



PRACHAND NEET



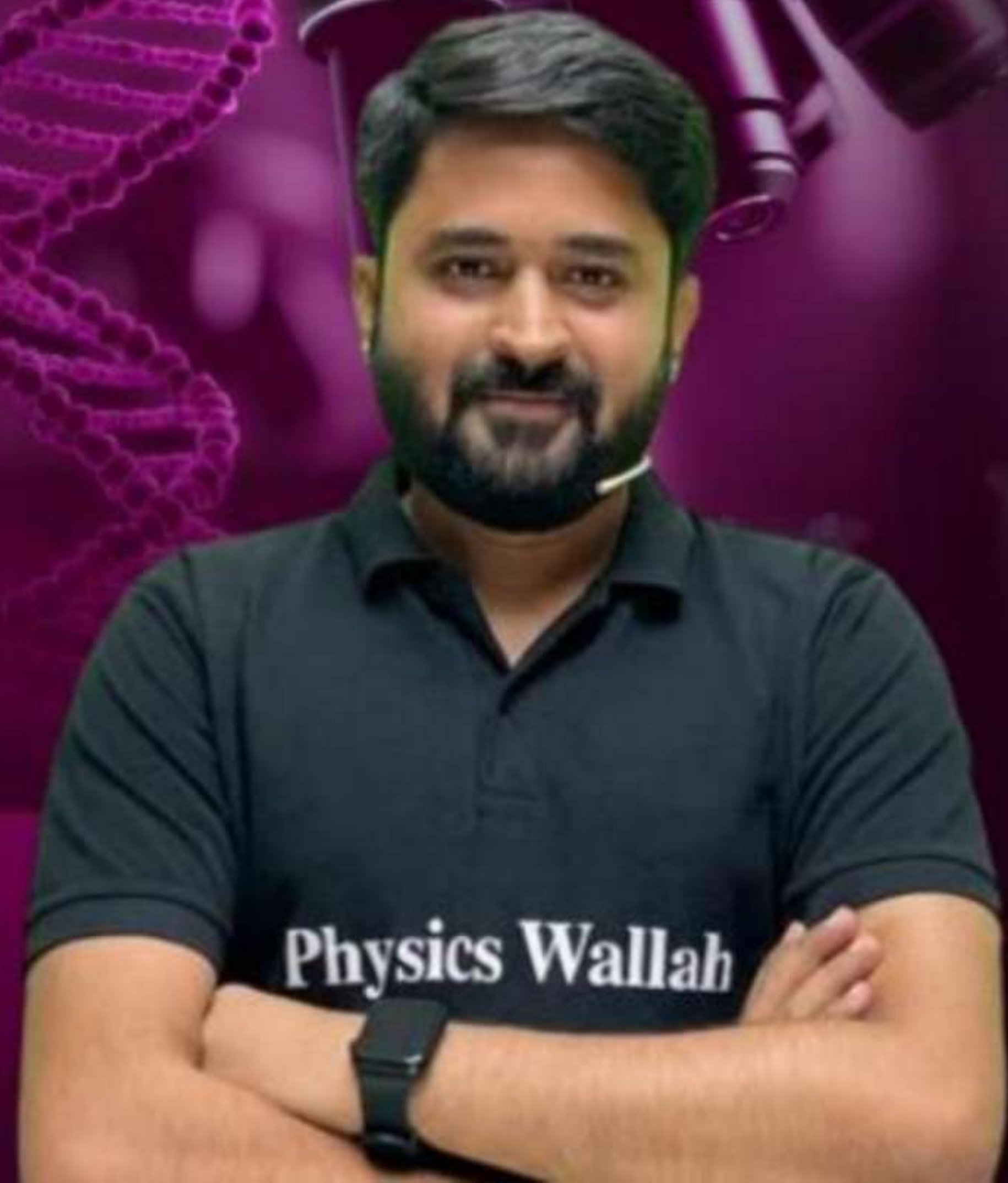
ONE SHOT



Inorganic Chemistry

Classification of Elements and Periodicity in Properties

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Topics

to be covered



Periodic Table One Shot



s-block																		p-block															
1 1A H Hydrogen	2 2A He Helium																	3 3A B Boron	4 4A C Carbon	5 5A N Nitrogen	6 6A O Oxygen	7 7A F Fluorine	8 8A Ne Neon										
3 Li Lithium	4 Be Beryllium																	5 Al Aluminium	6 Si Silicon	7 P Phosphorus	8 S Sulfur	9 Cl Chlorine	10 Ar Argon										
11 Na Sodium	12 Mg Magnesium																	13 Ga Gallium	14 Ge Germanium	15 As Arsenic	16 Se Selenium	17 Br Bromine	18 Kr Krypton										
19 K Potassium	20 Ca Calcium	21 Sc Scandium	22 Ti Titanium	23 V Vanadium	24 Cr Chromium	25 Mn Manganese	26 Fe Iron	27 Co Cobalt	28 Ni Nickel	29 Cu Copper	30 Zn Zinc	31 Ga Gallium	32 Ge Germanium	33 As Arsenic	34 Se Selenium	35 Br Bromine	36 Kr Krypton																
37 Rb Rubidium	38 Sr Strontium	39 Y Yttrium	40 Zr Zirconium	41 Nb Niobium	42 Mo Molybdenum	43 Tc Technetium	44 Ru Ruthenium	45 Rh Rhodium	46 Pd Palladium	47 Ag Silver	48 Cd Cadmium	49 In Indium	50 Sn Tin	51 Sb Antimony	52 Te Tellurium	53 I Iodine	54 Xe Xenon																
55 Cs Caesium	56 Ba Barium	57 La Lanthanum	72 Hf Hafnium	73 Ta Tantalum	74 W Tungsten	75 Re Rhenium	76 Os Osmium	77 Ir Iridium	78 Pt Platinum	79 Au Gold	80 Hg Mercury	81 Tl Thallium	82 Pb Lead	83 Bi Bismuth	84 Po Polonium	85 At Astatine	86 Rn Radon																
87 Fr Francium	88 Ra Radium	89 Ac Actinium	104 Rf Rutherfordium	105 Db Dubnium	106 Sg Seaborgium	107 Bh Bohrium	108 Hs Hassium	109 Mt Meitnerium	110 Ds Darmstadtium	111 Rg Roentgenium	112 Cn Copernicium	113 Nh Nihonium	114 Fl Flerovium	115 Mc Moscovium	116 Lv Livermorium	117 Ts Tennessine	118 Og Oganesson																
119 Uue Ununennium	120 Ubn Unbinilium	121 Ubu Unbiunium																															
		Lanthanides		57 La Lanthanum	58 Ce Cerium	59 Pr Praseodymium	60 Nd Neodymium	61 Pm Promethium	62 Sm Samarium	63 Eu Europium	64 Gd Gadolinium	65 Tb Terbium	66 Dy Dysprosium	67 Ho Holmium	68 Er Erbium	69 Tm Thulium	70 Yb Ytterbium	71 Lu Lutetium															
		Actinides		89 Ac Actinium	90 Th Thorium	91 Pa Protactinium	92 U Uranium	93 Np Neptunium	94 Pu Plutonium	95 Am Americium	96 Cm Curium	97 Bk Berkelium	98 Cf Californium	99 Es Einsteinium	100 Fm Fermium	101 Md Mendelevium	102 No Nobelium	103 Lr Lawrencium															
Alkali metals		Alkaline earth metals		Lanthanides		Actinides		Transition metals		Post-transition metals		Metalloids		Reactive nonmetals		Halogens		Noble gases		Unknowns													

3/IIIB

d-block

f-block

28 2p⁶
3s 3p
4s 3d 4p

32 elements

La

Ac

La

Ac

B

He

Mn

Te

Po

Cn

Ts

Og

Uue

Ubn

Ubu

Lanthanides

Actinides

Alkali metals Alkaline earth metals Lanthanides Actinides Transition metals Post-transition metals Metalloids Reactive nonmetals Halogens Noble gases Unknowns

actinoids





INTRODUCTION



It is a systematic arrangement of all known elements in increasing order of their atomic No. (Z) in horizontal rows (periods) & vertical columns (groups) so that the properties of the elements can be studied easily.

* Mendeleev's Periodic law :-

Physical & Chemical properties of
the elements are periodic function of their Atomic Weight.







MODERN PERIODIC TABLE

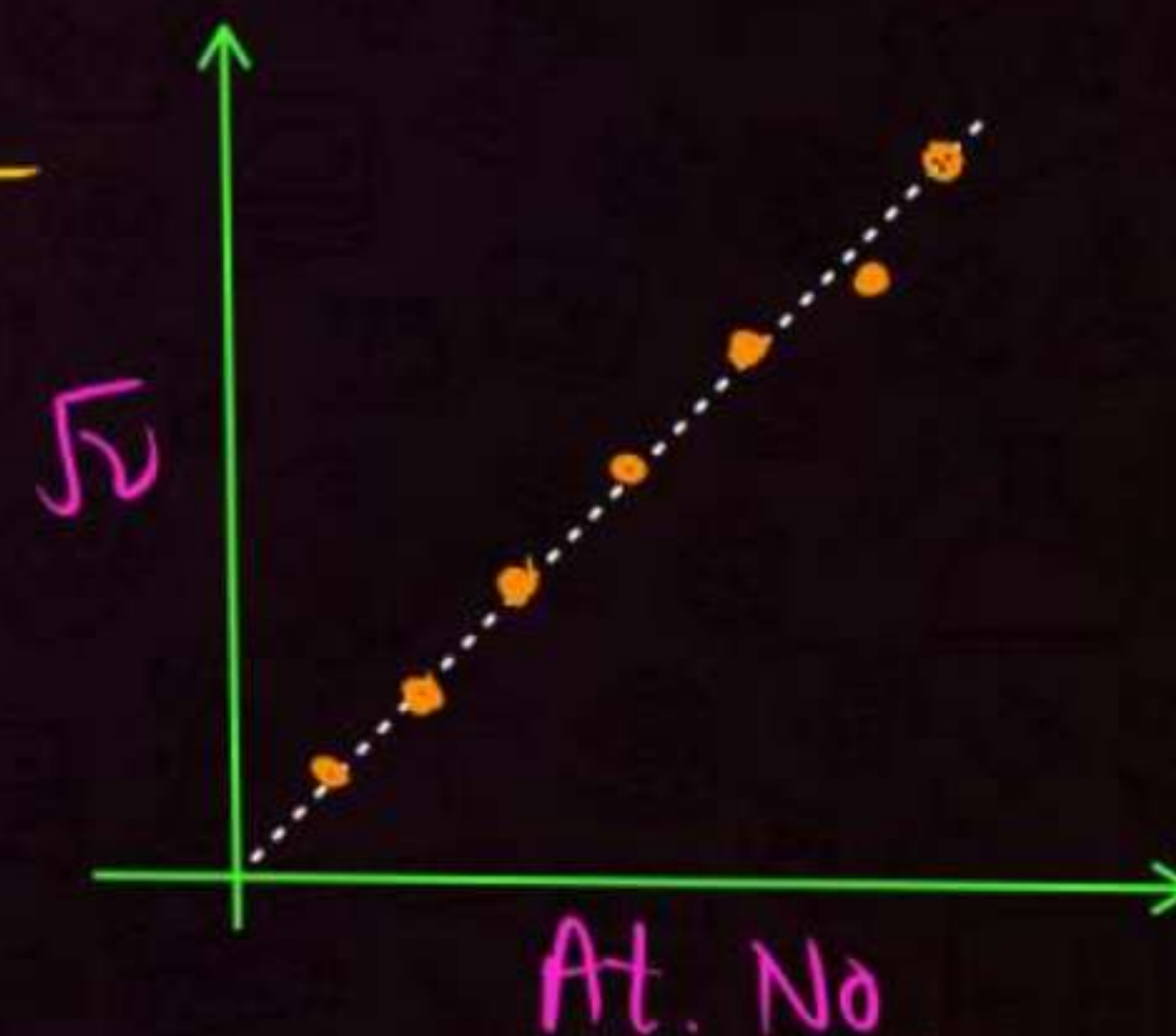
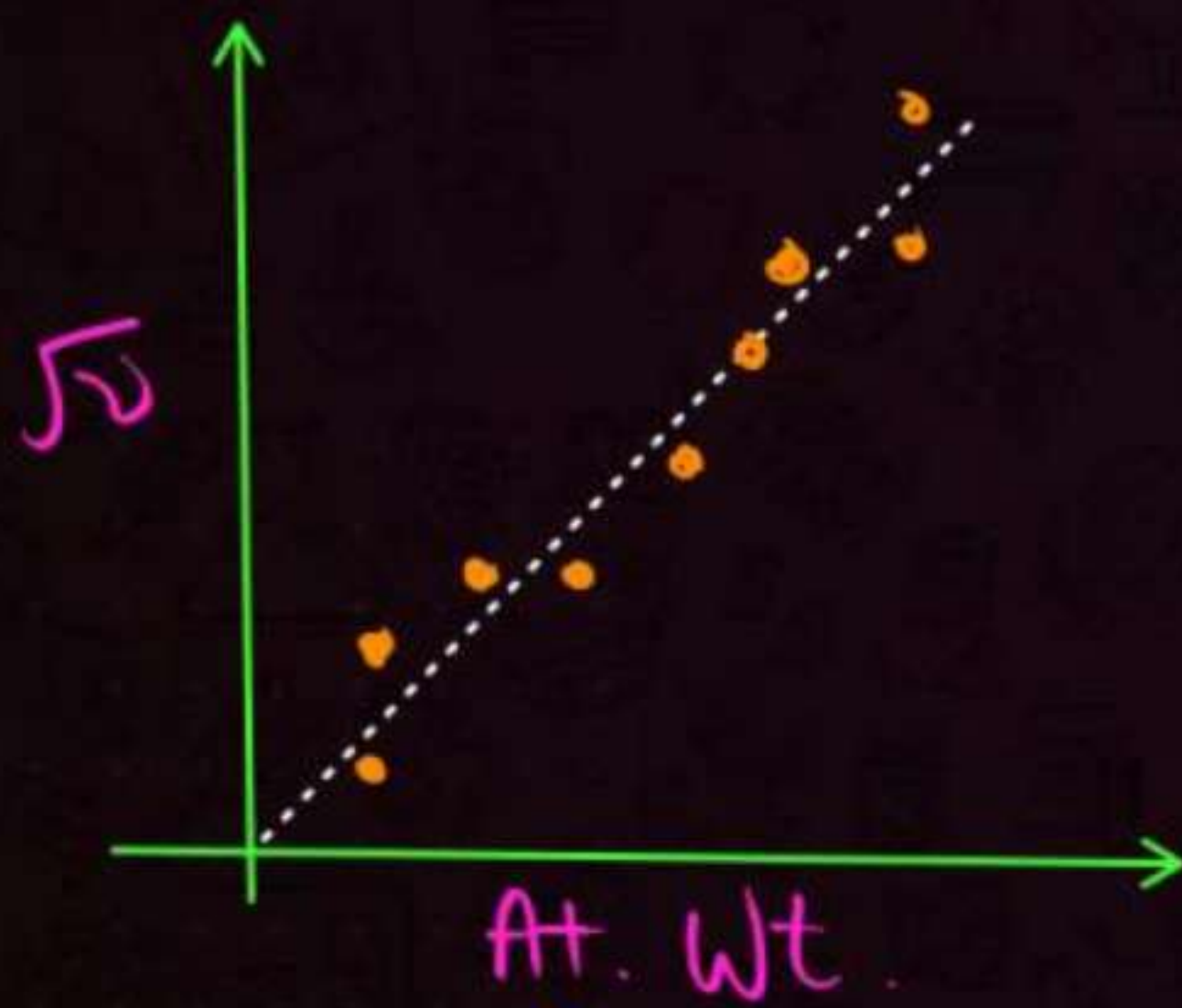


* Moseley's Experiment $\rightarrow$

β particles

X-Rays (ν)

M



* Modern Periodic law:-

Physical & Chemical properties of the elements are periodic function of their Atomic Numbers.





LONG FORM PERIODIC TABLE



periods = 7

Groups = 18 (1 to 18)

Blocks = 4 (s, p, d, f)

* Representative Elements $\rightarrow$ all s & p-block elements including Noble gases are called Representative Elements.
(Acc. to New NCERT)

* Transition Elements $\rightarrow$ d-block elements (except group -12) are c/d transition elements.



* Note $\rightarrow$ (n-1)d subshell incomplete either in their ground state or any stable oxidation state.

General e⁻ config of d-block elements : $n s^{1-2} (n-1) d^{1-10}$ ✓

Pd $\rightarrow [Kr] 5s^0 4d^{10}$

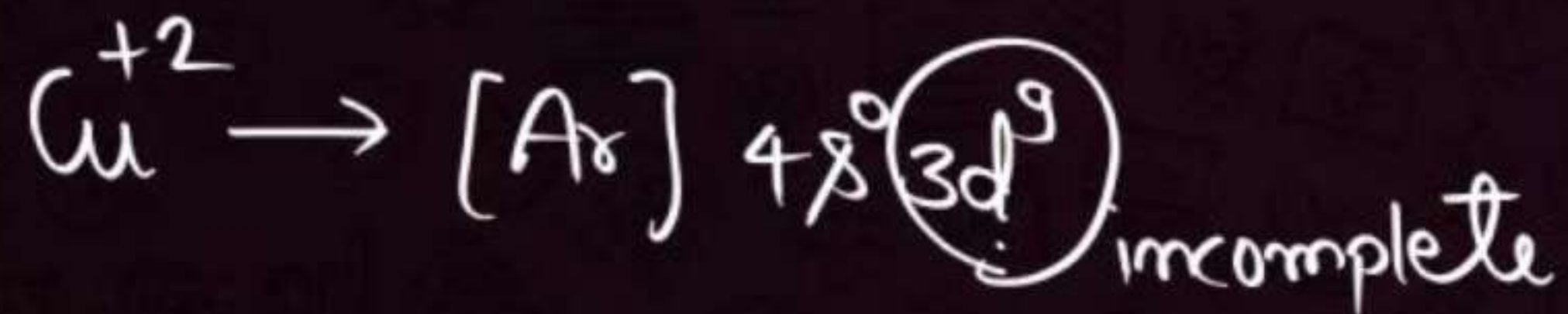
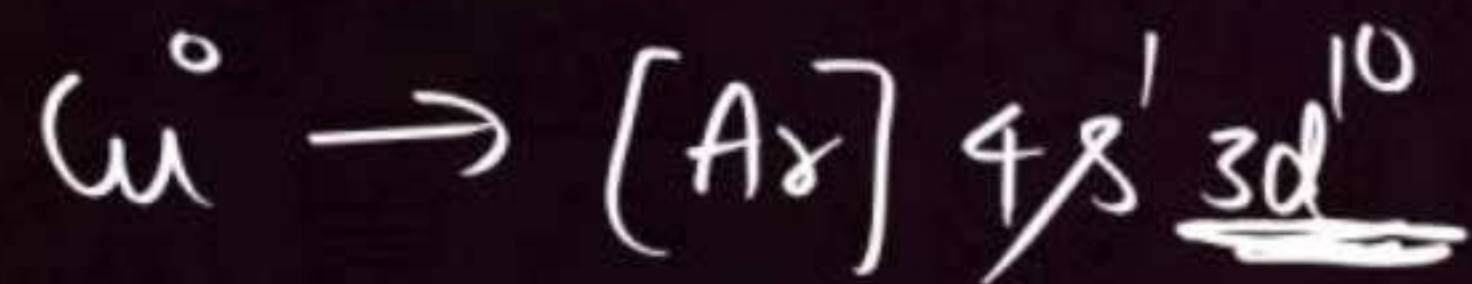
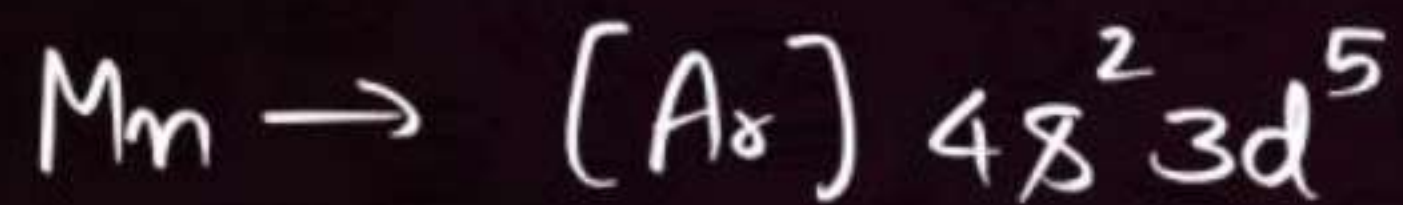
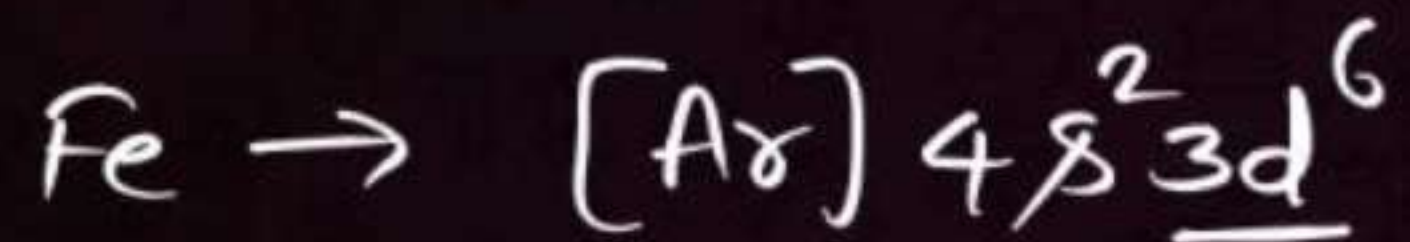
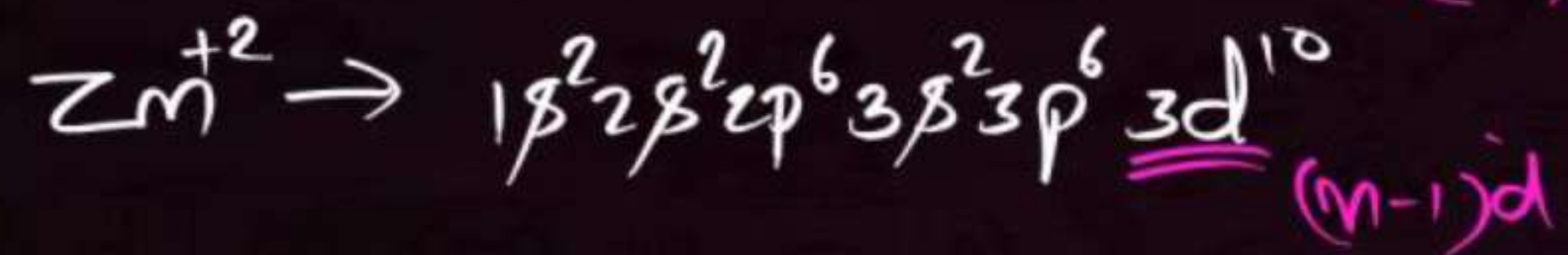
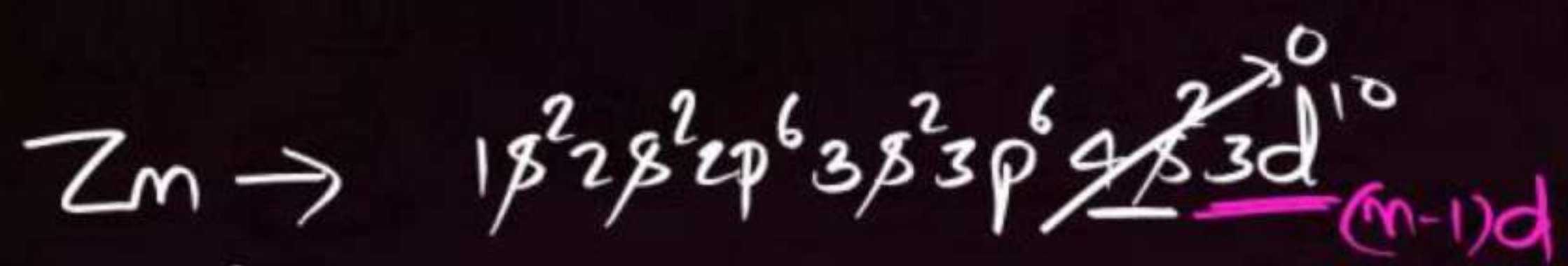
Q No. of transition elements in Long form periodic Table?

