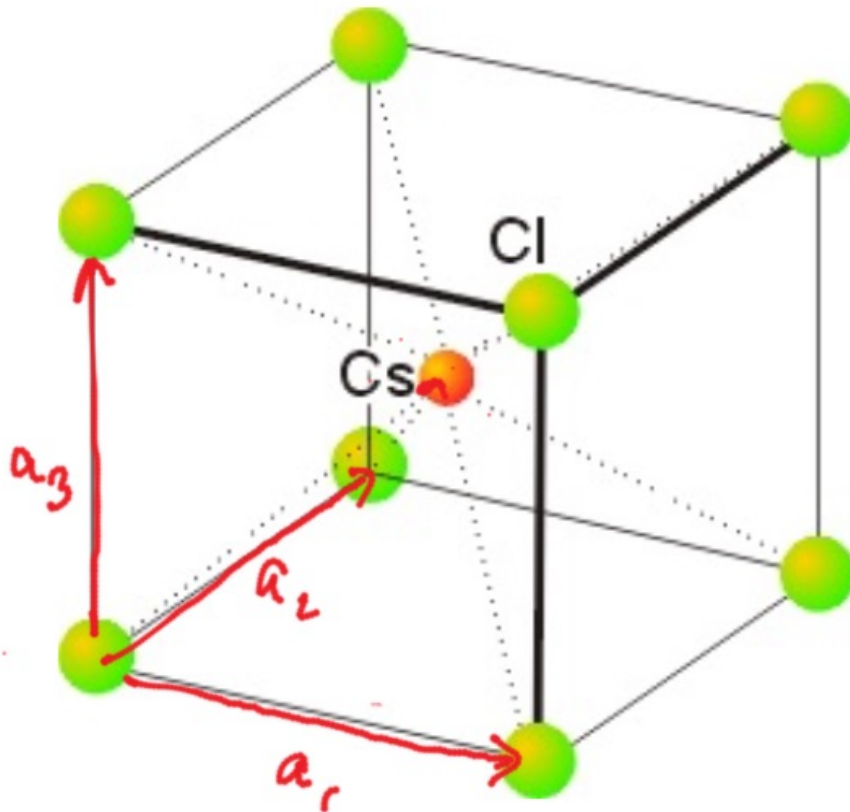
BASIS

$$(0, 0, 0)$$

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# QUIZ 1:



(1) SIMPLE CUBIC ✓

(2) BCC

(3) FCC

LATTICE = SC

BASIS

Cl @ (0, 0, 0)

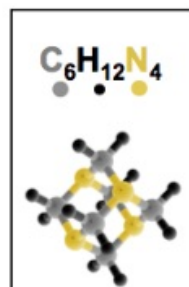
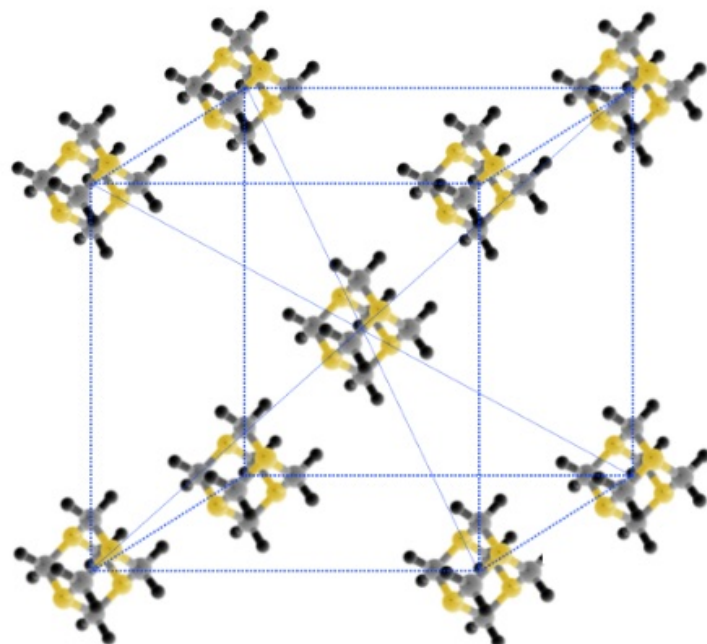
Cs @  $(\frac{1}{2}, \frac{1}{2}, \frac{1}{2})$

