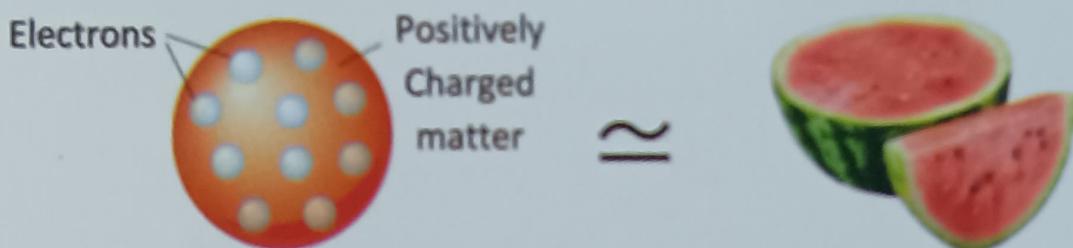


# STRUCTURE OF ATOM

e	$-1.6 \times 10^{-19}$ C	$9.1 \times 10^{-31}$ kg
p	$+1.6 \times 10^{-19}$ C	$1.672 \times 10^{-27}$ kg
n	Neutral	$1.674 \times 10^{-27}$ kg

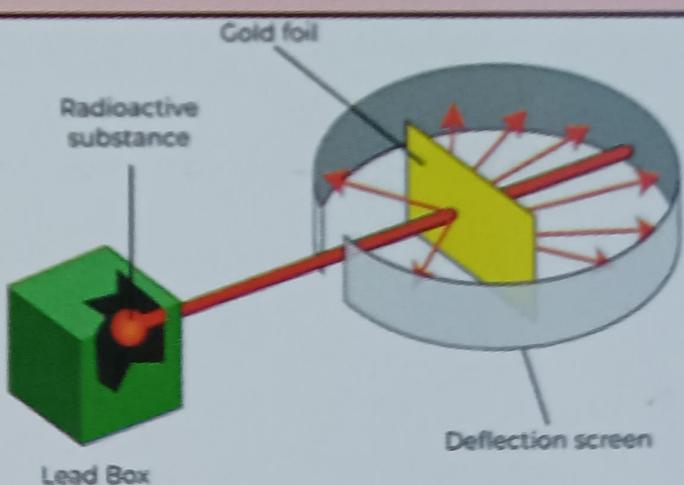
## ATOMIC MODELS

### THOMSON ATOMIC MODEL (Plum/Watermelon)



### RUTHERFORD'S ATOMIC MODEL

- Most of the space in the atom is empty.
- The positive charge is concentrated in a very small volume.
- The volume of the nucleus is very small compared to the total volume of the atom.



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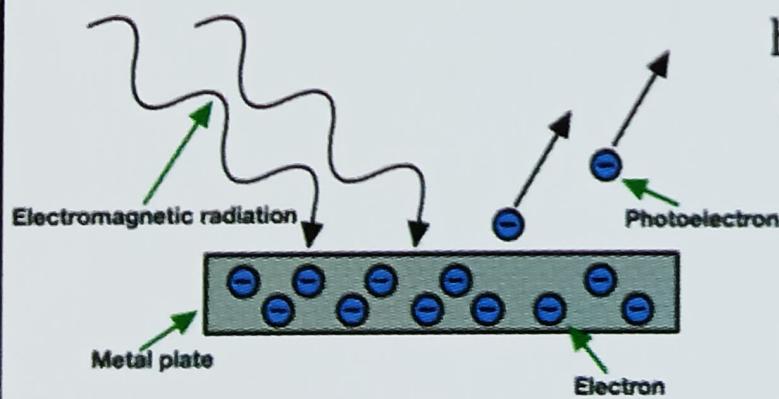
## Important terms

<b>Isoelectronic</b> : Same no. of electrons	CH <sub>4</sub> , NH <sub>3</sub> , H <sub>2</sub> O
<b>Isotopes</b> : Same atomic no. different mass no.	<sub>1</sub> H <sup>1</sup> , <sub>1</sub> H <sup>2</sup> , <sub>1</sub> H <sup>3</sup>
<b>Isobars</b> : same mass no. different atomic no.	<sub>18</sub> Ar <sup>40</sup> , <sub>20</sub> Ca <sup>40</sup>
<b>Isotones</b> : same number of neutrons	<sub>6</sub> C <sup>13</sup> , <sub>7</sub> N <sup>14</sup>

### Photoelectric Effect

- No. of Photoelectrons  $\propto$  Intensity of Light
- Kinetic Energy  $\propto$  Frequency of Light

Emission of electron from the surface of metal when a photon of certain  $v$  is incident to metal surface.