① 40

② 30

~~③ 36~~

④ 32



* Inner Transition Elements $\Rightarrow$ f-block Elements

* NEET 2025 *

General e⁻ config $\Rightarrow ns^2(n-1)d^{0-1}(n-2)f^{1-14}$



General Configuration Of a Period :

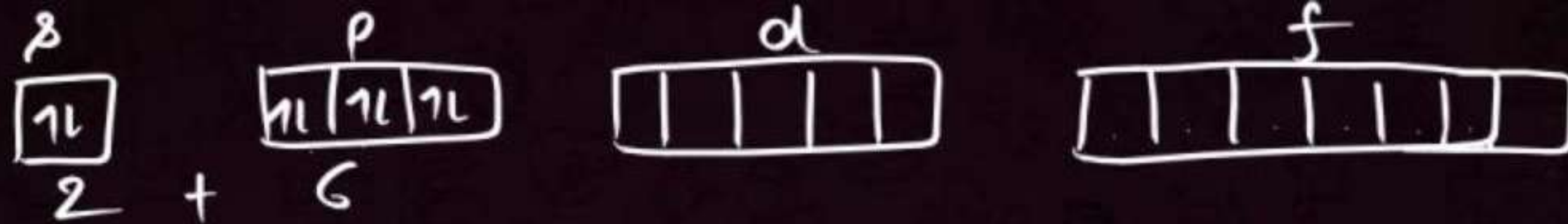


$n s (n-2) f (n-1) d n p$

$n=1 \Rightarrow \underline{1s} \quad \cancel{1p}$
 $n=2 \Rightarrow 2s \quad \cancel{1d} \quad 2p$
 $n=3 \Rightarrow \underline{3s} \quad \cancel{1f} \quad \cancel{2d} \quad 3p$
 $n=4 \Rightarrow 4s \quad \cancel{2f} \quad 3d \quad 4p$
 $n=5 \Rightarrow 5s \quad \cancel{3f} \quad 4d \quad 5p$
 $n=6 \Rightarrow 6s \quad 4f \quad 5d \quad 6p$

~~$\cancel{1s}$
 $n=1$
 $l=1$~~

~~$\cancel{2s}$~~
 $2p$
 $n=2$
 $l=1$
 $n > l$



PERIOD NO.	SUBSHELLS FILLED	ELEMENTS	NO. OF ELEMENTS	PERIOD NAME
1	$1s^2$		2	SHORTEST
2	$2s^2, 2p^6$		8	SHORT
3	$3s^2, 3p^6$		8	SHORT
4	$4s^2, 3d^{10}, 4p^6$		18	LONG
5	$5s, 4d, 5p$		18	LONG
6	$6s^2, 4f^{14}, 5d^{10}, 6p^6$		32	LONGEST
7	$7s, 5f, 6d, 7p$		32	LONGEST*

Q If an orbital can accommodate $3e^-$ then find the no. of elements in 3rd period?

$$n=3 \Rightarrow 3s^3 \quad 3p^9 \quad \underline{3+9=12}$$



$$4 \times 2 = 8$$

$$4 \times 3 = 12$$

$$4 \times 5 = \underline{\underline{20}}$$



$$\underline{\underline{n=8}}$$

$n s \dots (n-4) h \quad (n-3) g \quad (n-2) f \quad (n-1) d \quad n p$

$$\frac{8s}{1}$$

$$\cancel{3i} \quad \cancel{4h} \quad \cancel{5g} \quad \frac{6f}{7} \quad \frac{7d}{5} \quad \frac{8p}{3}$$

$$\cancel{n=4} \\ \cancel{l=5}$$

$$n=5 \\ l=4$$

$$\boxed{25 \times 2 = \underline{\underline{50}}}$$

$$\begin{aligned} &\underline{\underline{l}} \\ s &\rightarrow 0 \\ p &\rightarrow 1 \\ d &\rightarrow 2 \\ f &\rightarrow 3 \\ \boxed{g} &\rightarrow 4 \\ h &\rightarrow 5 \\ i &\rightarrow 6 \\ j &\rightarrow 7 \\ k &\rightarrow 8 \\ & \\ & \end{aligned}$$

Q If an orbital can accommodate $3e^-$ then no of elements present in 8th period?

$$\begin{array}{rcl} 25 \times 2 & = & 50 \\ & \times 3 & = 75 \\ 25 \times 5 & = & 125 \end{array}$$



QUESTION



The electronic configuration of an element is $1s^2, 2s^2, 2p^6, 3s^2, 3p^6, 3d^{10}, 4s^2, 4p^6, 4d^{10}, 5s^2, 5p^3$. From which group of the periodic table it belongs?

(Karnataka CET 2010)

A 2^{nd}

B 5^{th}

C 3^{rd}

D 7^{th}

Question



Find the no. of elements present in 7th period of the periodic table ?

A 18

C 50

~~**B** 32~~

D 72

Question



Find the no. of elements present in 3rd period of the periodic table ?

A 18

C 2

B 32

~~**D** 8~~

Question



Find the no. of elements present in 8th period of the periodic table ?

A 18

B 32

~~**C** 50~~

D 72

Question



Find the no. of elements present in 8th period of the periodic table if an orbital can accommodate 3 e- ?

A 18

C 50

B 32

~~**D** 75~~

Question

How?



Find the no. of elements present in 9th period of the periodic table if an orbital can accommodate 3 e- ?

A

32

B

50

C

75

D

72

Problem 3.2

How would you justify the presence of 18 elements in the 5th period of the Periodic Table?

NCERT

18

Problem 3.3

The elements $Z = 117$ and 120 have not yet been discovered. In which family / group would you place these elements and also give the electronic configuration in each case.



IUPAC Naming of Elements



106
Uuh

Uuu

Uuu

Uue
i i g
Uuo
i i p
Uuo

104

Un + nil + quad + "ium"



Uus
~~107~~

Un mil quadium

(Uug)

Table 3.4 Notation for IUPAC Nomenclature of Elements

Digit	Name	Abbreviation
0	<u>ni</u> l	n
1	<u>u</u> n	u
2	<u>bi</u>	b
3	<u>tri</u>	t
4	quad	q
5	pent	p
6	hex	h
7	<u>sept</u>	s
8	<u>oct</u>	o
9	<u>enn</u>	e

Question

Unbinilium



What would be the IUPAC name of the element with atomic number 120?

[Kerala PMT-2015]

- A** Unnilbium
- B** Ununbium
- C** Unnilunium
- D** Unbinilium

Question



The IUPAC name of the element with atomic number 119 is : [NEET - 2022]

Ununennium

A

Unnilbium

B

Ununennium

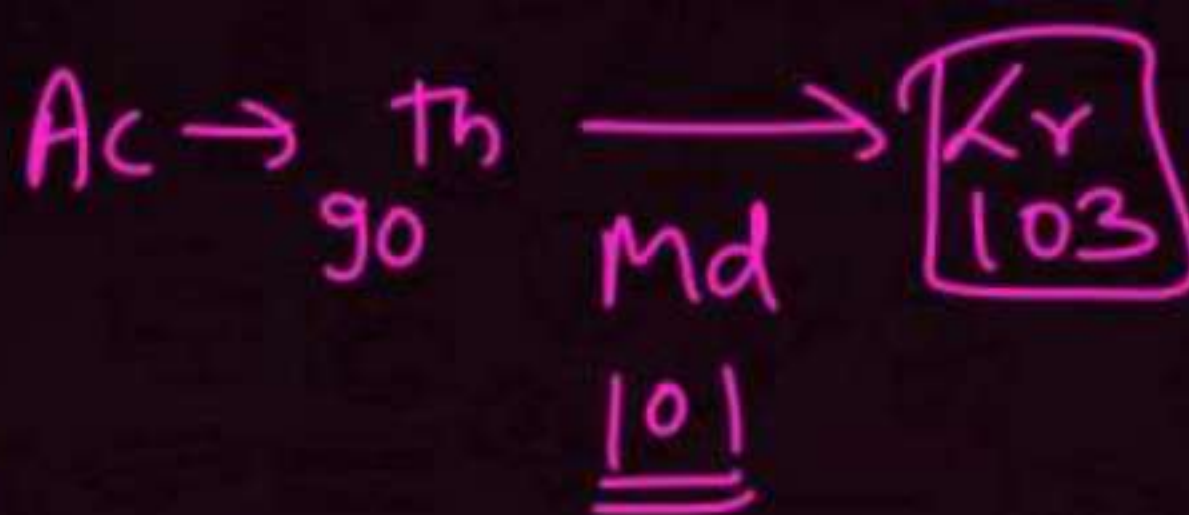
C

Unnilunium

D

Unbinilium

Question



Identify the incorrect match

[NEET - 2020]

NAME	IUPAC NAME
(a) Unnilunium ✓	(i) Mendelevium (101)
(b) Unniltrium ✓	(ii) Lawrencium (103)
(c) Unnilhexium ✓	(iii) Seaborgium (106)
(d) Unununium ✗	(iv) Darmstadtium (110)



A (b) – (ii)



B (c) – (iii)



C (d) – (iv)



D (a) – (i)

Group & Period Identification



① When Atomic No. is given :

$$\text{Group No.} = 18 + \text{given At. No.} - \text{At. No. of Next Noble gas}$$

$$\text{Period No.} = \text{Period No. of Next Noble gas.}$$

Ex : $Z = 34$

$$G = 18 + 34 - 36 = 16$$

$$G = ?$$

$$G = 16$$

$$P = ?$$

$$P = 4$$

$2\text{He} \rightarrow 1$
 $10\text{Ne} \rightarrow 2$
 $18\text{Ar} \rightarrow 3$
 $36\text{Kr} \rightarrow 4$
 $54\text{Xe} \rightarrow 5$
 $86\text{Rn} \rightarrow 6$
 $118\text{Og} \rightarrow 7$

$$\underline{\underline{118 \geq Z \geq 104}}$$

$$P = 7^{\text{th}}$$

G = last 2 digits of At. No.



Ex: $\rightarrow$

116

$$\Rightarrow G = 16$$

$$P = 7$$

$$Z = \underline{\underline{119}}$$

$$G = 1 \quad ?$$
$$P = 8$$

106

$$\Rightarrow G = 6$$

$$P = 7$$

117

$$\Rightarrow$$

$$G = 17$$

$$P = 7$$

$$Z = 120$$

$$G = 2$$

$$P = 8$$

Case-2 $\therefore \rightarrow$ if e-config is given.

Period no. = value of n



$$\underline{P = 6}$$

$$G =$$





Screening Effect & Zeff



Zeff → A.A.P.O.N

$$\text{Nuclear charge} = Z(\text{no. of } p^+)$$

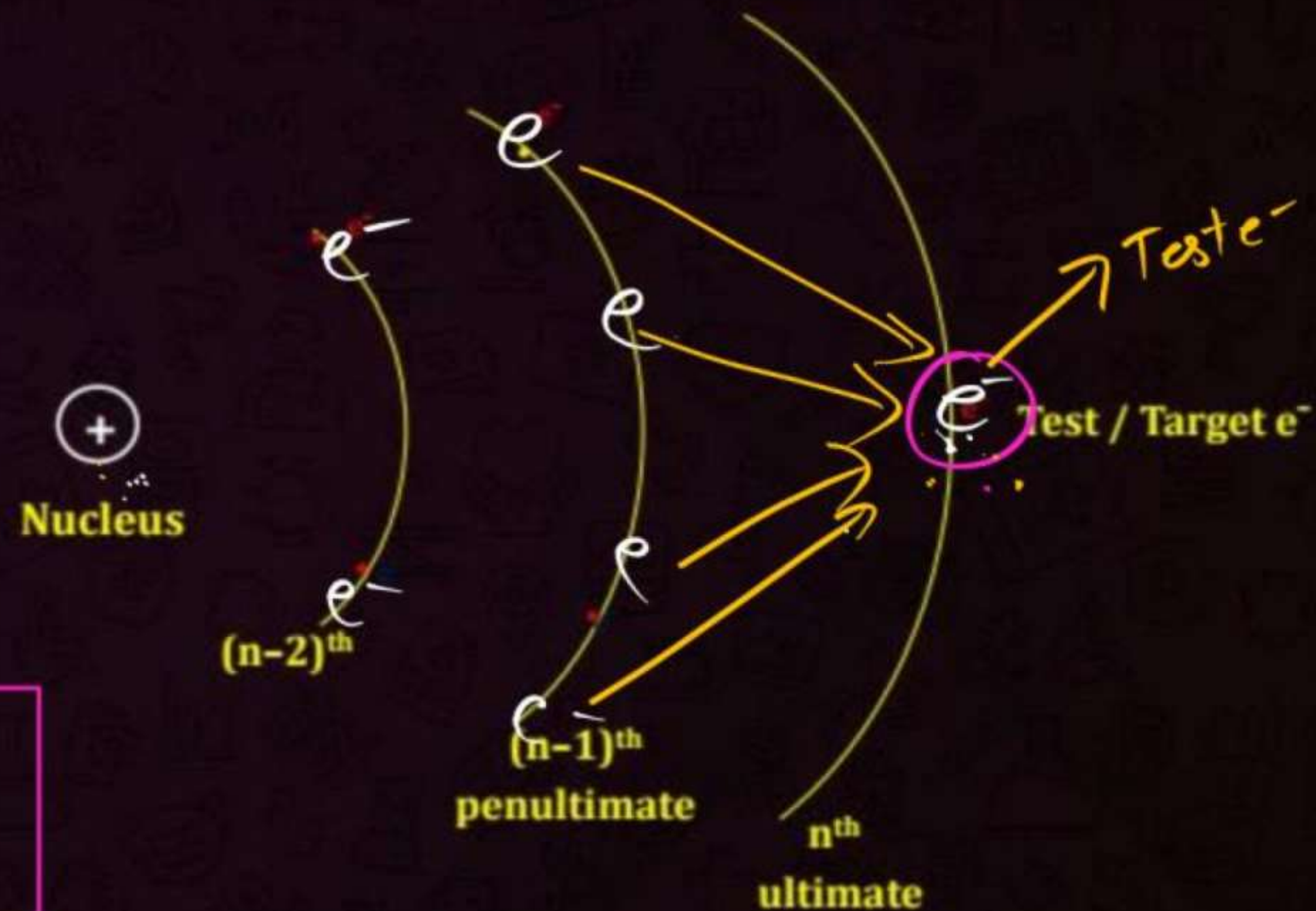


$$\text{Actual } Att_{\gamma}^m = Att_{\gamma}^m - R_{ep}$$

$$Z_{eff} = Z - \sigma$$

σ = screening / shielding const.

order of screening = $s > p > d > f$



* Z_{eff} $\rightarrow$ in a period $Z_{eff} \uparrow$ from L to R

* Isoelectronic species $\rightarrow$

EX: F^{-} O^{2-} N^{3-} $-ve \uparrow Z_{eff} \downarrow$

Na^{+} Mg^{2+} Al^{3+} $+ve \uparrow Z_{eff} \uparrow$

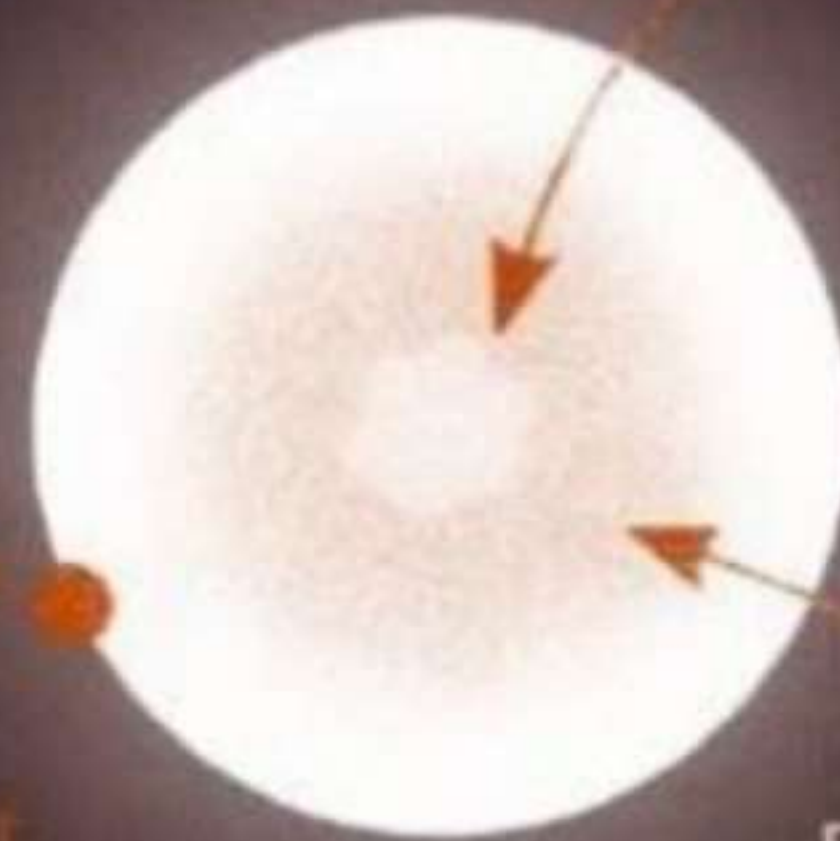
$$Z_{eff} \propto \frac{+ve}{-ve}$$

Electrons outside
have no effect for
electron of interest

Positively charged
nucleus

Electron of interest

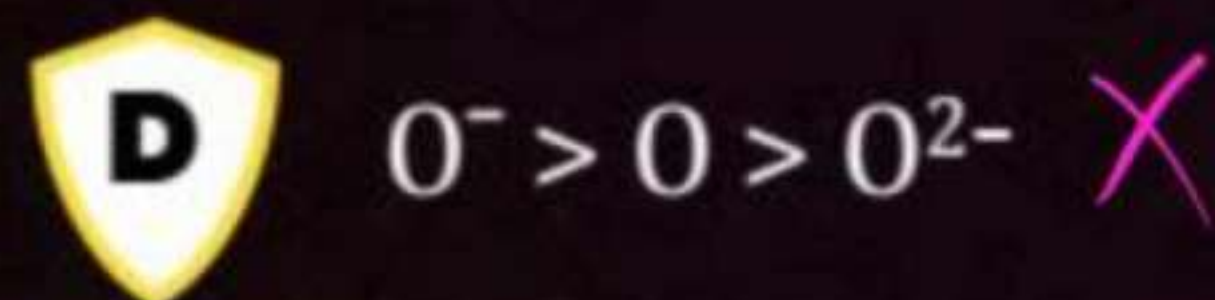
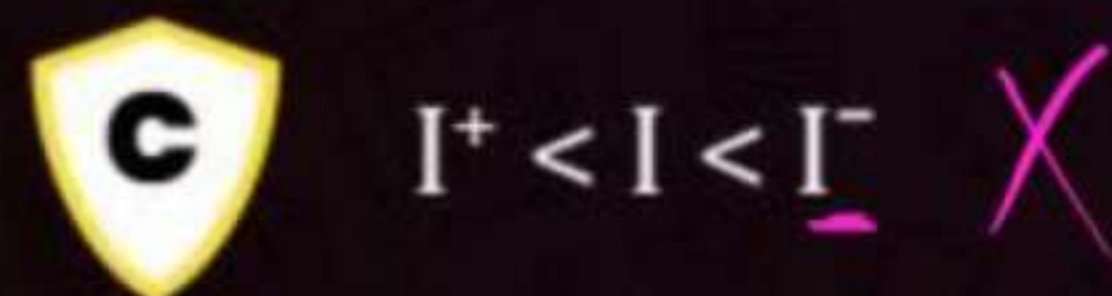
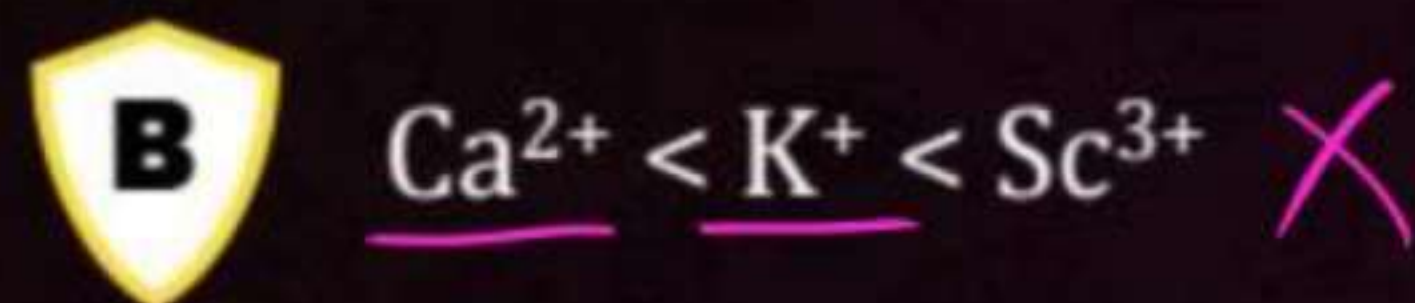
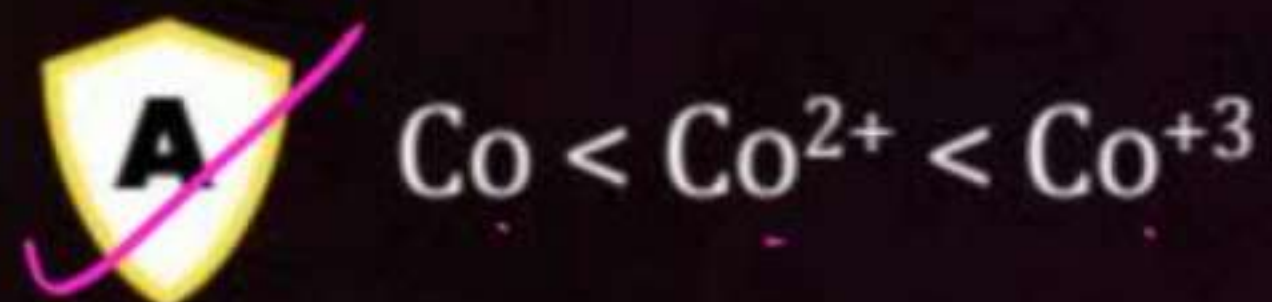
Electrons in between
cancel some of
the nuclear charge



Question



Determine the correct order of Z_{eff} ?