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# QUIZ 2



(1) SIMPLE CUBIC

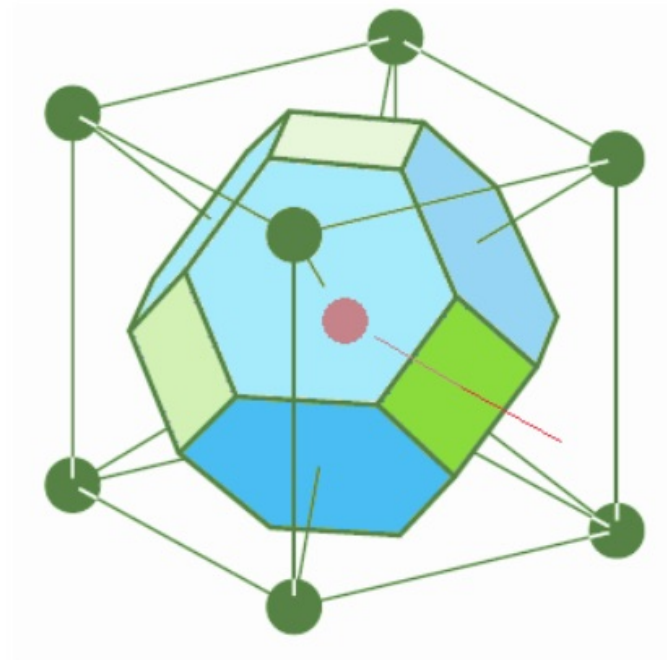
(2) BCC ✓

(3) FCC

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# WIGNER-SEITZ CELL:

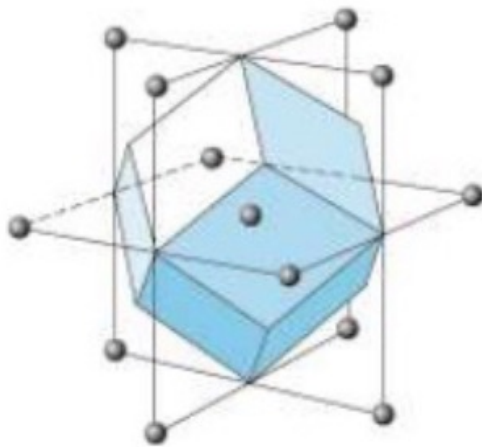


Which lattice?