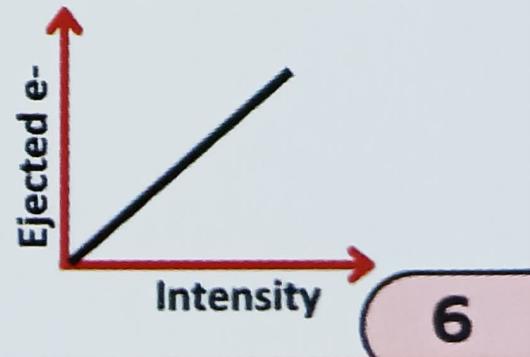
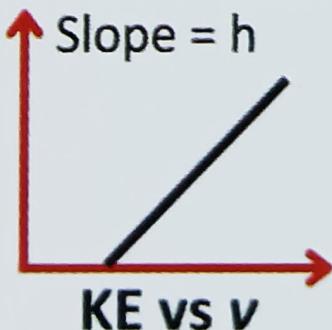
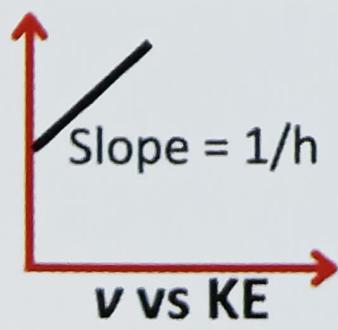


$$h\nu = h\nu_0 + \frac{1}{2}mv^2$$

$$h\nu = w_0 + \frac{1}{2}mv^2$$

w<sub>0</sub> = Work Function  
ν<sub>0</sub> = Threshold Freq.

After Threshold, it's all Kinetic Energy



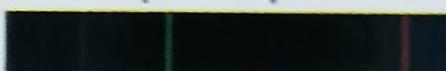
## Emission Spectrum of Hydrogen Atom



Continuous spectrum



Absorption spectrum



Emission spectrum

$$\bar{v} = \frac{1}{\lambda} = R_H \left[ \frac{1}{n_1^2} - \frac{1}{n_2^2} \right] \times Z^2$$

$n_1$  &  $n_2$  = energy levels of transitions  
 $R_H$  = Rydberg Constant  
 $Z$  = atomic number

Number of Spectral Lines

$$\frac{(n_2 - n_1)(n_2 - n_1 + 1)}{2}$$

### Spectrum Lines of Hydrogen Atom

Series	$n_1$	$n_2$	Spectral Reg.
Lymann	1	2,3..	U.V.
Balmer	2	3,4..	Visible
Paschen	3	4,5..	IR
Brackett	4	5,6..	IR
Pfund	5	6,7..	IR

### Bohr's Theory

- Electron in circular orbits
- Energy of electron doesn't change in an orbit.
- Energy is absorbed or emitted in the difference of  $\Delta E = h\nu = hc/\lambda$

## Bohr model Formulas

**Angular Momentum is quantised; n = 1,2,3...**

$$mvr = n \frac{h}{2\pi}$$

**Bohr's Radius**

$$r_n = 0.529 \times \frac{n^2}{Z} \text{ Å}^\circ$$

**Energy of Electron**  
**T.E. = -K.E. = P.E/2**

$$E_n = -13.6 \times \frac{Z^2}{n^2} \text{ eV}$$

$$E_n = -2.18 \times 10^{-18} \times \frac{Z^2}{n^2} \text{ J}$$

## De-Broglie Hypothesis

$$\lambda = \frac{h}{p} = \frac{h}{mv} = \frac{h}{\sqrt{2m(KE)}} = \frac{h}{\sqrt{2mqV}} = \frac{12.24}{\sqrt{V}} \text{ Å}^\circ$$

V = Potential; q = charge; KE = Kinetic Energy

## Heisenberg Uncertainty Principle

$$\Delta x \cdot \Delta p \geq \frac{h}{4\pi}$$

$\Delta x$  = Change in Position

$\Delta p$  = Change in Momentum

Wolfgang Pauli  
 & Neil Bohr  
 OBSERVING  
 SPIN 1954

# Quantum Numbers

## Principal quantum number (n)

- Positive integer with values  $n = 1, 2, 3, 4, \dots$
- $n$  denotes the **shell number, energy of electron, size of the shell.**
- Maximum number of electron in shell =  $2n^2$
- Maximum number of orbitals in shell =  $n^2$

## Azimuthal Quantum number (l)

- Positive integer with values  $l = 0, 1, 2, 3, \dots$
- $l$  denotes the **subshell, suborbit, sub energy level.**
- value of  $l$  ranges from 0 to  $n-1$

$l$	0	1	2	3
orbital	s	p	d	f

## Magnetic quantum Number (m)

- Positive integer depending upon value of  $l$  as  $m$  ranges from  $-l$  to  $+l$  (total =  $2l+1$  values)
- It tells about the orientation of the orbital.  
eg : p orbital has  $l = 1$  ( $m = -1, 0, +1$ )

## Spin Quantum Number (s)

- Electron rotates on its own axis in clockwise or anticlockwise direction. So the spin quantum number can have two values of  $+1/2$  and  $-1/2$ .

## Energy in different electron systems

Monoelectronic	Polyelectronic
<p>Degeneracy = <math>n^2</math></p>	<p>No Degeneracy</p>

## Electronic configuration rules

**Aufbau Principle :** Ground state electrons filled into atomic orbitals in the increasing order of orbital energy level.

**Hunds Rule :** Before the double occupation of any orbital, every orbital in the sub level is singly occupied.

**Pauli's Principle :** In single atom no two electrons will have an identical set of quantum numbers ( $n$ ,  $l$ ,  $m$ , and  $m_s$ ).

## Exceptions

- Chromium : [Ar]4s<sup>1</sup>3d<sup>5</sup> (Half Filled configuration)
- Copper : [Ar]4s<sup>1</sup>3d<sup>10</sup> (Fully Filled configuration)

**Nodes** ( $n$ =Principle Qno. ;  $l$  = Azimuthal Qno.)

Radial node =  $n-l-1$  ; Angular Node =  $l$

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