Question



Screening effect is not observed in

A He^+

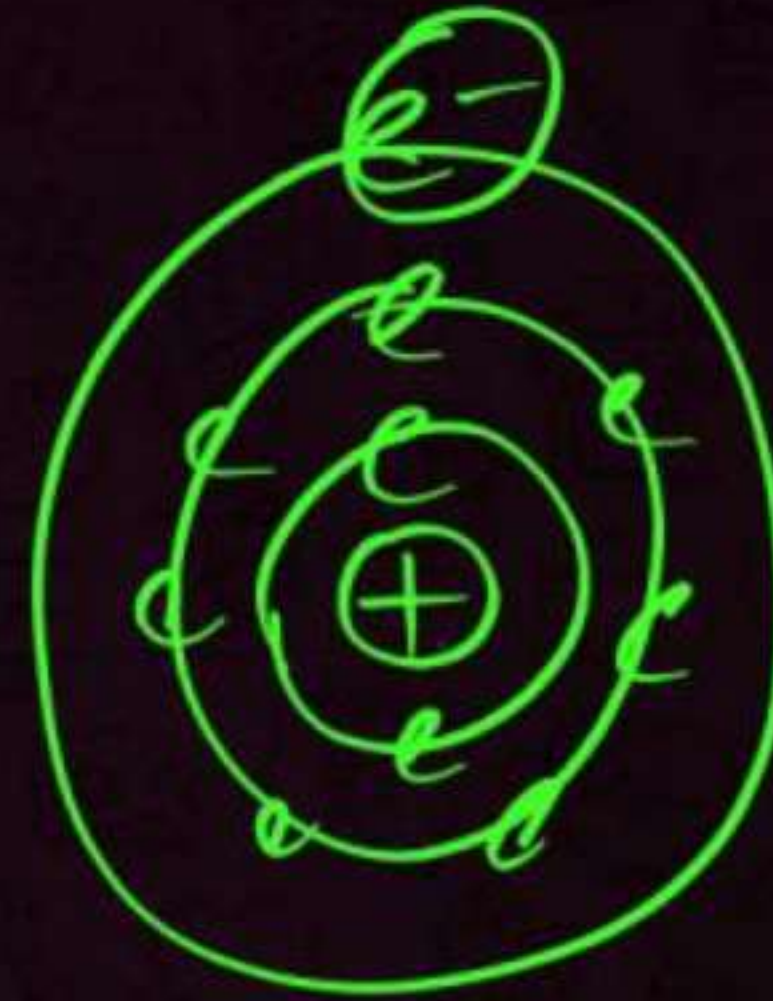
B Li^{+2}

C H

D All of these



$$\sigma = 0$$





ATOMIC RADIUS



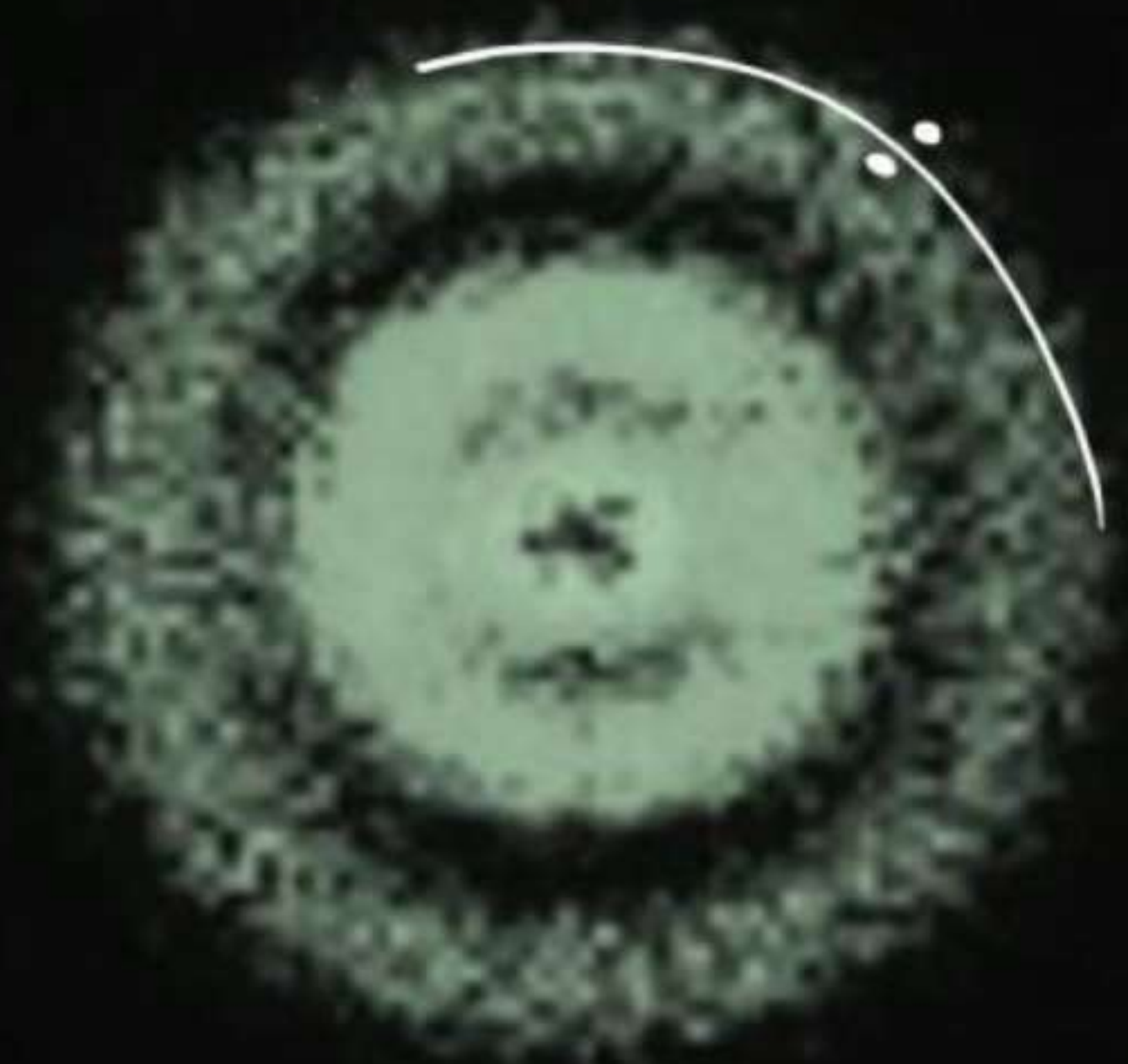
It is the distance b/w centre of the nucleus and the outermost e^- in an isolated gaseous atom

Difficulties :

(ओके ला / single)

- ① size of atom is very small
- ② boundary of an atom is not defined
- ③ Isolation of an atom is very difficult.

YOU CAN CALL ME



“ATOM”

Types Of Atomic Radii

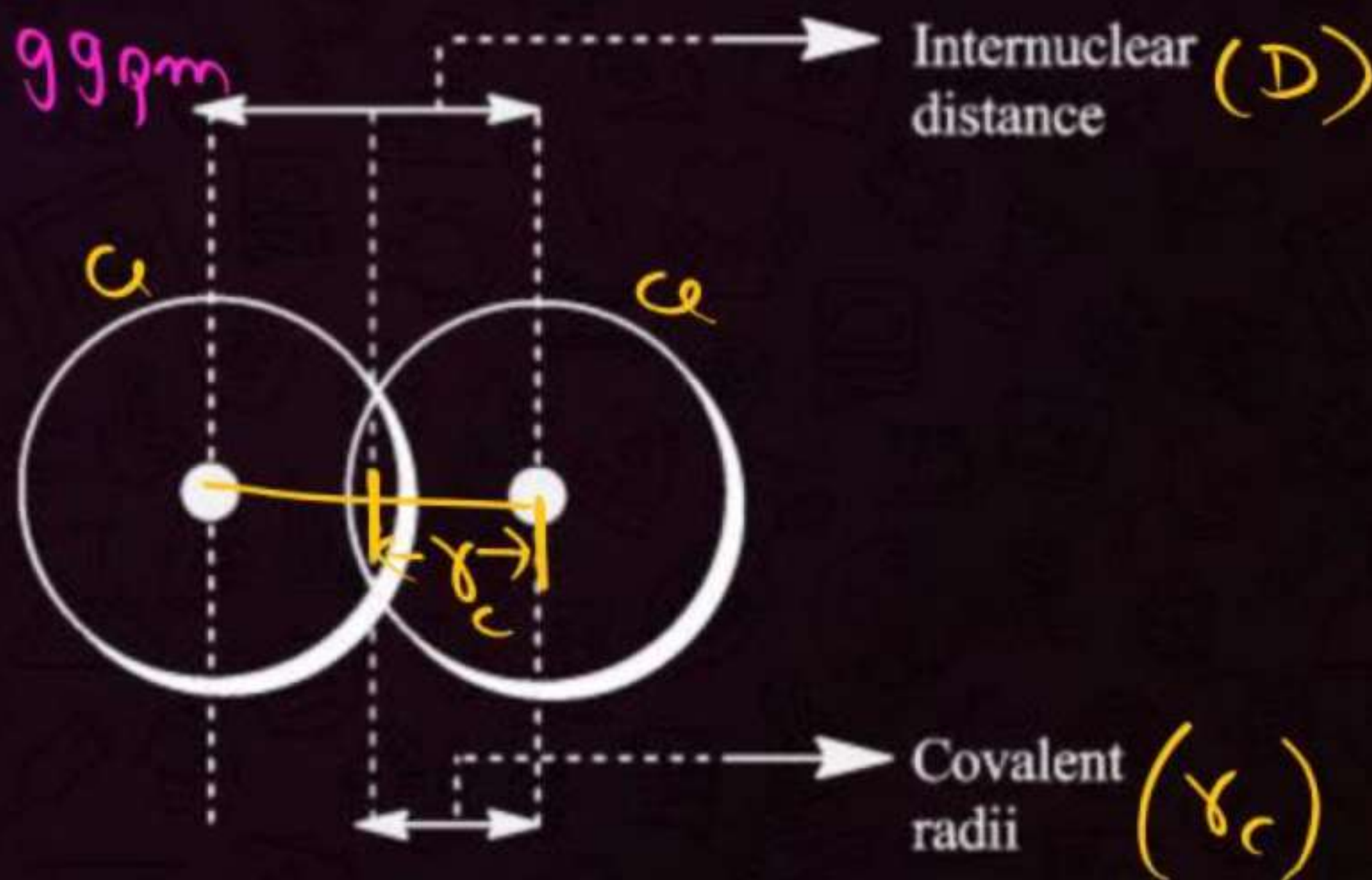
(i) Covalent Radius :

Cl_2 $D_{\text{Cl-Cl}} = 198 \text{ pm}$
 $r_c = \frac{198}{2} = 99 \text{ pm}$

$D \rightarrow$ Internuclear distance / B.L.

$$r_c = \frac{D}{2}$$

$$= \frac{\text{B.L.}}{2}$$

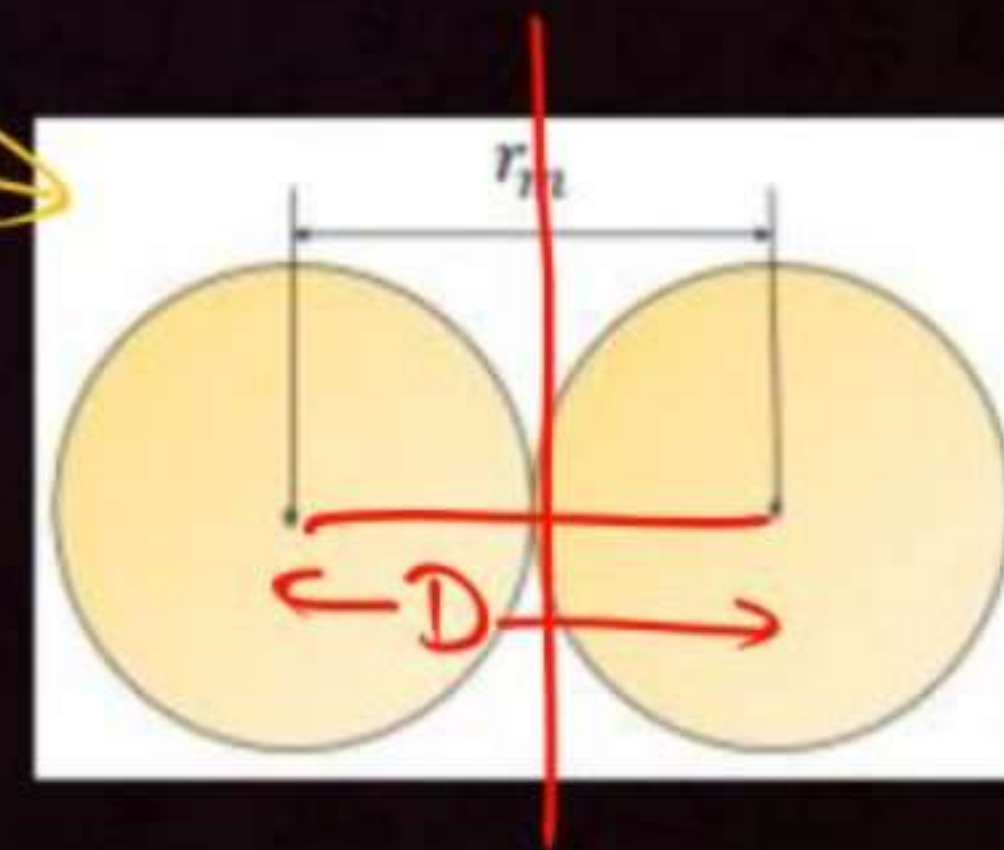
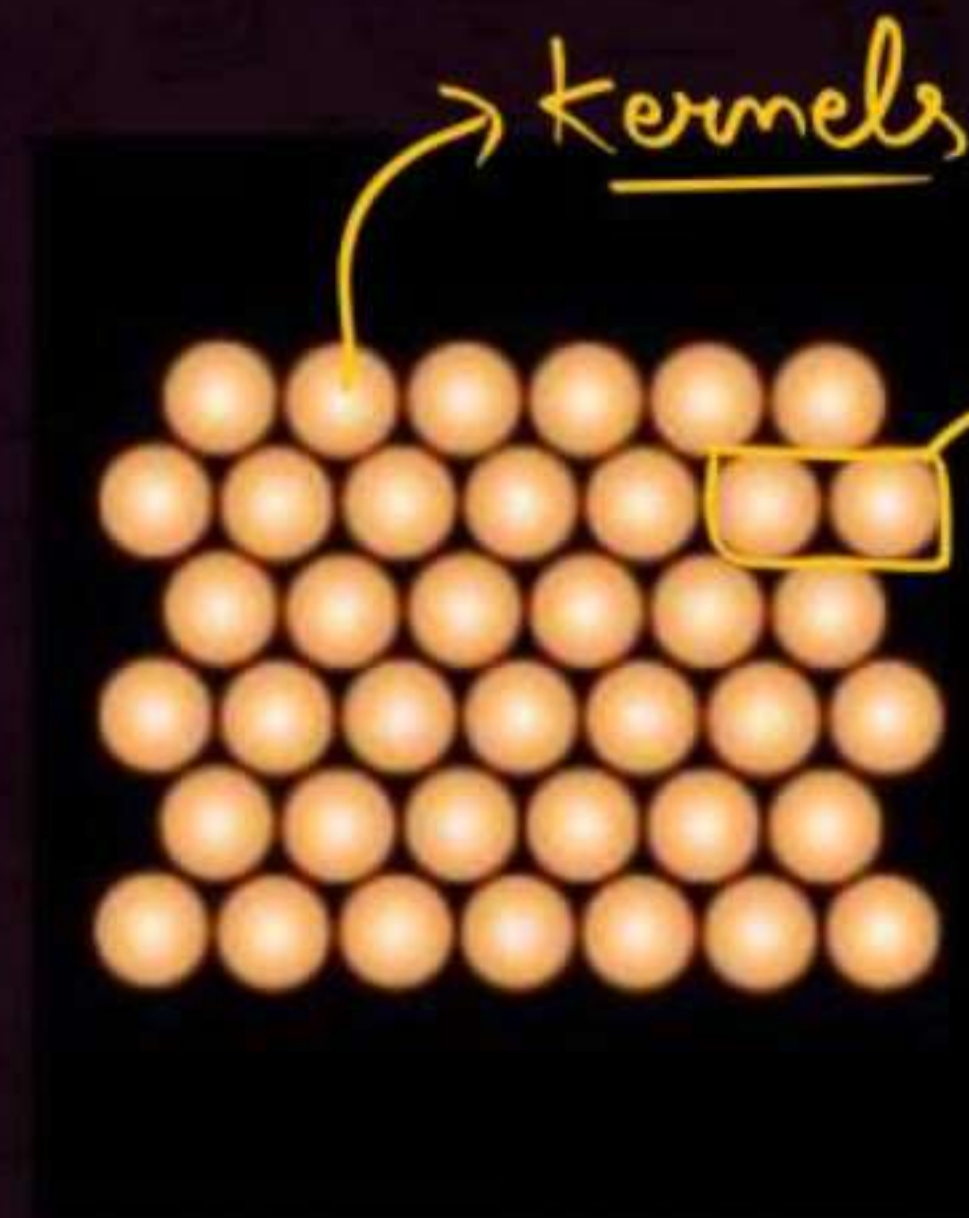


(ii) Metallic Radius : (Crystal Radius)

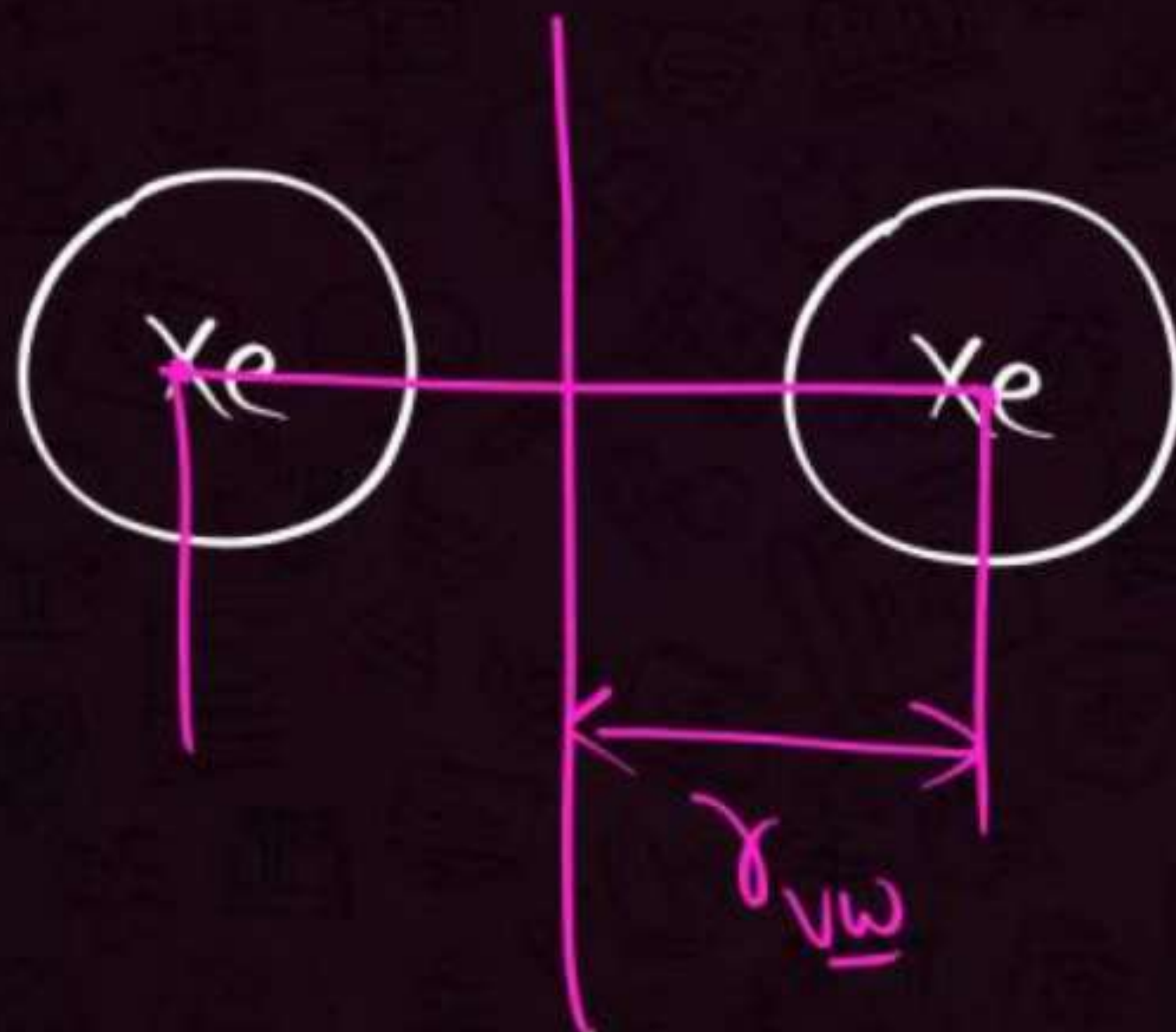
$$r_m = \frac{D}{2}$$

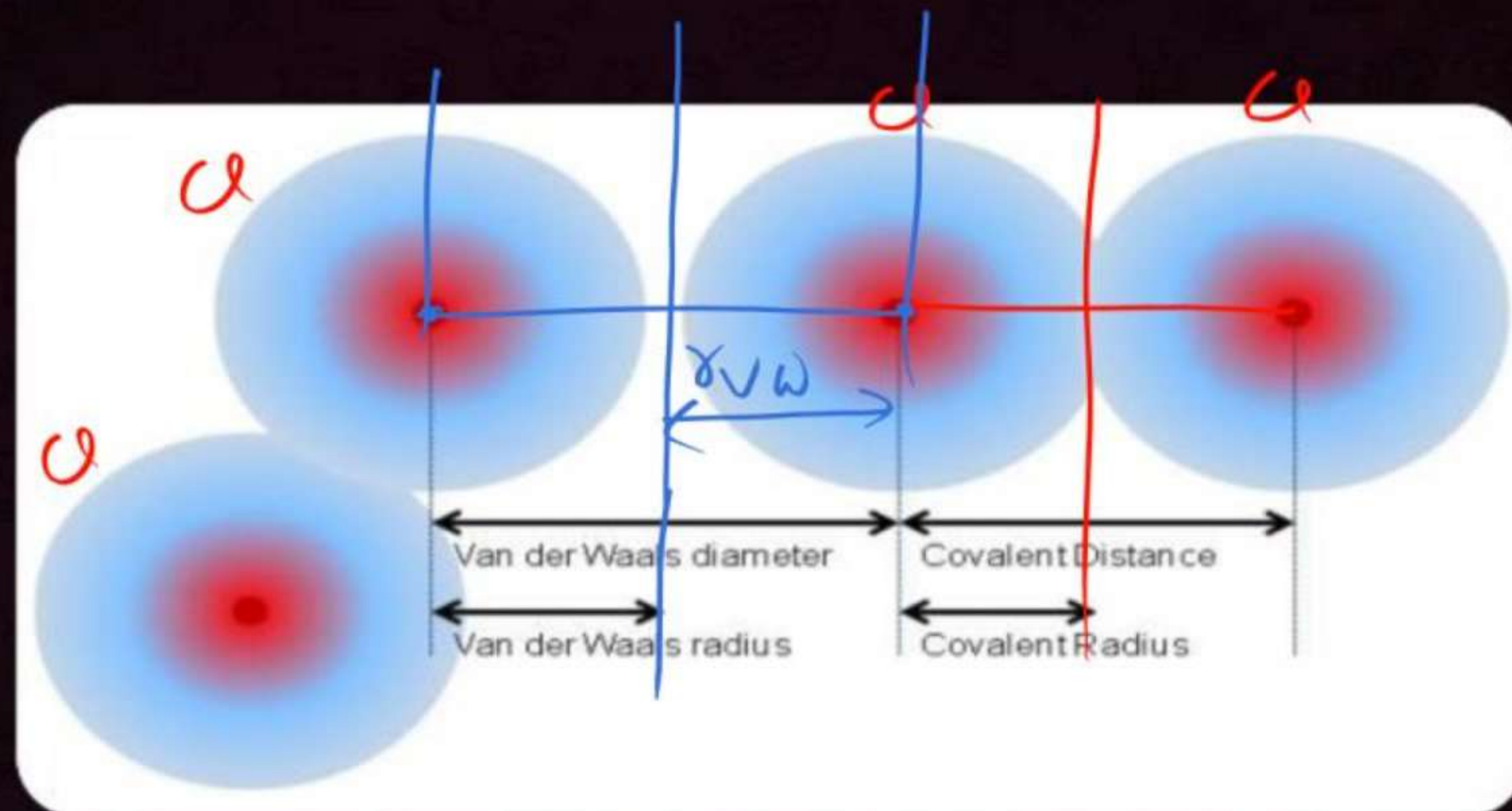
$$r_m \propto \frac{1}{\text{Met. Bond Strength}}$$