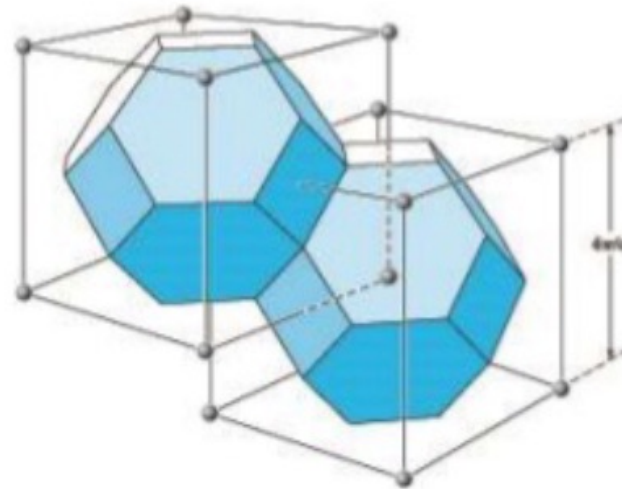
Created with Doceri



## Wigner-Seitz Cell - 3D



fcc wigner-seitz cell



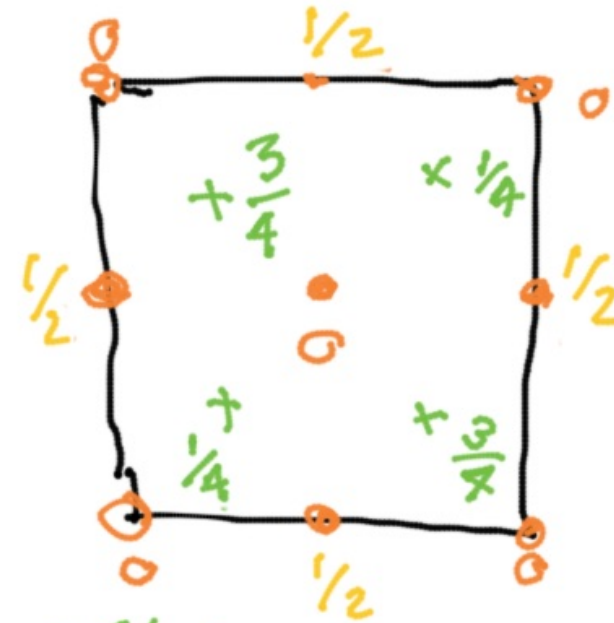
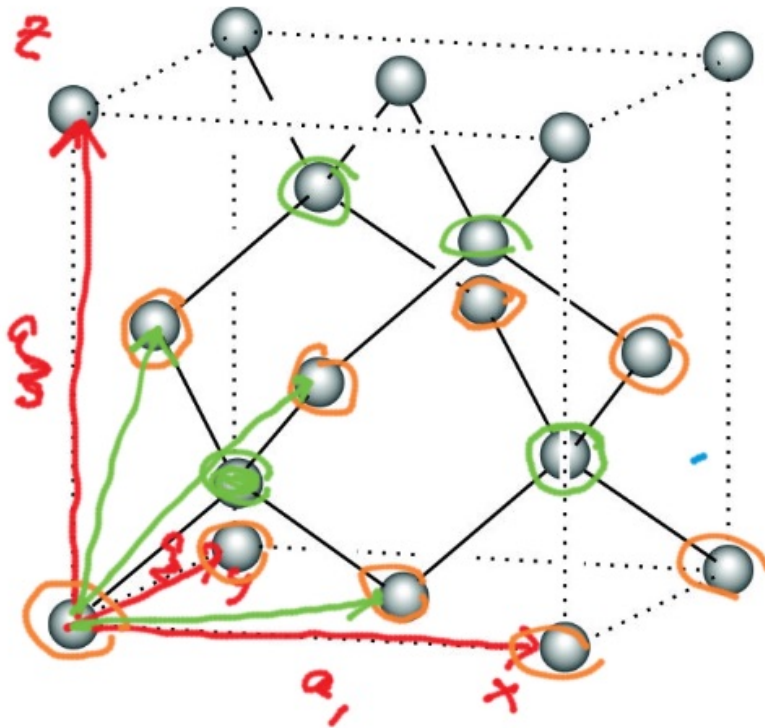
bcc wigner-seitz cell

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Solid State Physics

# DIAMOND STRUCTURE:



$$a_1 = (\frac{1}{2}, \frac{1}{2}, 0)$$

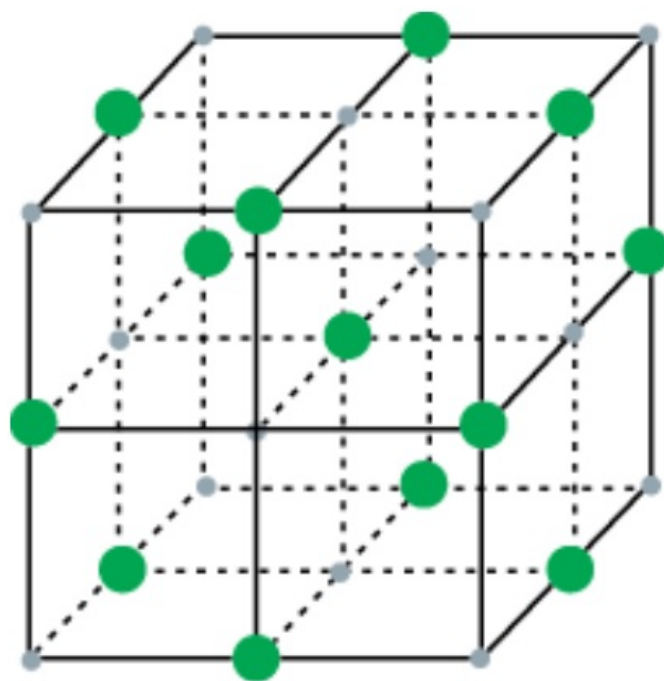
$$a_2 = (\frac{1}{2}, 0, \frac{1}{2})$$

$$a_3 = (0, \frac{1}{2}, \frac{1}{2})$$

BASIS = C at  $(0,0,0)$  and  $(\frac{1}{4}, \frac{1}{4}, \frac{1}{4})$



# TABLE-SALT STRUCTURE:



● Cl<sup>-</sup>    • Na<sup>+</sup>

FCC - Lattice

TWO ATOM  
BASIS

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