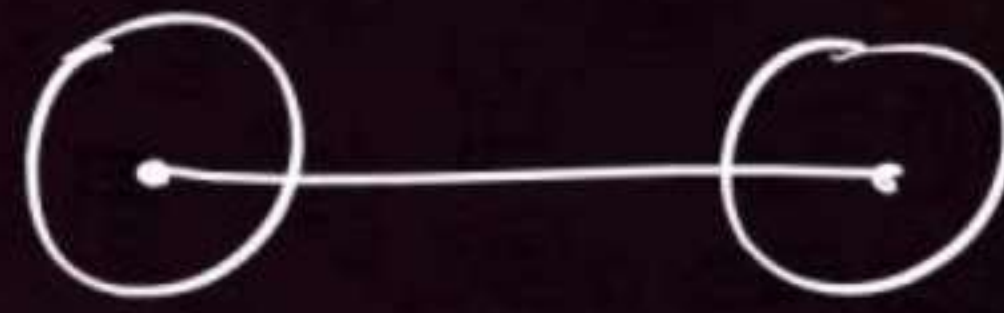
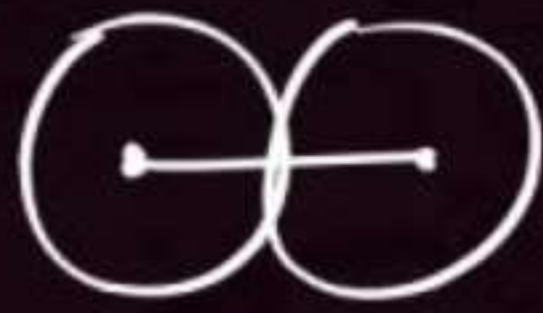
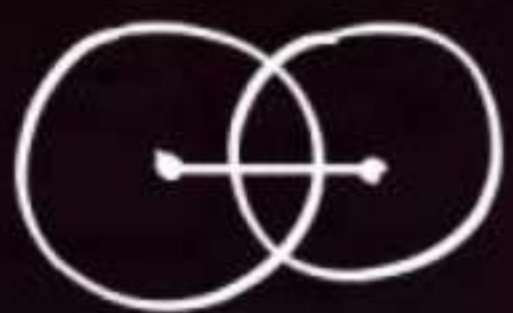
$C_r < M_m$
 ↓
Weak Met. Bond



(iii) Vander Waal Radius : (Noble gas Radius)



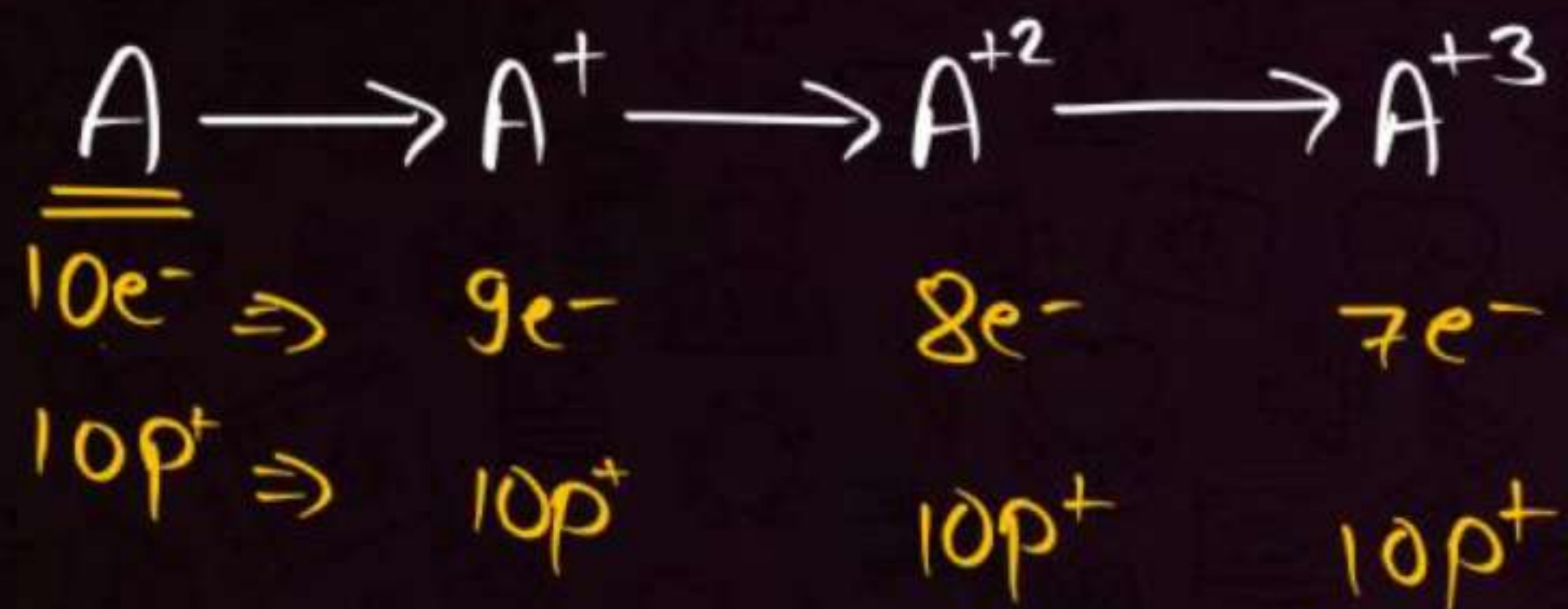


 χ_c $<$ χ_m $<$ χ_{vw}

IONIC RADII



Cationic Radius



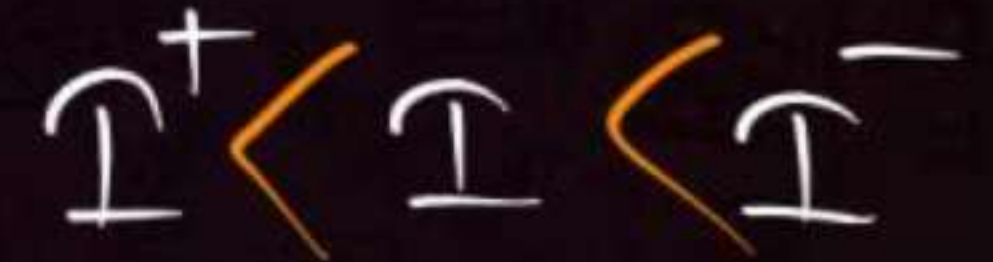
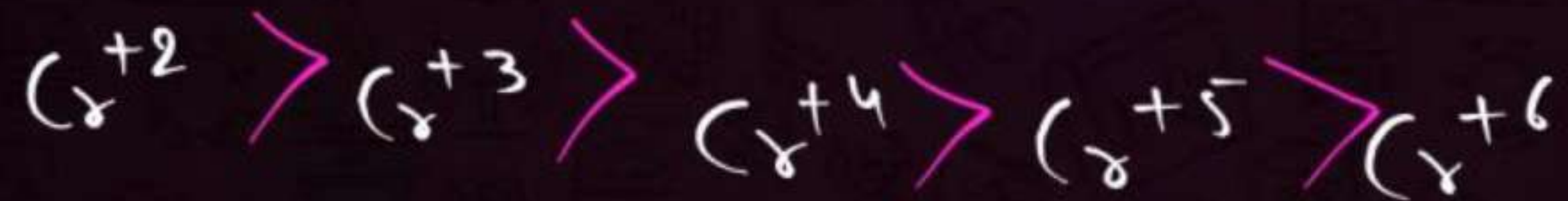
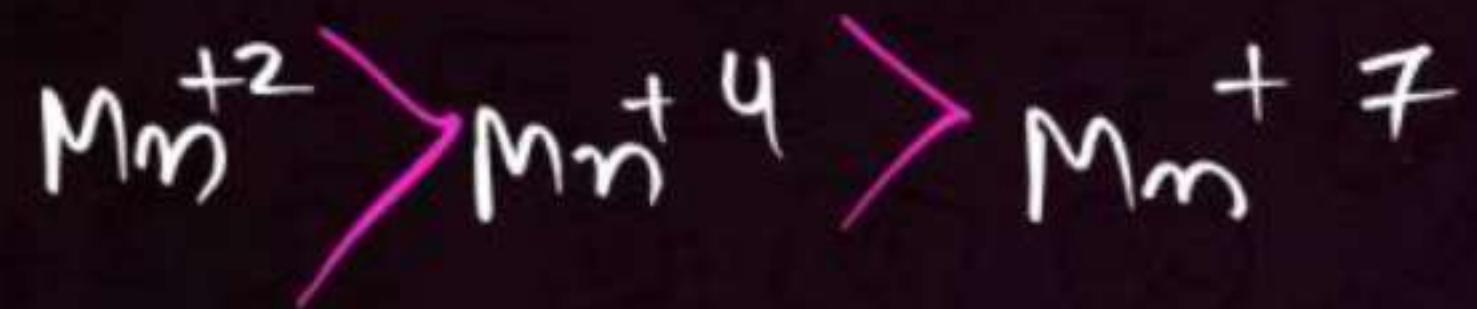
size $\propto \frac{1}{+ve}$

Anionic Radius

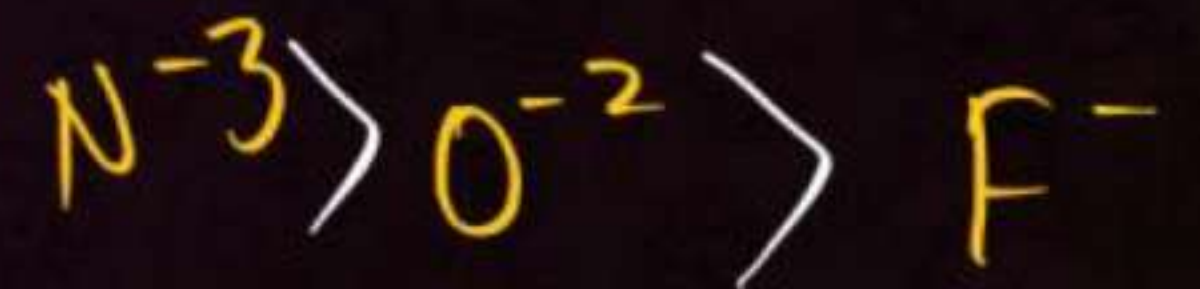
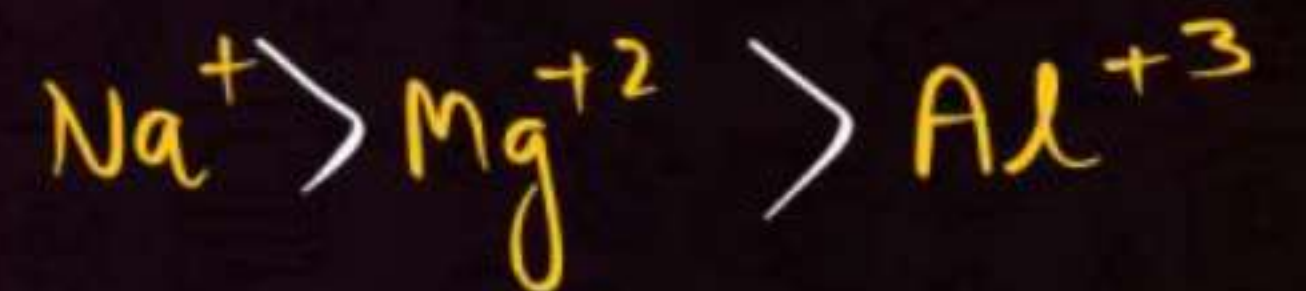


size $\propto -ve$

size $\propto$ $\frac{-ve}{+ve}$



Cation < Neutral < Anion



Variation Of Size in Group & Period



Period $\Rightarrow$ $L \xrightarrow[\text{size} \downarrow]{z_{\text{eff}} \uparrow} R$

$$\text{size} \propto \frac{1}{z_{\text{eff}}}$$

$\text{Li} > \text{Be} > \text{B} > \text{C} > \text{N} > \text{O} > \text{F} < \text{Ne}$
V.W.R.

Group $\Rightarrow$ $\begin{matrix} \text{Top} \\ \downarrow \\ \text{Bottom} \end{matrix}$ $\begin{matrix} \text{no. of shells} \uparrow \\ \text{size} \uparrow \end{matrix}$

$\text{Li} < \text{Na} < \text{K} < \text{Rb} < \text{Cs}$

$\text{Be} < \text{Mg} < \text{Ca} < \text{Sr} < \text{Ba}$

Table 3.6(a) Atomic Radii/pm Across the Periods

Atom (Period II)	Li	Be	B	C	N	O	F
Atomic radius	152	111	88	77	74	66	64
Atom (Period III)	Na	Mg	Al	Si	P	S	Cl
Atomic radius	186	160	143	117	110	104	99

Table 3.6(b) Atomic Radii/pm Down a Family

Atom (Group I)	Atomic Radius	Atom (Group 17)	Atomic Radius
Li	152	F	64
Na	186	Cl	99
K	231	Br	114
Rb	244	I	133
Cs	262	At	140

QUESTION



Increasing order or ionic size for the ions, F^{-} , O^{2-} , Na^{+} , Al^{3+} is:

(AMU 2009)

- ☐ A $O^{2-} < F^{-} < Na^{+} < Al^{3+}$
- ☒ B $Al^{3+} < Na^{+} < F^{-} < O^{2-}$
- ☐ C $O^{2-} < Na < F^{-} < Al^{3+}$
- ☐ D $Al^{3+} > F^{-} < Na^{+} < O^{2-}$

Ans. (D)

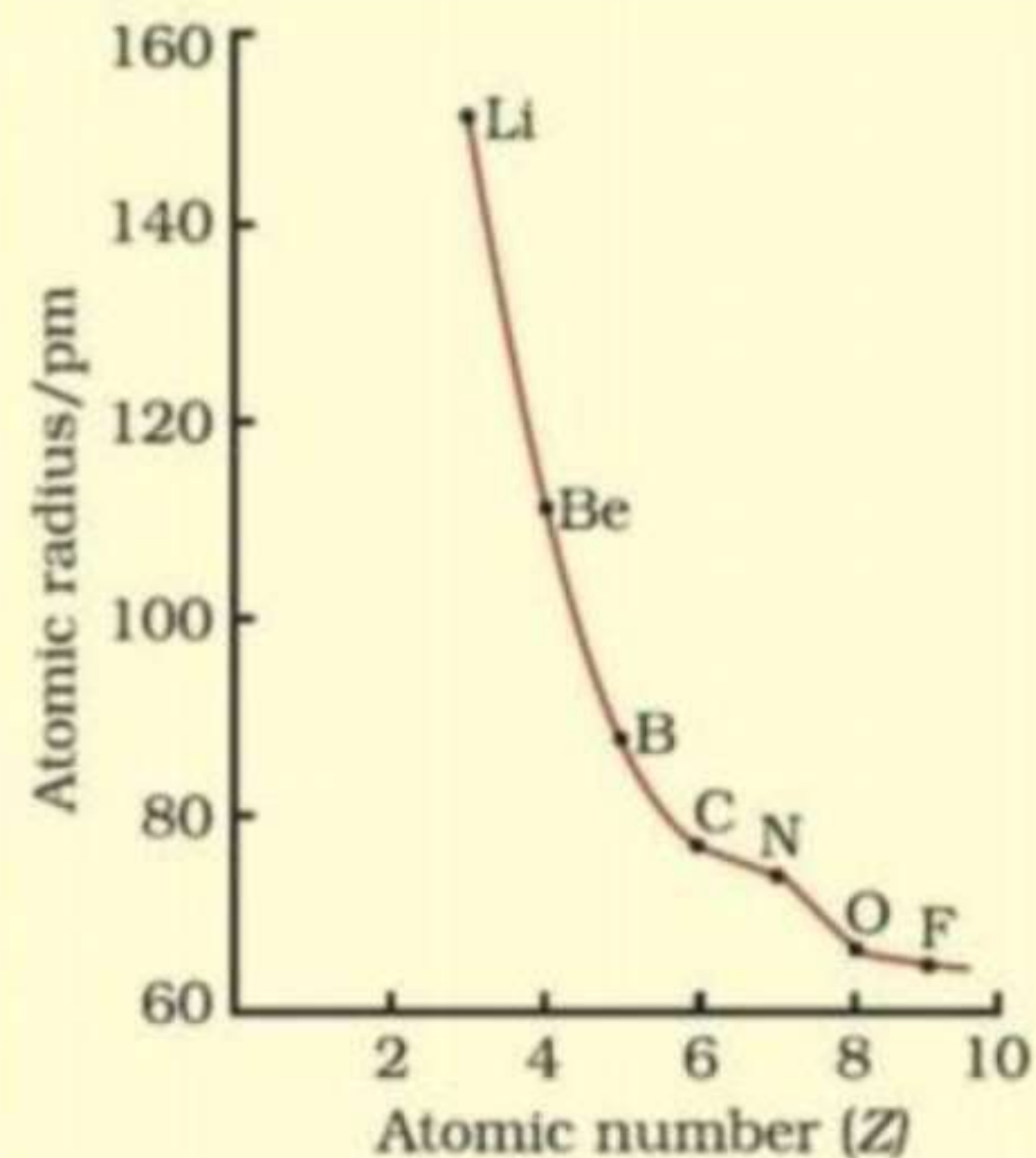


Fig. 3.4 (a) Variation of atomic radius with atomic number across the second period

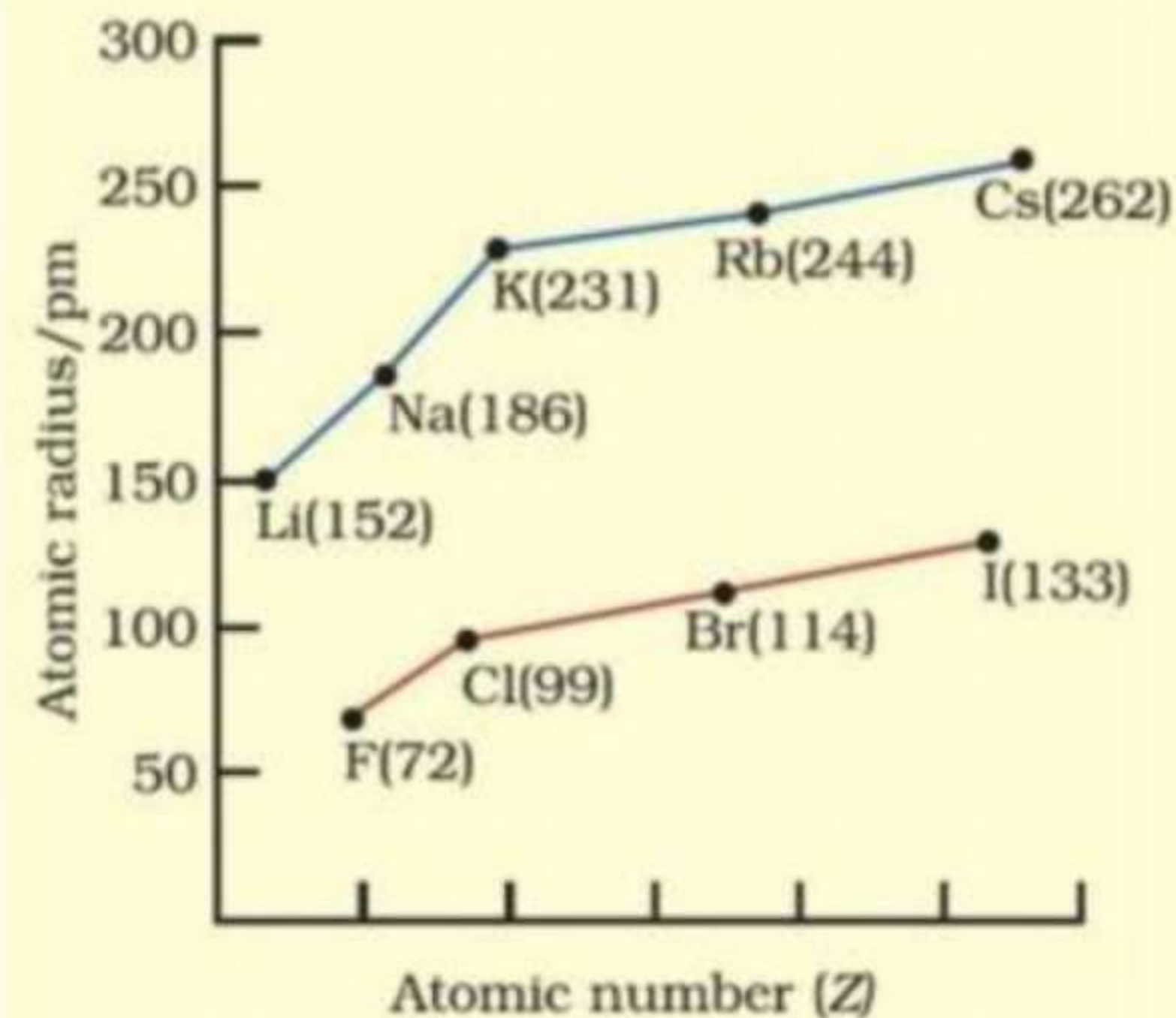


Fig. 3.4 (b) Variation of atomic radius with atomic number for alkali metals and halogens

NCERT CORNER

In general, the ionic radii of elements exhibit the same trend as the atomic radii. A cation is smaller than its parent atom because it has fewer electrons while its nuclear charge remains the same. The size of an anion will be larger than that of the parent atom because the addition of one or more electrons would result in increased repulsion among the electrons and a decrease in effective nuclear charge. For example, the ionic radius of fluoride ion (F^-) is 136 pm whereas the atomic radius of fluorine is only 64 pm. On the other hand, the atomic radius of sodium is 186 pm compared to the ionic radius of 95 pm for Na^+ .

Problem 3.5

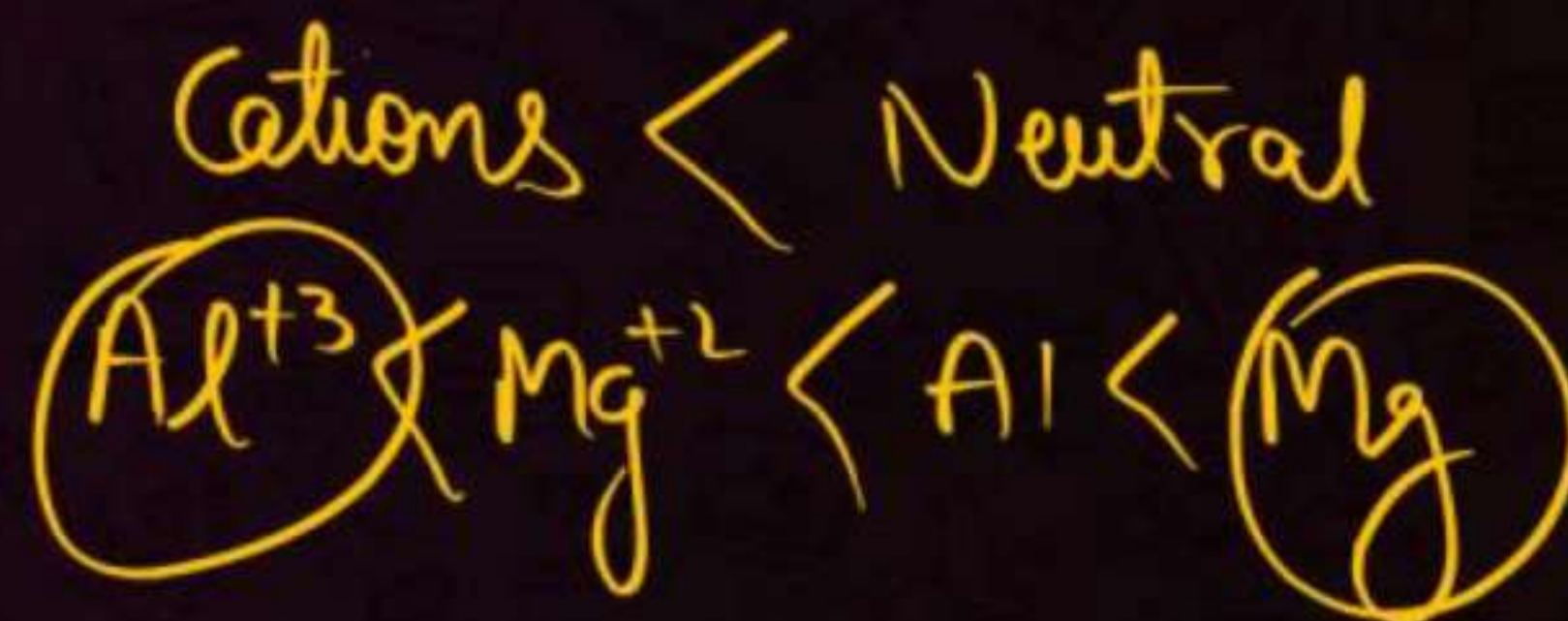
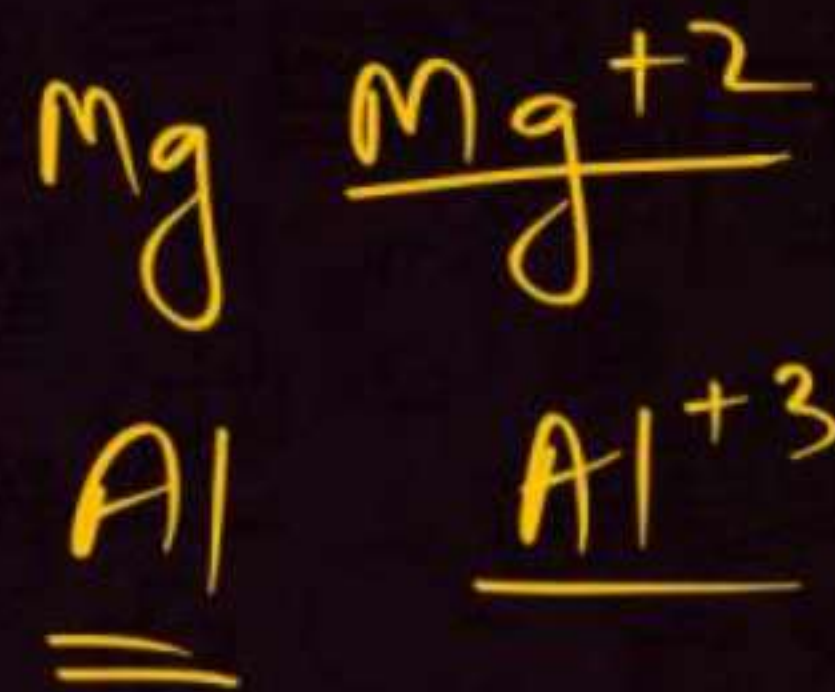
Which of the following species will have the largest and the smallest size?

Mg, Mg^{2+} , Al, Al^{3+} .

Solution

Atomic radii decrease across a period. Cations are smaller than their parent atoms. Among isoelectronic species, the one with the larger positive nuclear charge will have a smaller radius.

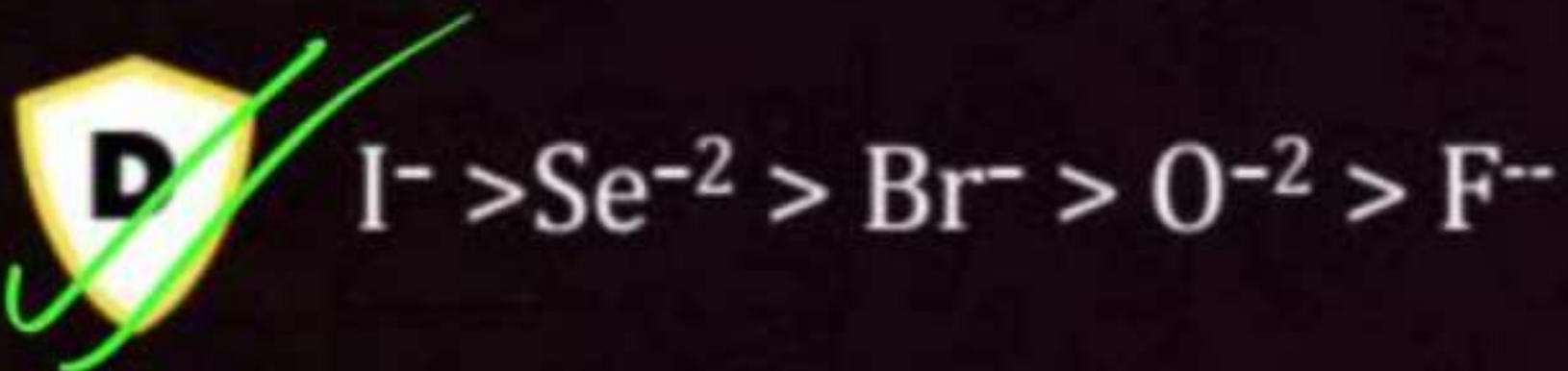
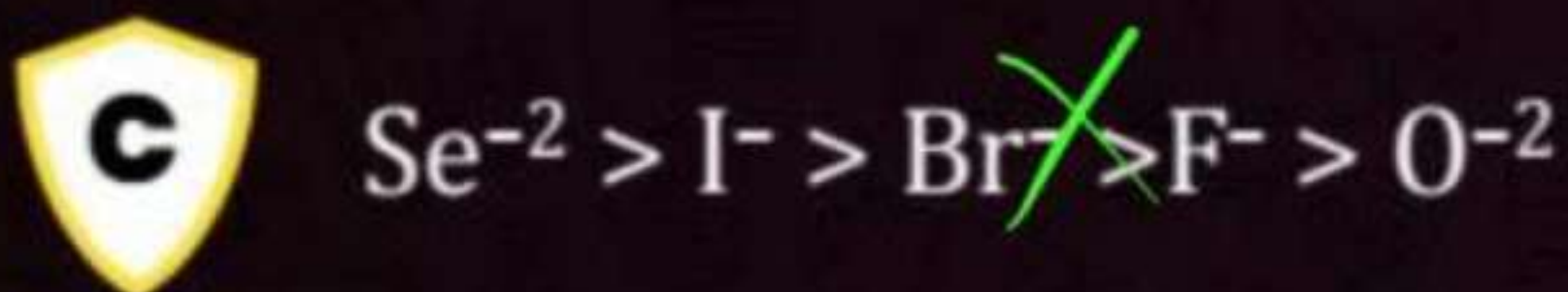
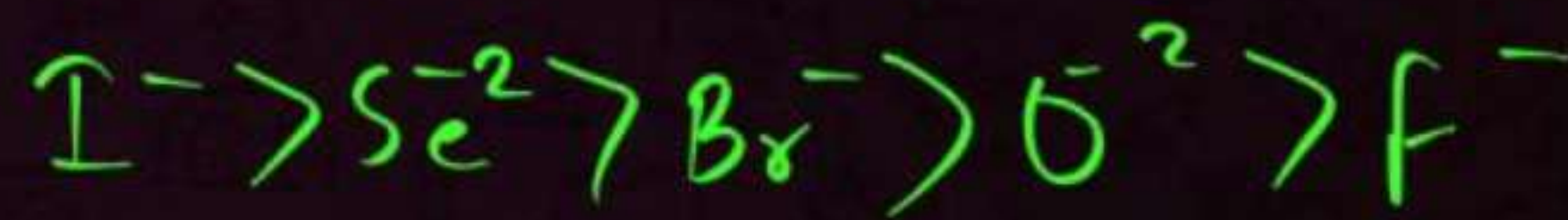
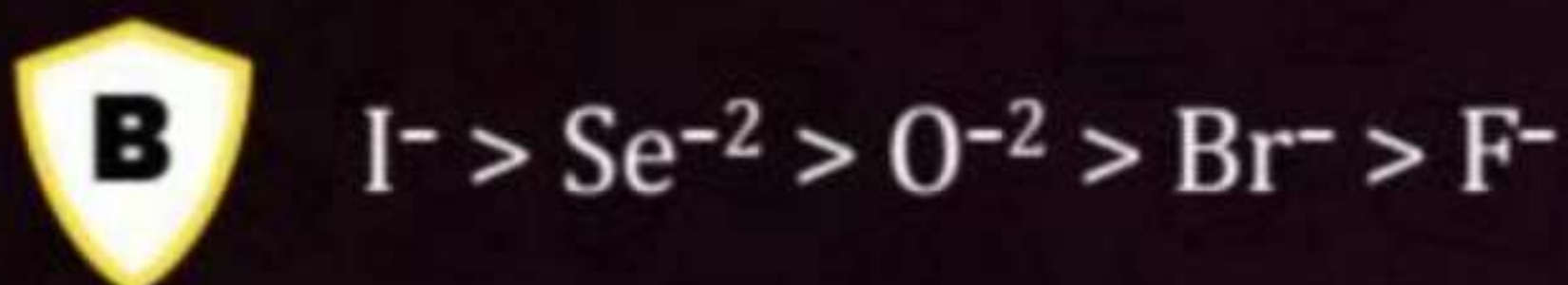
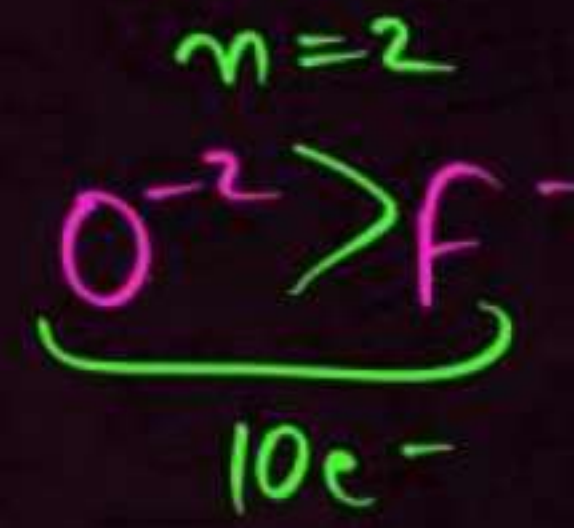
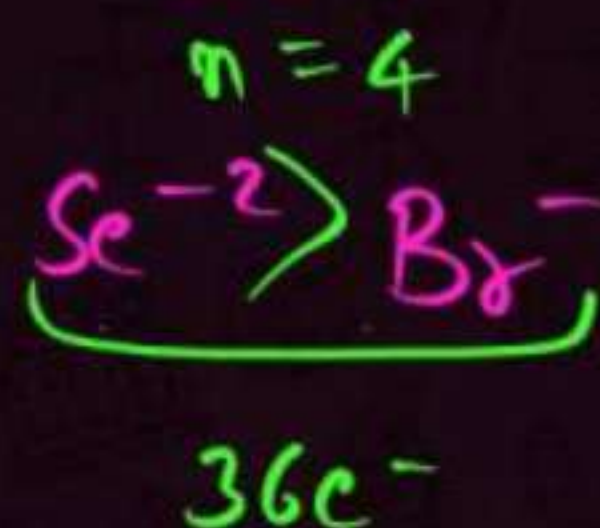
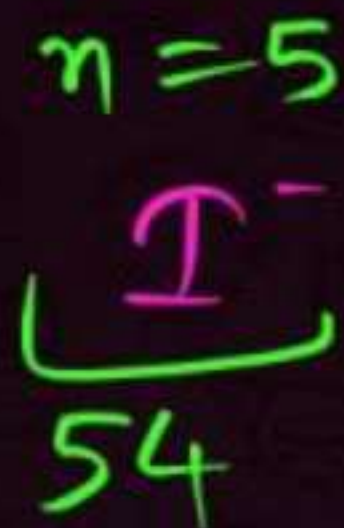
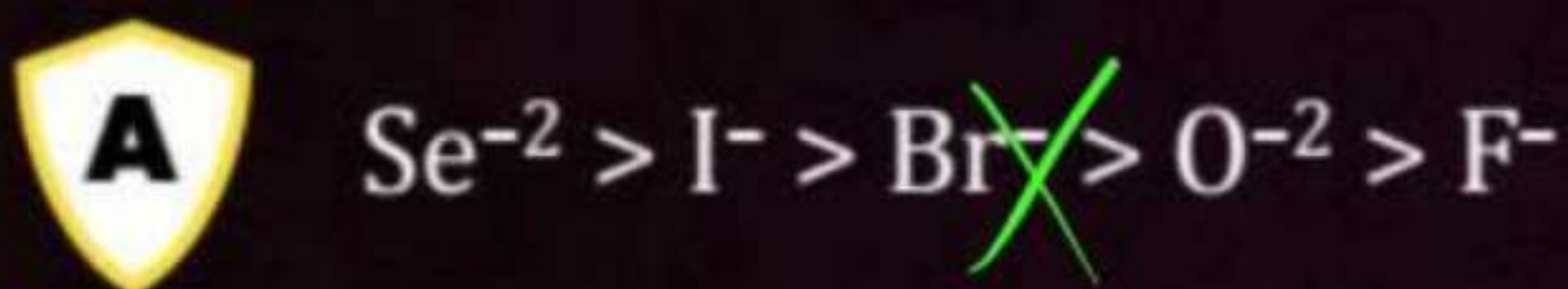
Hence the largest species is Mg; the smallest one is Al^{3+} .



Question



Which of the following sequences is correct for decreasing order of ionic radius –



Order of size in P-block :



$Al > Ga$
S.C/T.C.

3de- $\Rightarrow$ poor $\sigma \Rightarrow$ zeff $\uparrow$





$\text{He} < \text{Ne} < \text{Ar} < \text{Kr} < \text{Xe} < \text{Rn} < \text{Og}$

Order of size in D-block :



3d series

h h h h h



z_{eff} dominating

$z_{\text{eff}} \approx 0$

σ dominating

$$\underline{sc < y < da}$$



$$Ti < Zr \approx Hf$$

L.C.

$$4f^{14}$$

very poor σ

$$Z_{eff} \uparrow$$

size $\downarrow$

$$3d < 4d \approx 5d$$

Ionic Radius (f-block) :

$z_{eff} \uparrow$ size $\downarrow$

$\text{La}^{+3} \rangle \text{Ce}^{+3} \rangle \text{Pr}^{+3} \rangle \text{Nd}^{+3} \rangle \text{Pm}^{+3} \rangle \text{Sm}^{+3} \rangle \text{Eu}^{+3} \rangle \text{Gd}^{+3} \rangle \text{Tb}^{+3} \rangle \text{Dy}^{+3} \rangle \text{Ho}^{+3} \rangle \text{Er}^{+3} \rangle \text{Tm}^{+3} \rangle \text{Yb}^{+3} \rangle$

Lu^{+3}

$4f e^- \uparrow$

$\sigma \downarrow$

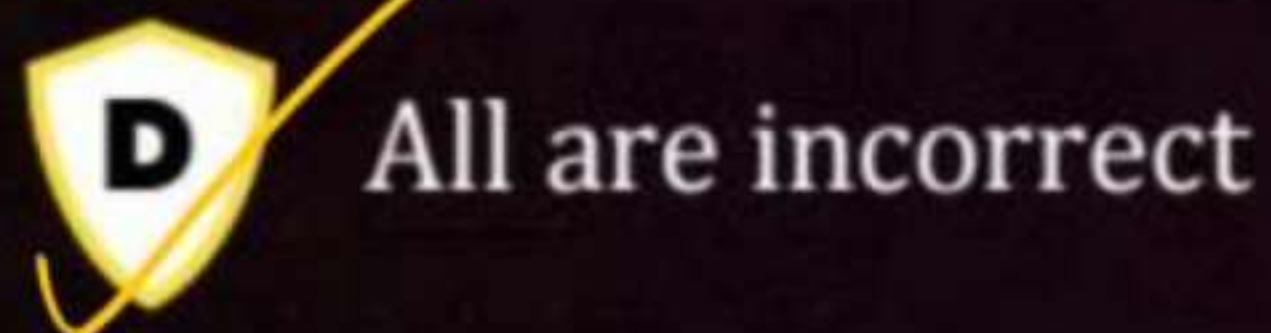
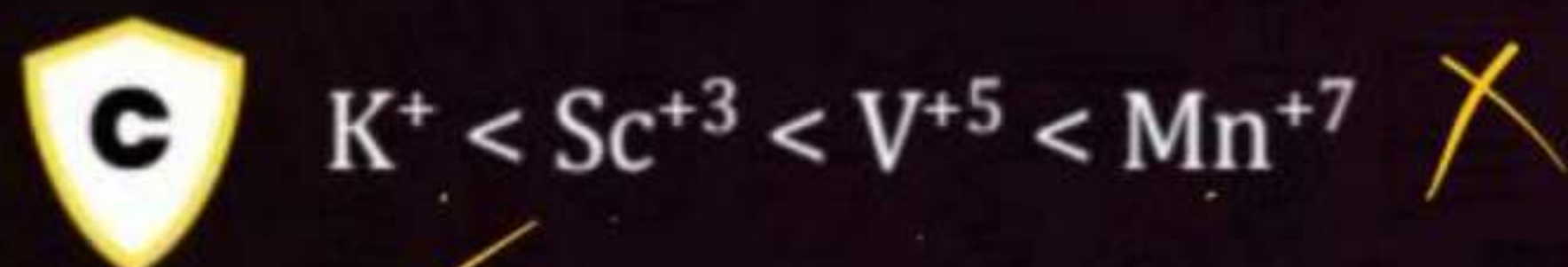
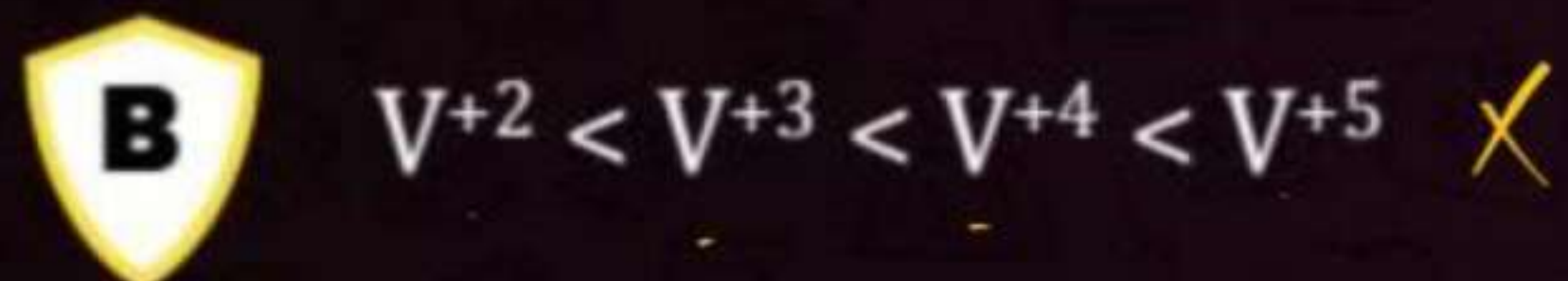
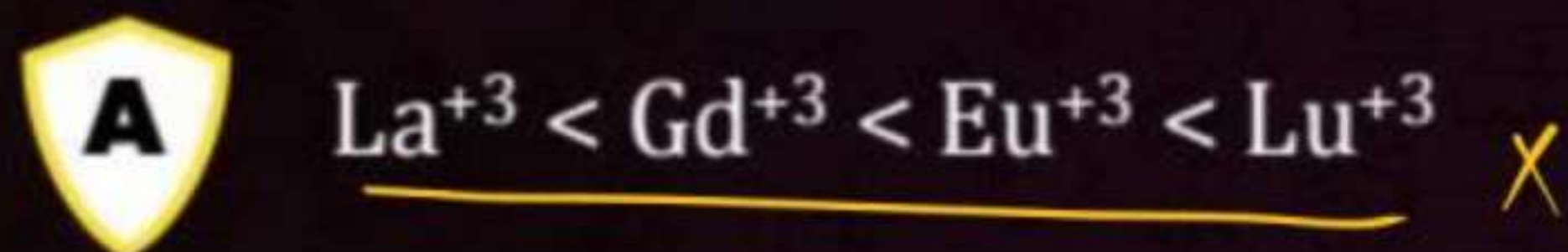
$z_{eff} \uparrow$

contraction $\uparrow$

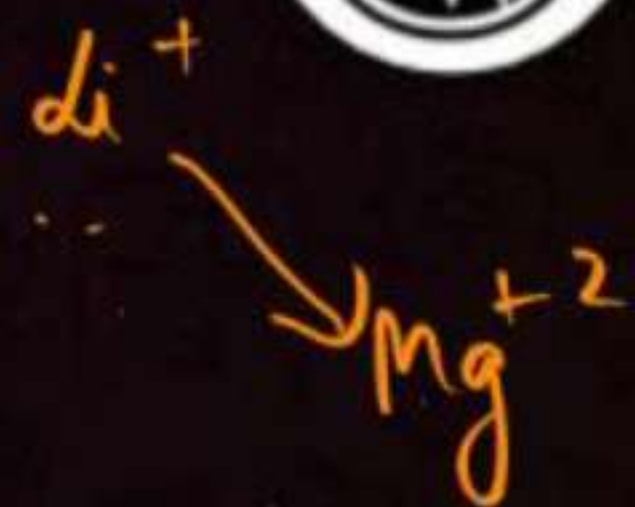
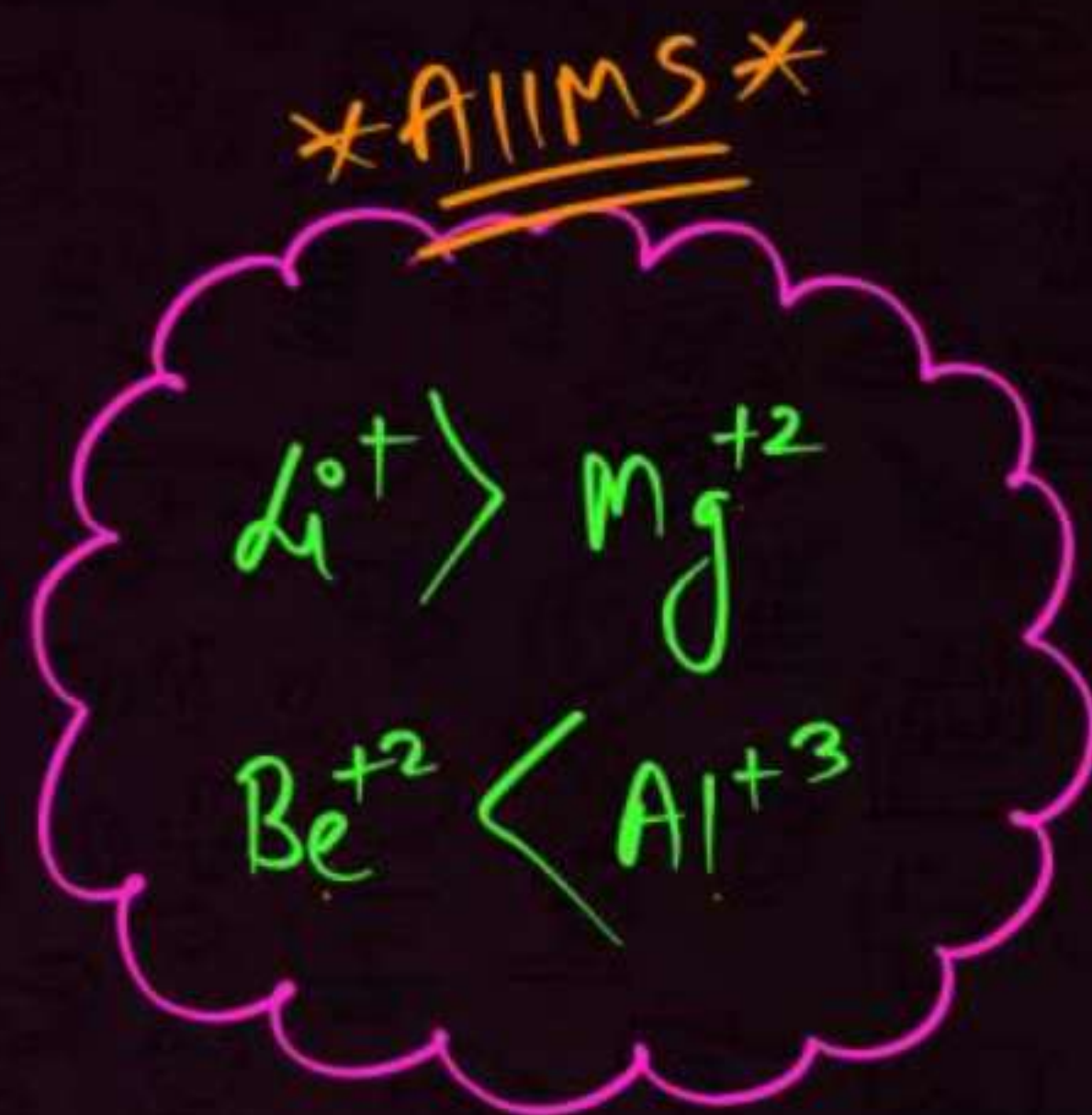
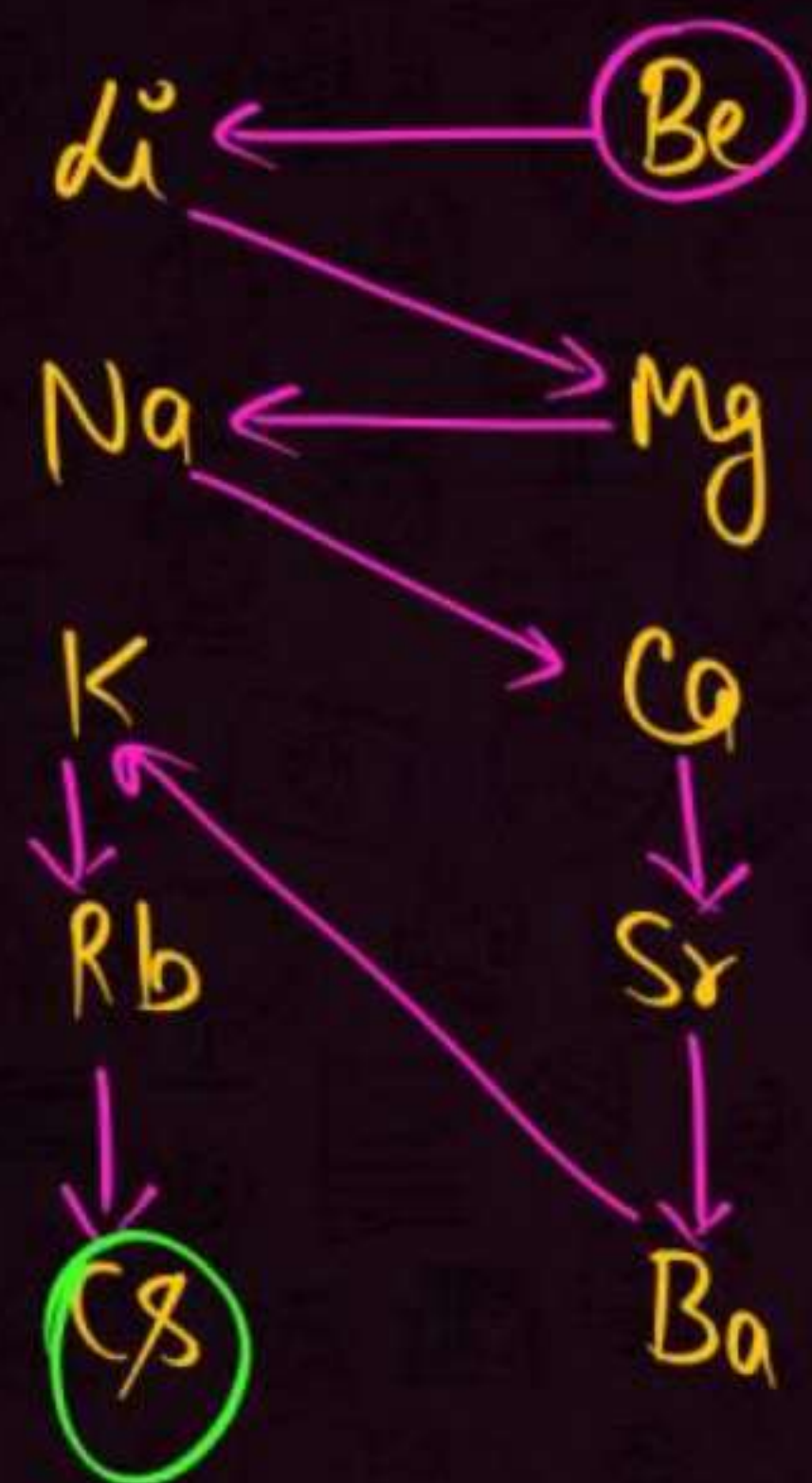
Question



Incorrect order of ionic size is:



Order of size in s-block :



$\phi = \underline{\text{Similar}}$

$$\left(\frac{\text{Charge}}{\text{Size}} \right)$$

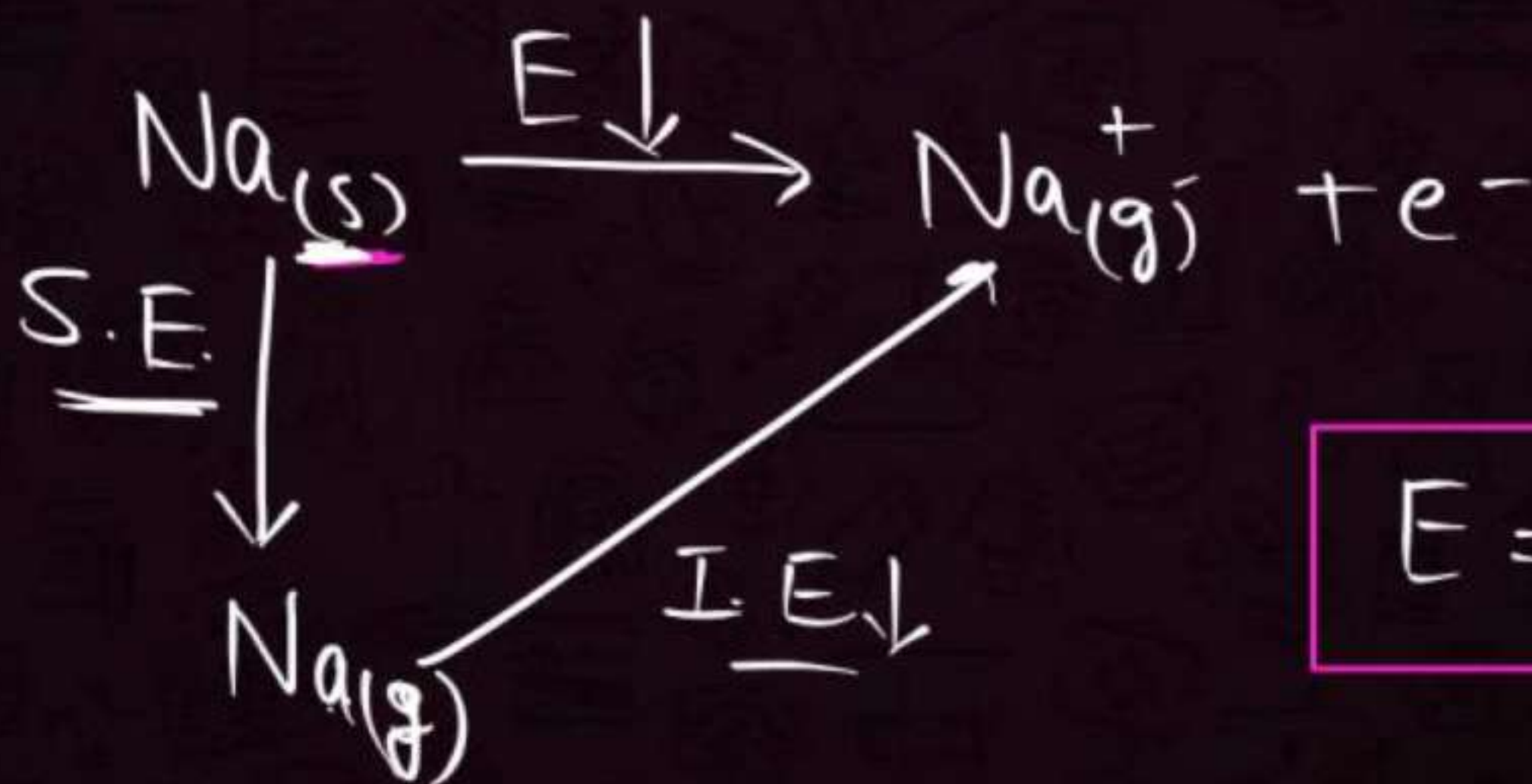
Ionic
Pot.



IONIZATION ENERGY



The amount of energy required to remove an e^- from an isolated gaseous atom is called Ionization Energy.

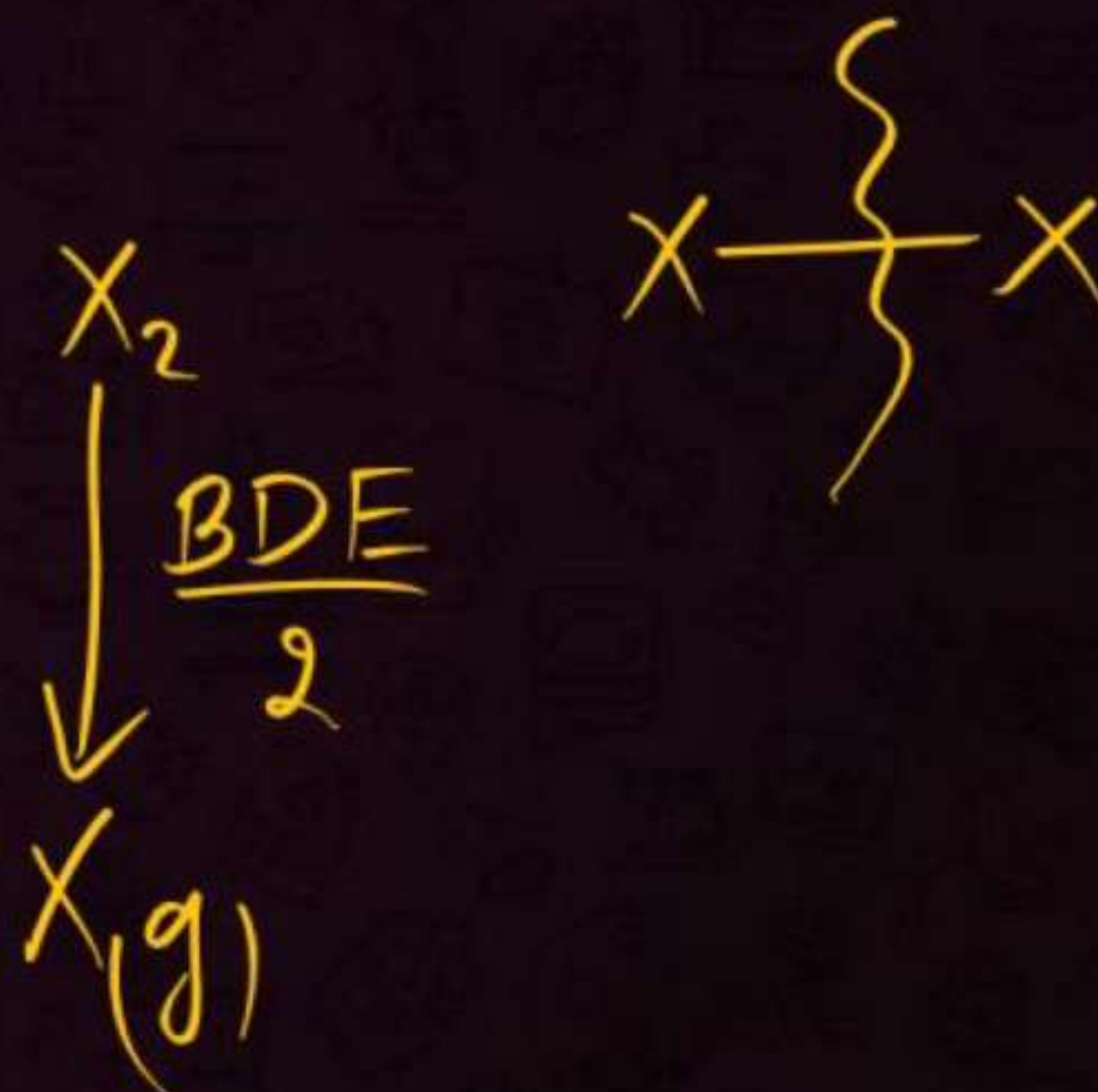
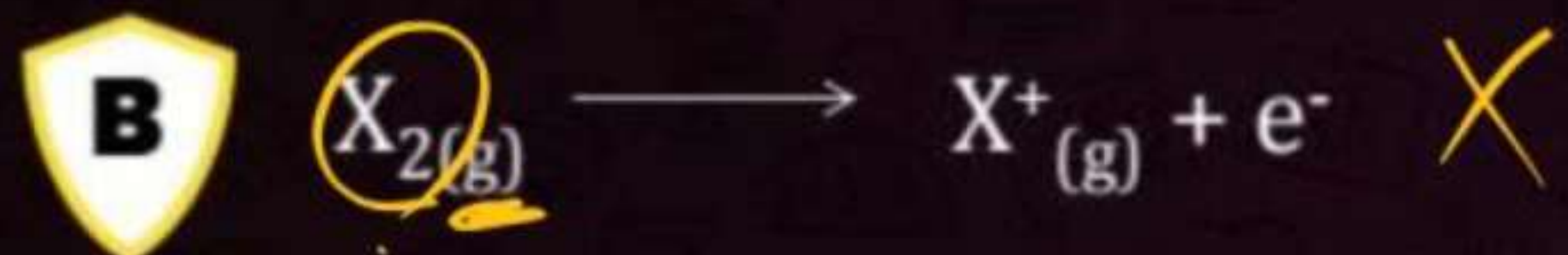


$$E = \text{S.E.} + \text{I.E.}$$

Question

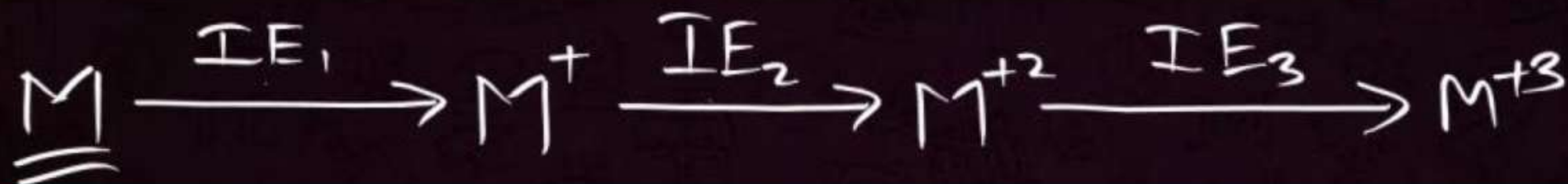


In which of the following process the energy change corresponds to Ionisation Energy ?





SUCCESSIVE IONIZATION ENERGY



$$IE_1 < IE_2 < IE_3 < IE_4 \dots IE_{(n-1)} < IE_n$$

$$IE_2 \text{ of } M = IE_1 \text{ of } M^+$$

$$IE_3 \text{ of } M = IE_1 \text{ of } M^{+2} = IE_2 \text{ of } M^+$$

* Note : $\rightarrow$ Successive I.E. are always higher.

NCERT CORNER

Energy is always required to remove electrons from an atom and hence ionization enthalpies are always positive. The second ionization enthalpy will be higher than the first ionization enthalpy because it is more difficult to remove an electron from a positively charged ion than from a neutral atom. In the same way the third ionization enthalpy will be higher than the second and so on. The term “ionization enthalpy”, if not qualified, is taken as the first ionization enthalpy.

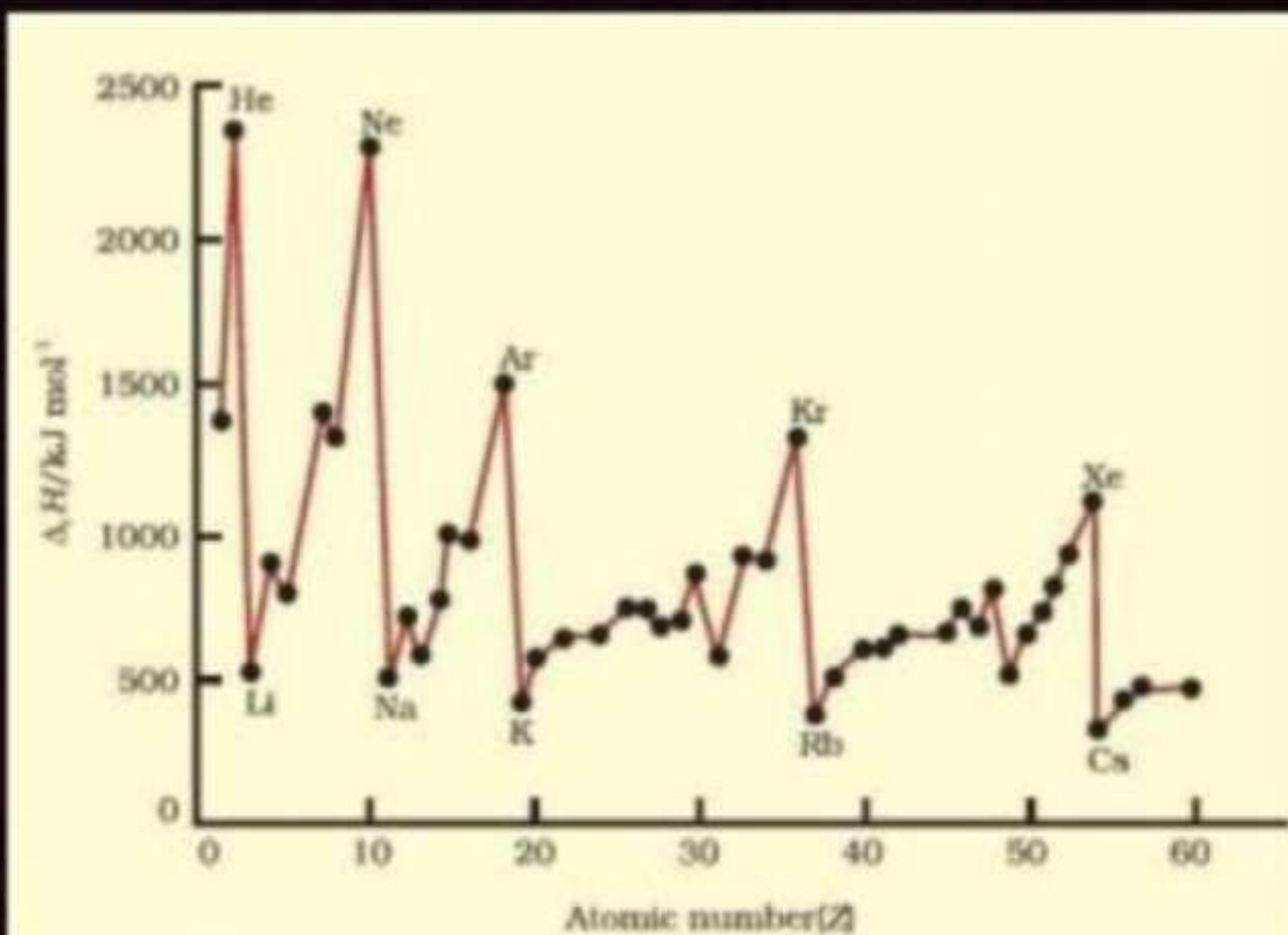


Fig. 3.5 Variation of first ionization enthalpies ($\Delta_i H$) with atomic number for elements with $Z = 1$ to 60



FACTORS AFFECTING I.E.



① Z_{eff} $\rightarrow$

$$I.E \propto Z_{eff}$$

Ex: $B < C < N$
 $Al < Si < P$

② $Size$ $\rightarrow$

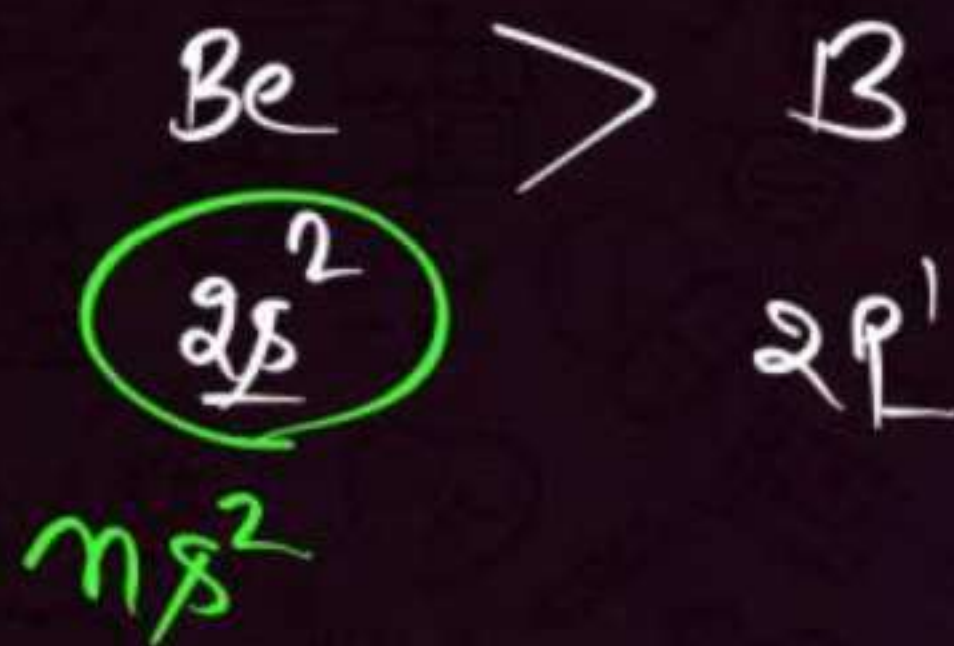
$$I.E \propto \frac{1}{Size}$$

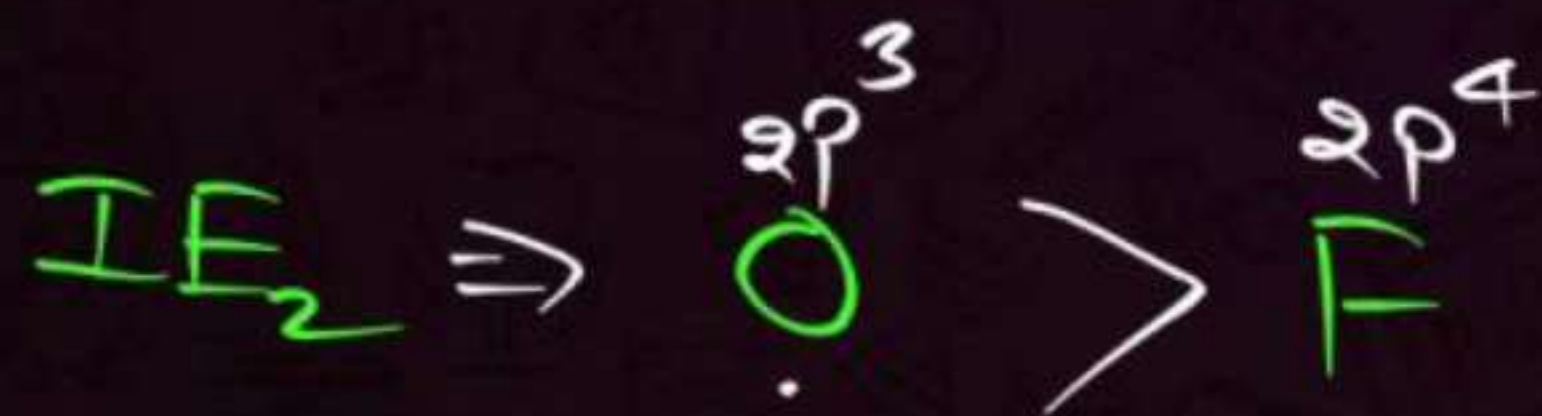
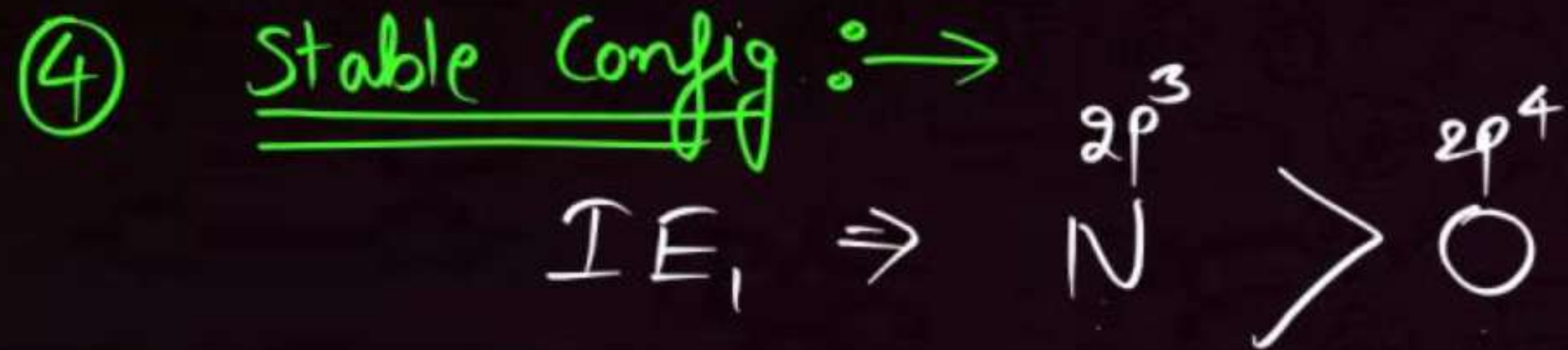
Ex: $Li > Na > K > Rb > Cs$
 $Be > Mg > Ca > Sr > Ba$

③ Penetration Power of Subshells $\rightarrow$

$\downarrow$
closeness from the nucleus

$s > p > d > f$





NCERT CORNER

Page 89

You will appreciate that the ionization enthalpy and atomic radius are closely related properties. To understand these trends, we have to consider two factors : (i) the attraction of electrons towards the nucleus, and (ii) the repulsion of electrons from each other. The effective nuclear charge experienced by a valence electron in an atom will be less than

the actual charge on the nucleus because of “**shielding**” or “**screening**” of the valence electron from the nucleus by the intervening core electrons. For example, the 2s electron in lithium is shielded from the nucleus by the inner core of 1s electrons. As a result, the valence electron experiences a net positive charge which is less than the actual charge of +3. In general, shielding is effective when the orbitals in the inner shells are completely filled. This situation occurs in the case of alkali metals which have single outermost *ns*-electron preceded by a noble gas electronic configuration.

2 Neg
2025

2nd period

Li < Be > B < C < N > O < F < Ne

(Li) < B < Be < C < O < N < F < (Ne)

Max. I.E.

(He)

Min. I.E.

(Cs)

3rd period

⇒ Na < Al < Mg < Si < S < P < Cl < Ar.

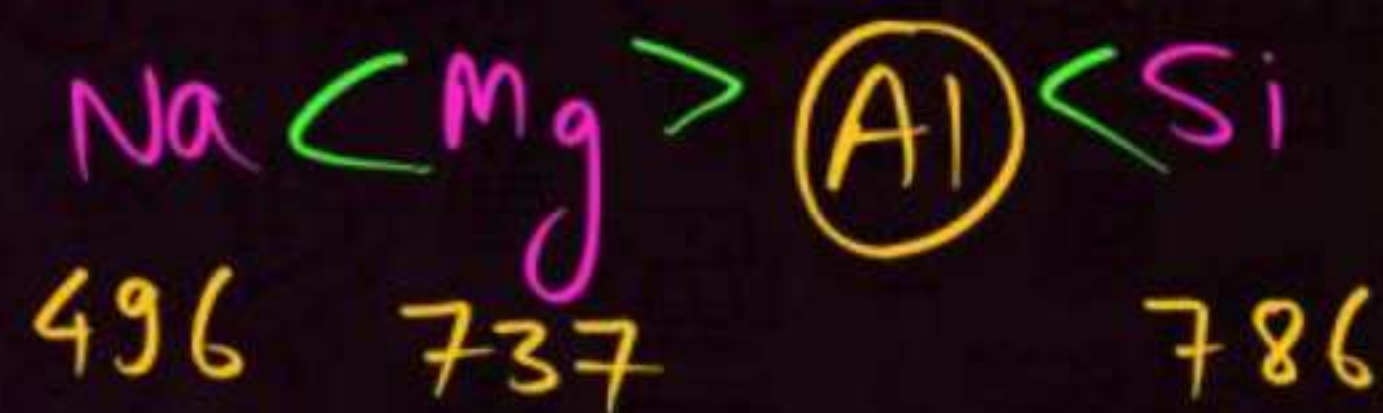
* NEET 2025 *

Problem 3.6

The first ionization enthalpy ($\Delta_i H$) values of the third period elements, Na, Mg and Si are respectively 496, 737 and 786 kJ mol⁻¹. Predict whether the first $\Delta_i H$ value for Al will be more close to 575 or 760 kJ mol⁻¹? Justify your answer.

575

760



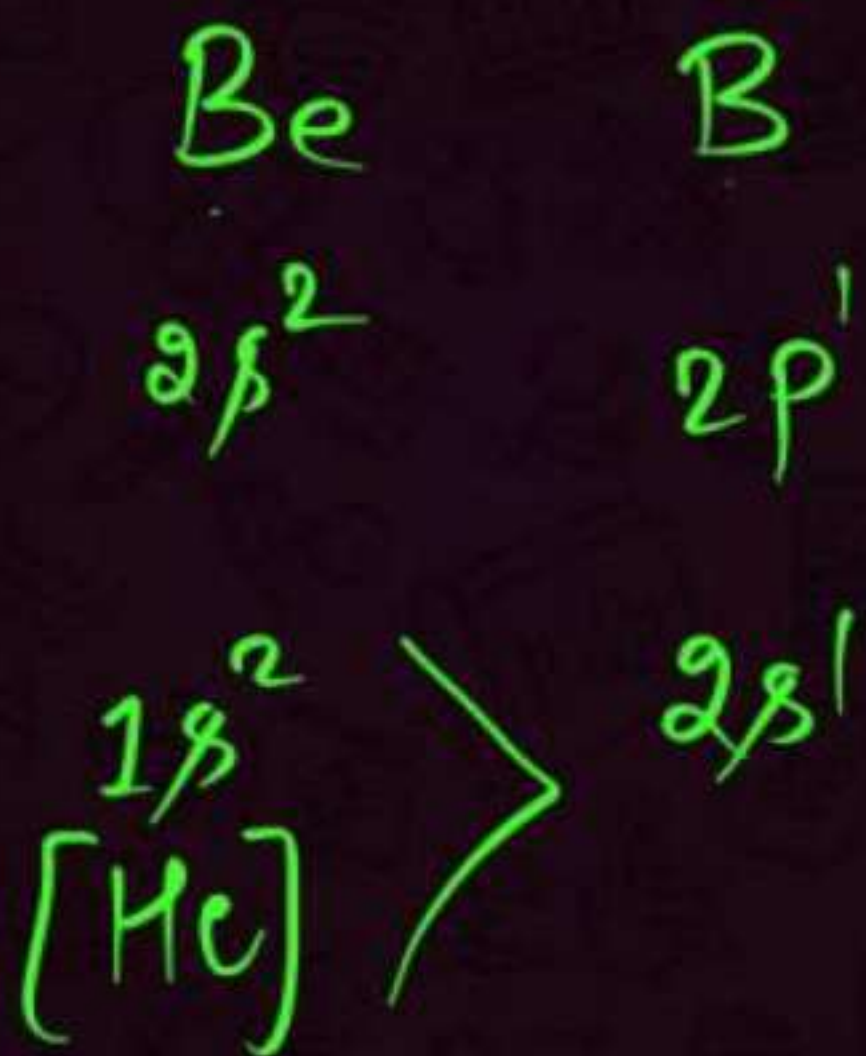
QUESTION



Among the following, third ionisation energy is highest for:

(AMU 2010)

- ☐ A magnesium
- ☐ B boron
- ☒ C beryllium
- ☐ D aluminium

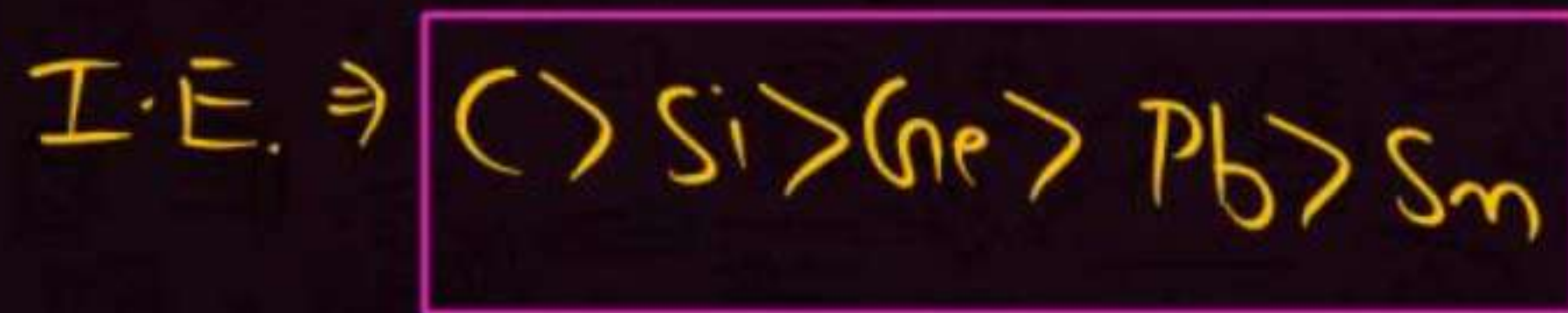


NCERT CORNER

Page 89

From Fig. 3.6(a), you will also notice that the first ionization enthalpy of boron ($Z = 5$) is slightly less than that of beryllium ($Z = 4$) even though the former has a greater nuclear charge. When we consider the same principal quantum level, an s -electron is attracted to the nucleus more than a p -electron. In beryllium, the electron removed during the ionization is an s -electron whereas the electron removed during ionization of boron is a p -electron. The penetration of a $2s$ -electron to the nucleus is more than that of a $2p$ -electron; hence the $2p$ electron of boron is more shielded from the nucleus by the inner core of electrons than the $2s$ electrons of beryllium. Therefore, it is easier

to remove the $2p$ -electron from boron compared to the removal of a $2s$ -electron from beryllium. Thus, boron has a smaller first ionization enthalpy than beryllium. Another “anomaly” is the smaller first ionization enthalpy of oxygen compared to nitrogen. This arises because in the nitrogen atom, three $2p$ -electrons reside in different atomic orbitals (Hund’s rule) whereas in the oxygen atom, two of the four $2p$ -electrons must occupy the same $2p$ -orbital resulting in an increased electron-electron repulsion. Consequently, it is easier to remove the fourth $2p$ -electron from oxygen than it is, to remove one of the three $2p$ -electrons from nitrogen.



$N > P > As > Sb > Bi$

$O > S > Se > Te > Po$

$F > Cl > Br > I > At$

* d-block $\rightarrow$

$3d > 4d < 5d$

($z_{eff} \uparrow$)
L.C.

$5d > 3d > 4d$

Ex: $Cu > Ag < Au$



APPLICATIONS OF I.E.



① Metallic/Non-Metallic Character :

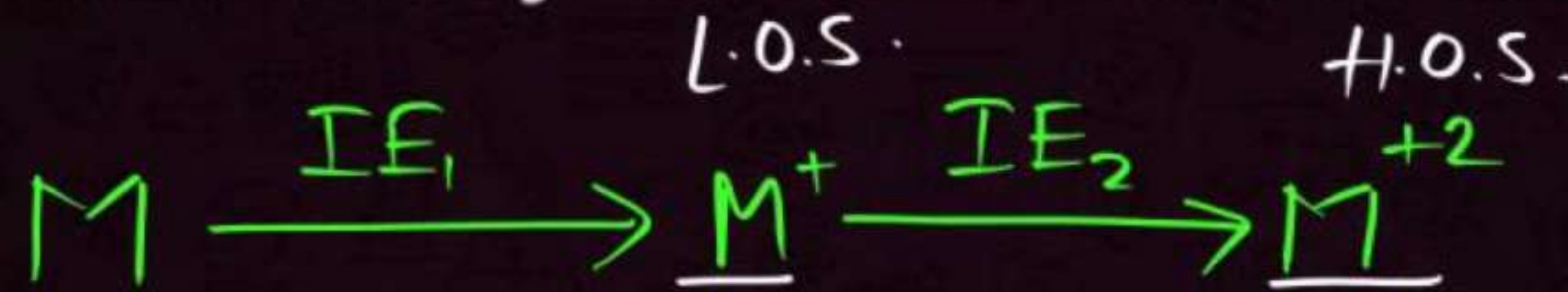
$$\text{Metallic Nature} \propto \frac{1}{\text{I.E.}}$$

Ex:



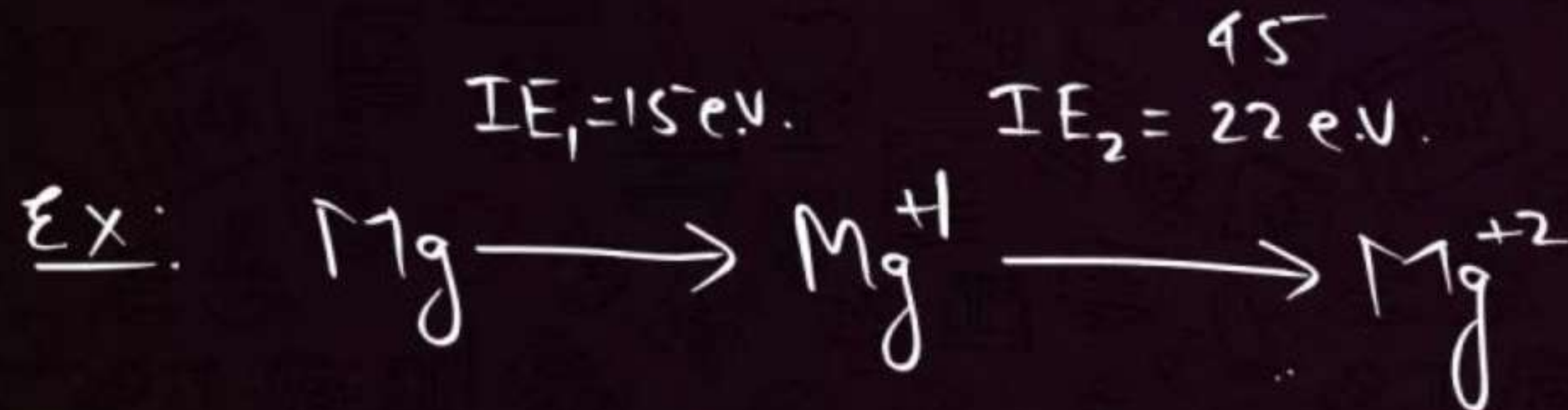
Size $\uparrow$ I.E. $\downarrow$ Metallic Nature $\uparrow$

② Comparison in stability of oxidation states :



$$\Delta IE = (IE_2 - IE_1) < 11 \text{ e.V.} \Rightarrow \text{H.O.S. More Stable}$$

$$> 16 \text{ e.V.} \Rightarrow \text{L.O.S. More Stable}$$



$$\Delta IE = 22 - 15 = \underline{\underline{7 \text{ e.V.}}}$$

③ ^{*Imp} No. of valence e^- $\rightarrow$

A $\rightarrow$ 15 e.v., 22 e.v., 45 e.v.,

165 e.v., 195 e.v. ...

Sudden large
jump b/w I.E values

Valence $e^- = 3e^-$

Most stable O.S. = +3

oxide = A_2O_3

Nitride = AN

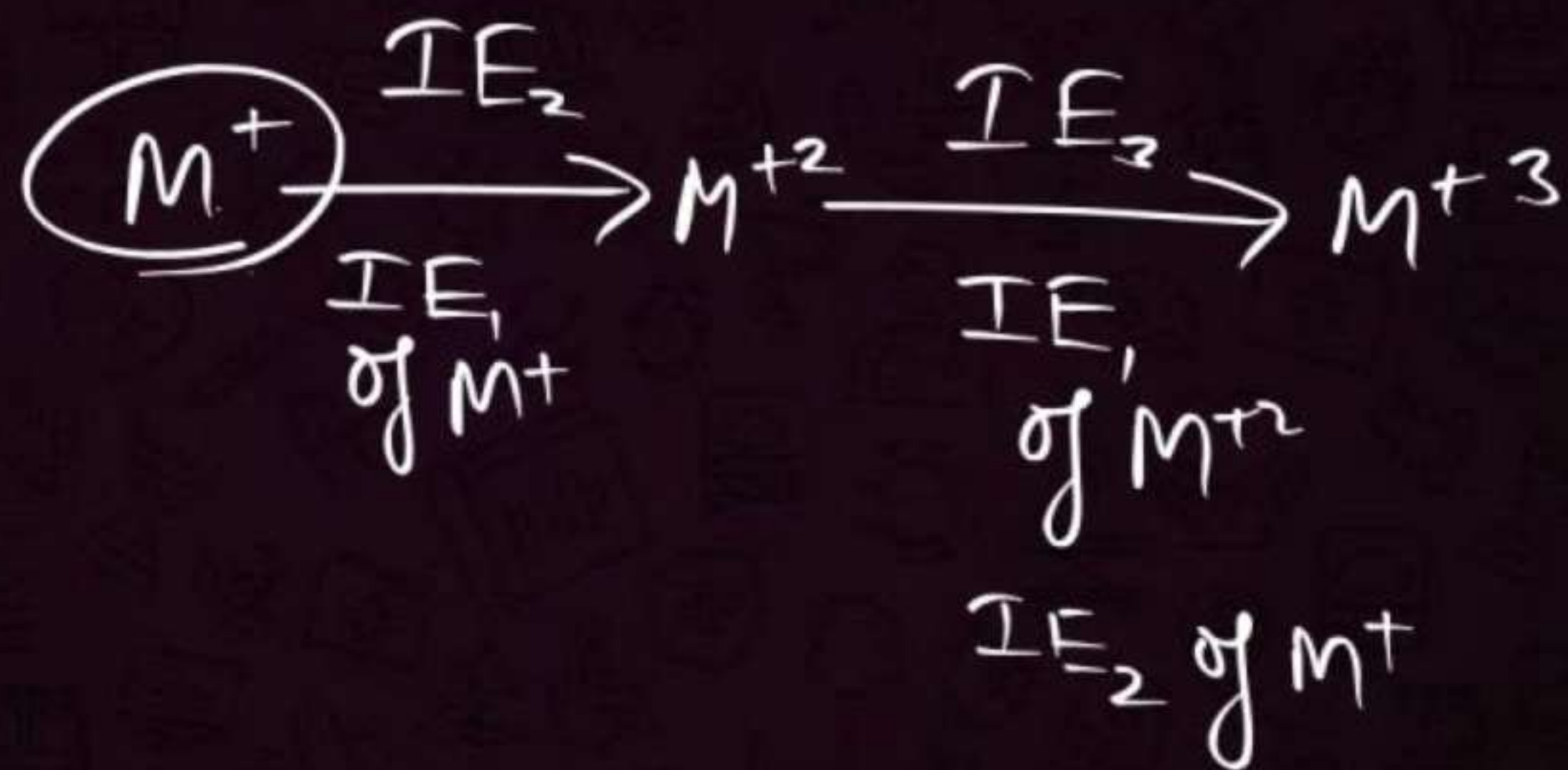
Halide = AX_3

Hydride = AH_3

Q Element A?

① Mg ~~② Al~~

③ Si ④ N



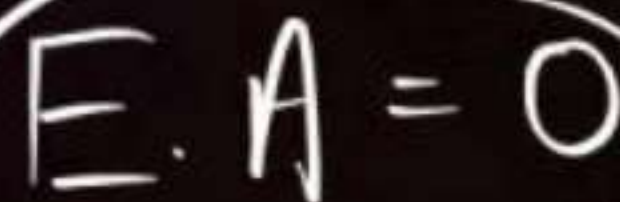
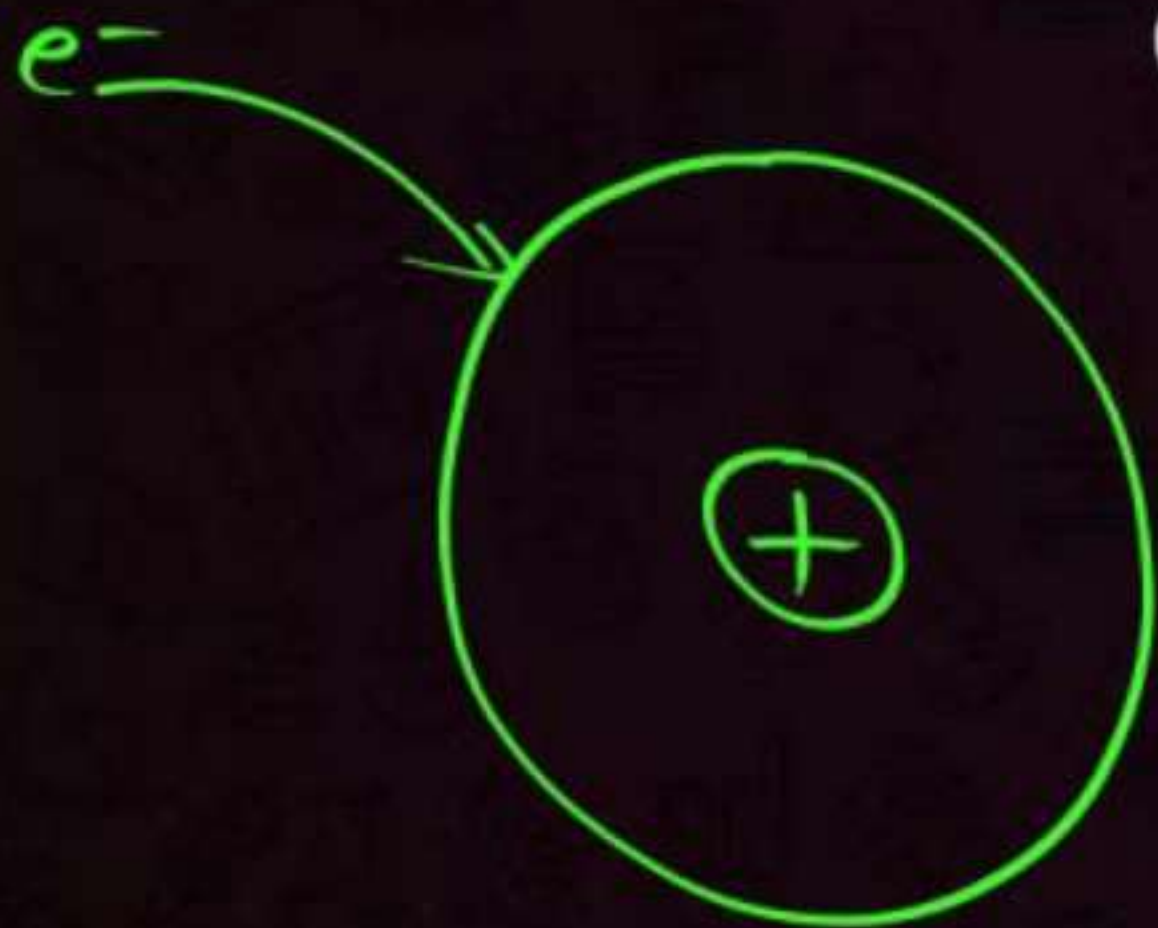


ELECTRON AFFINITY

चाहत

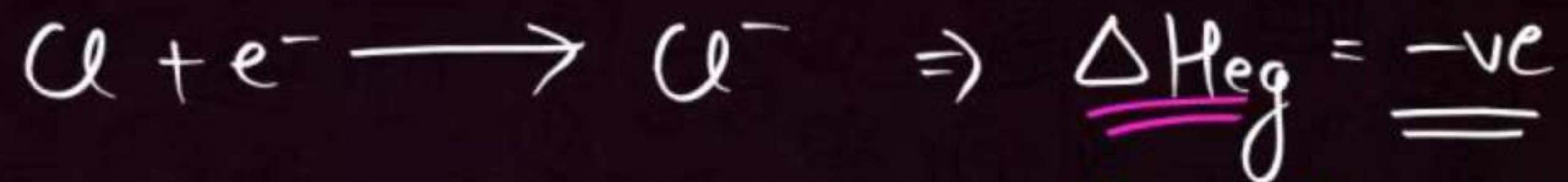


The amt. of energy released when an e^- is added in an isolated gaseous atom.





e^- gain enthalpy (ΔH_{eg})



More the energy released
then more -ve will be the
value of ΔH_{eg} (higher the
E.A.)



More +ve value of ΔH_{eg}
represents lesser the tendency
to accept e^- & hence lesser will be
the E.A.

$$\Delta H_{eg} = -E \cdot A.$$



$$\Delta H_{eg} = -A_e + \frac{5}{2}RT$$

* Successive E.A. $\therefore \downarrow$



$$\Delta H_{eg_1} = \underline{\underline{-144 \text{ KJ/mol}}}$$



$$\Delta H_{eg_2} = \underline{\underline{+744 \text{ KJ/mol}}}$$



$$\Delta H_{\text{total}} = \underline{\underline{+600 \text{ (endo.)}}}$$



} endo.



FACTORS AFFECTING ELECTRON AFFINITY

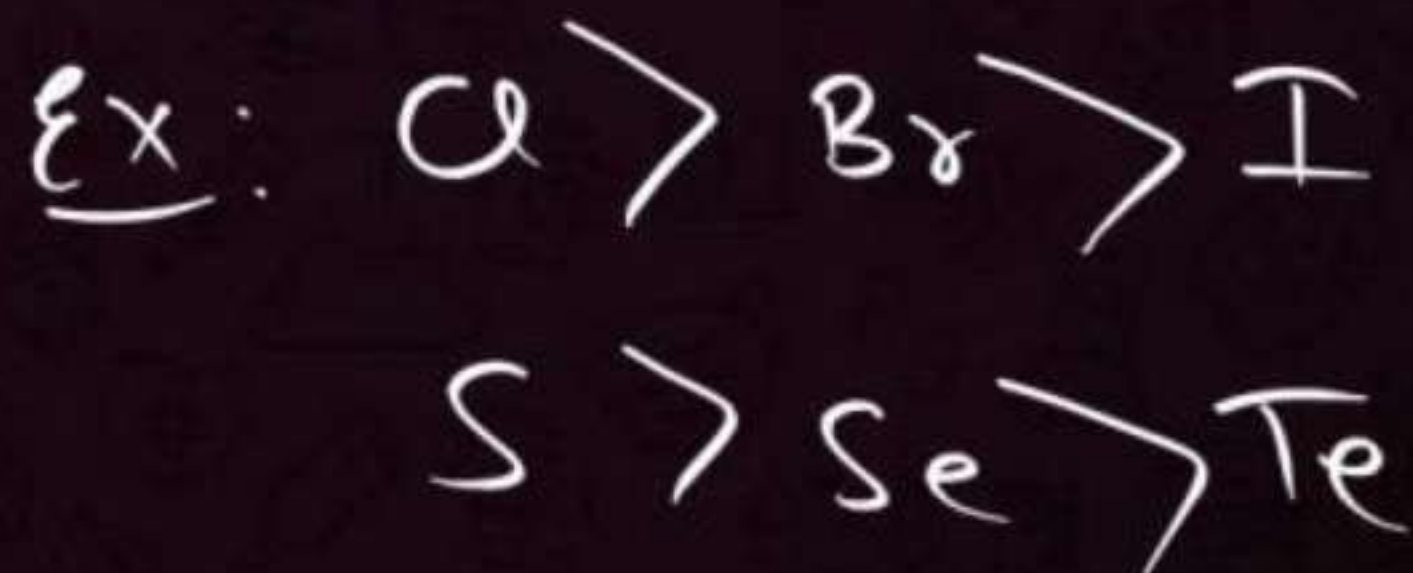


① z_{eff} $\rightarrow$

$$E.A \propto z_{eff}$$

② size $\rightarrow$

$$E.A. \propto \frac{1}{size}$$



2nd period

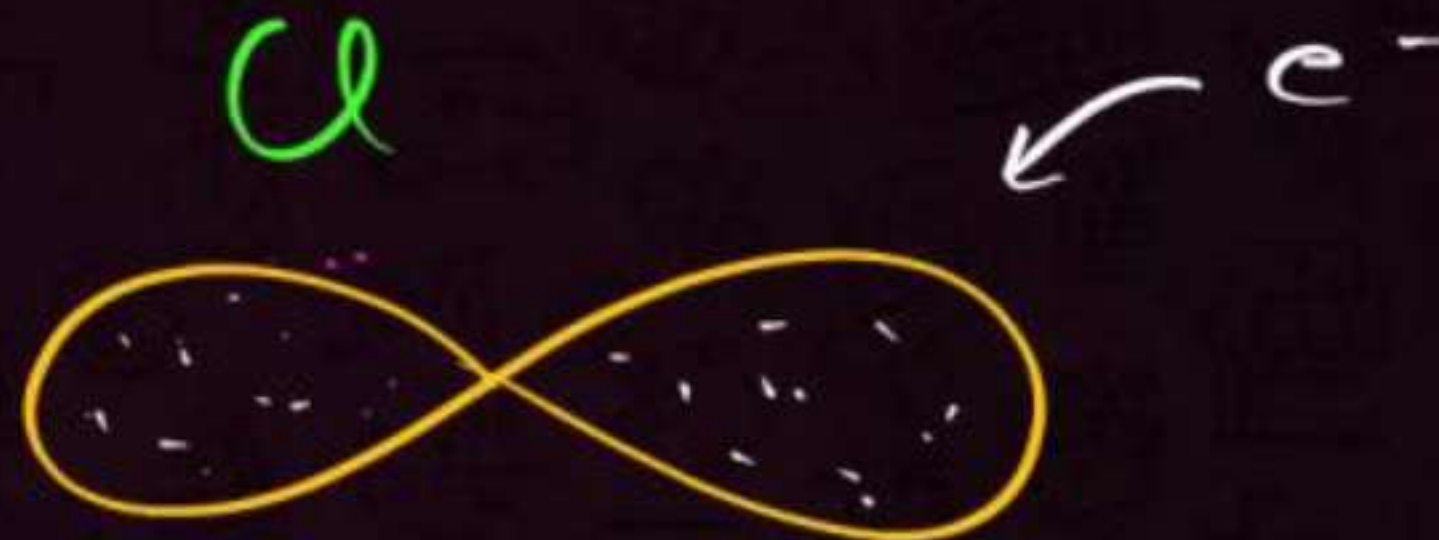
F



(inter e⁻ rep.
more)

3rd period

Cl



F < Cl

O < S

N < P

C < Si

ex: F < Cl > Br > I

O < S > Se > Te

~~imp~~

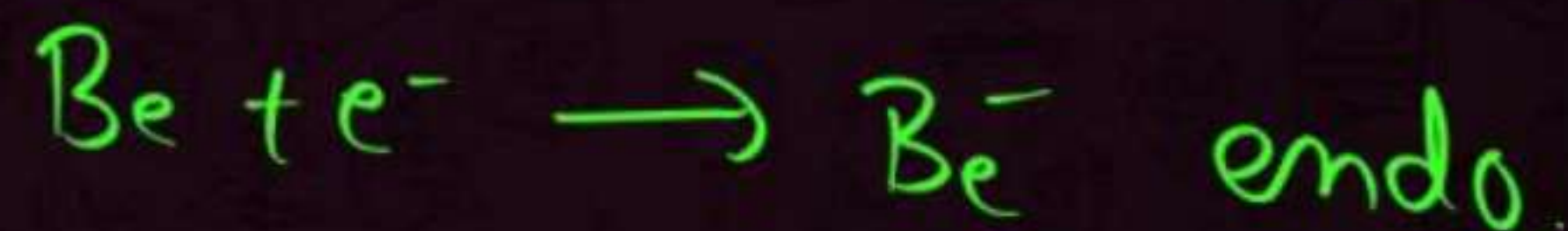
Cl > F > Br > I

S > Se > Te > Po > O

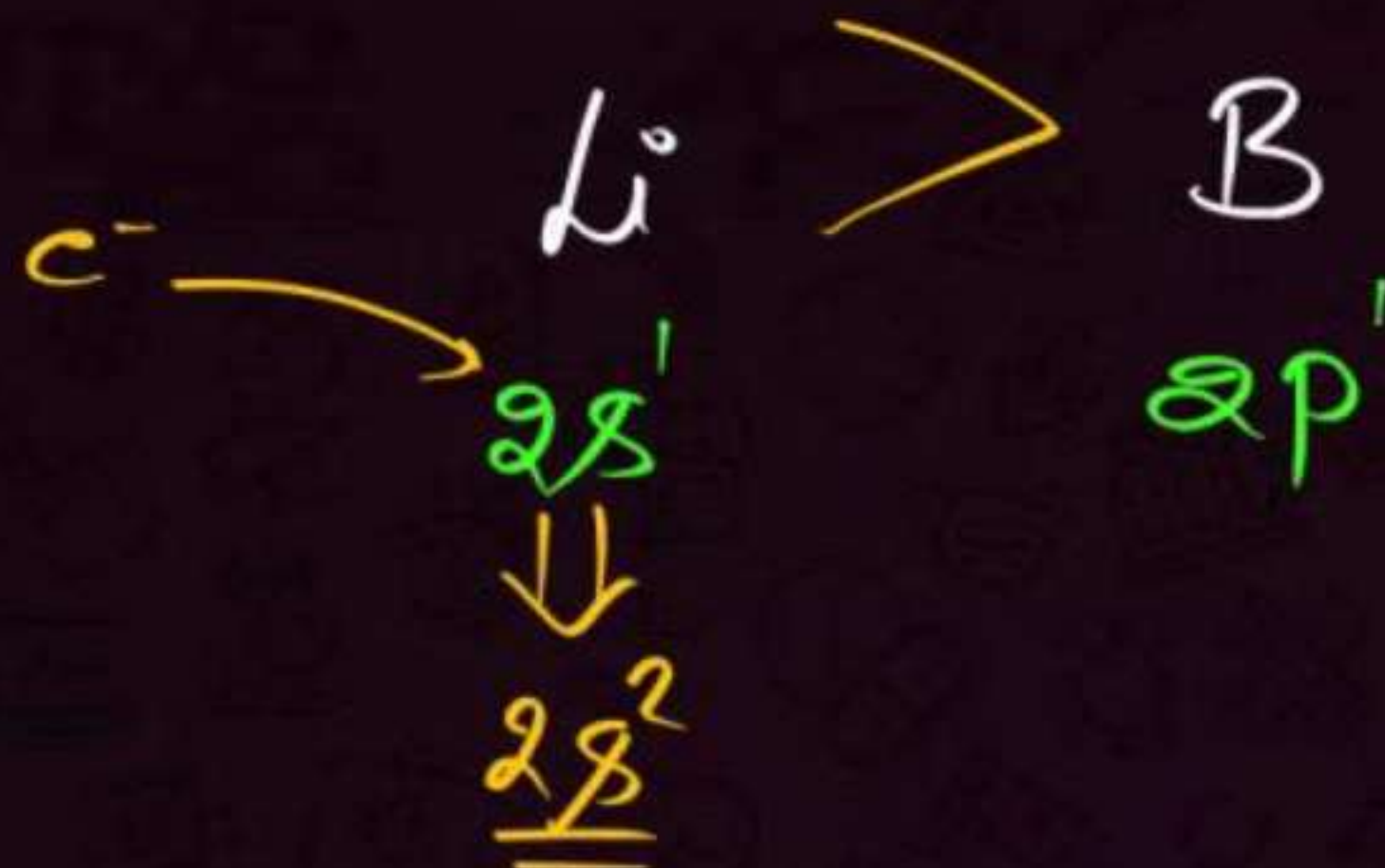


③ Stable Config. $\therefore \rightarrow$

$\Delta H_{eg,1} = +ve$ for Noble gases $\underbrace{Be, Mg, N}_{ns^2} \underbrace{N}_{p^3}$



④ Penetration Power $\rightarrow$



2nd period $\rightarrow$



* 3rd period $\Rightarrow$



H.W.

$$\text{Max. E.A} = \underline{\underline{0}}$$

QUESTION

Which of the following processes involves absorption of energy?

- A** $S(g) + e^{-} \rightarrow S^{-}(g)$
- B** $S^{-} + e^{-} \rightarrow S^{2-}(g)$
- C** $Cl(g) + e^{-} \rightarrow Cl^{-}(g)$
- D** None of these

Ans. 0

QUESTION



Which of following will have the most negative electron gain enthalpy and which the least negative ? F, P, S, Cl

A ~~P, Cl~~

B Cl, F

C Cl, S

~~**D**~~ Cl, P

Ans. 0

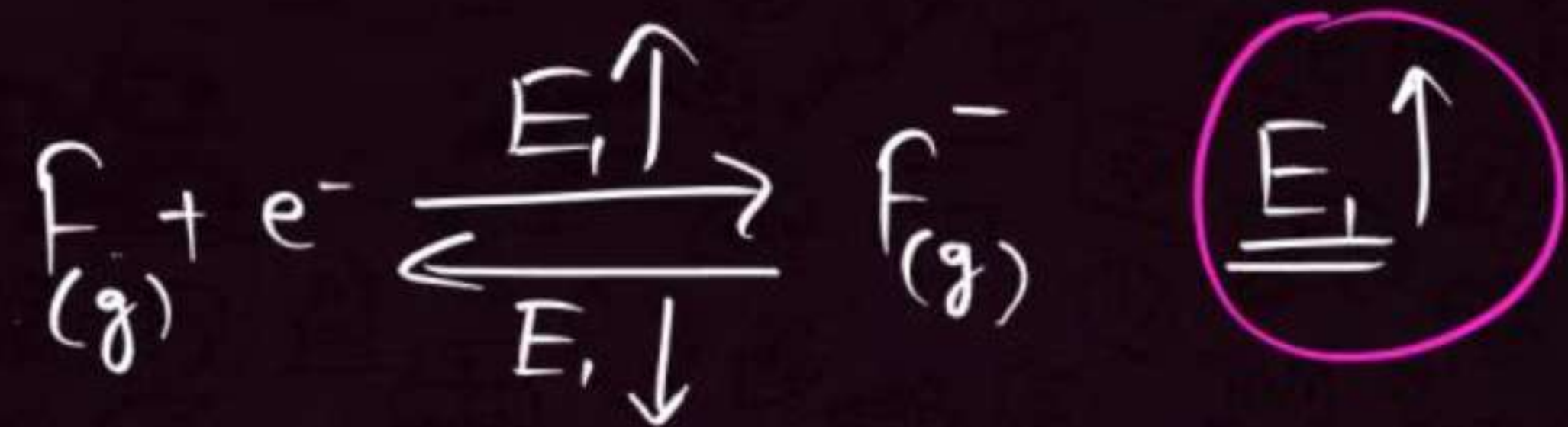
$N, P, S, O, U, F, \underline{C}, \underline{Si}$

* Hitsick *

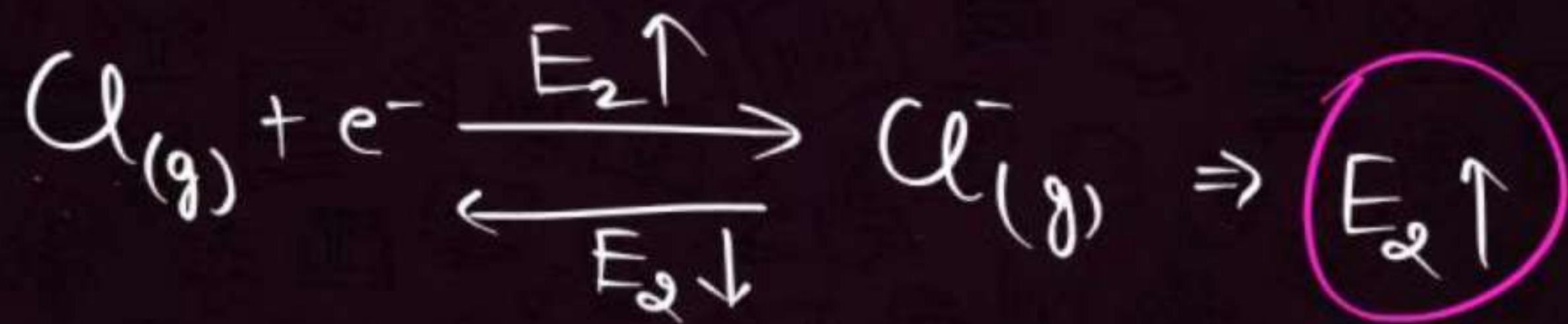
$N\text{-family} < C\text{-family} < O\text{-family} < X\text{-family}$

$N < P < C < Si < O < S < F < U$

Relation b/w I.E & E.A : $\rightarrow$



$$E_1 = \text{E.A. of } F_{(g)} = \text{I.E. of } F_{(g)}^-$$



$$E_2 = \text{E.A. of } Cl_{(g)} = \text{I.E. of } Cl_{(g)}^-$$

I.E. of $\underline{a^-} > \underline{F^-}$

I.E. of $a < F$

I.E. order

F, a, F^-, a^-

$F > a > a^- > F^-$

QUESTION



Which of the following is correct order of Ionization energy of $\{O, O^{-1}, S, S^{-1}\}$?

$$O > S > \underline{S^{-1}} > \underline{O^{-1}}$$

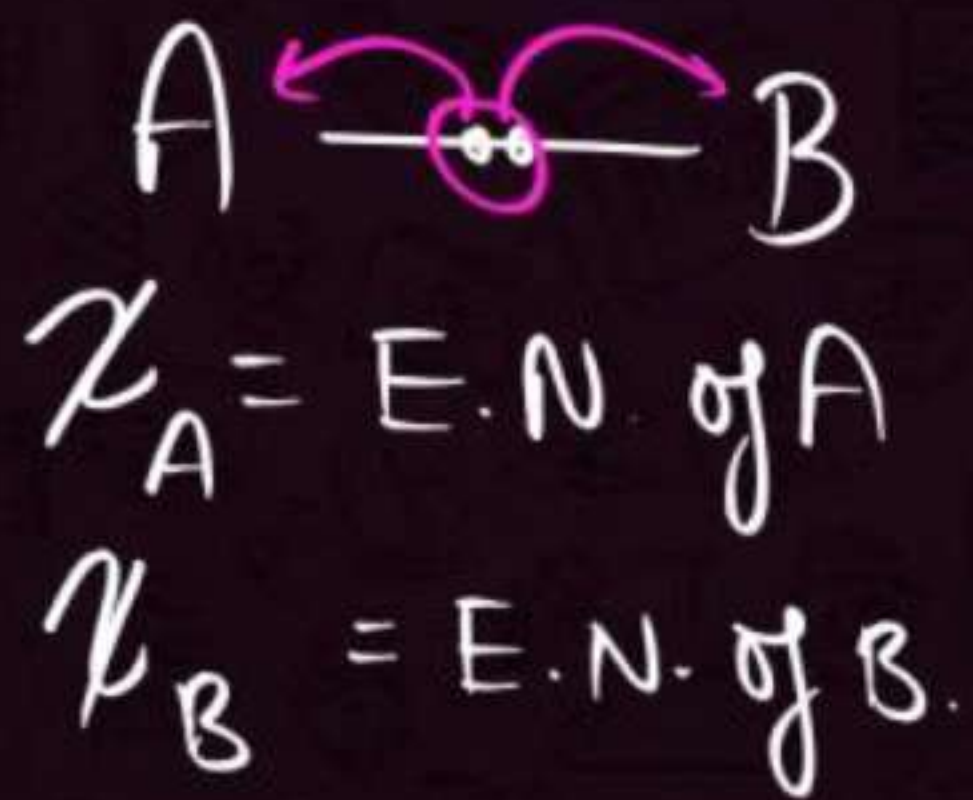
- A** $O > S > O^{-1} > S^{-1}$
- ~~**B** $O > S > S^{-1} > O^{-1}$~~
- C** $S^{-1} > O^{-1} > O > S$
- D** $S^{-1} > S^{+} > O^{-1} > O$

Ans. 0

ELECTRONEGATIVITY



It is the tendency of a covalently bonded atom to attract shared pair of e^-



if $\chi_A > \chi_B$



* Note $\rightarrow$ E.N. does not depends on e^- config



SCALES OF E.N.



① Mullikan's Scale $\therefore \rightarrow$

$$E.N. = \frac{E.A + I.E.}{2}$$

② Pauling Scale $\therefore \rightarrow$ based on Bond Energy.

$$\Delta EN \propto \sqrt{\Delta}$$

$\rightarrow$ ionic Resonance Energy

(F) $\rightarrow$ Reference
4.0

Li 1.0	Be 1.5	B 2.0	C 2.5	N 3.0	O 3.5	F 4.0
Na 0.9	Mg 1.2	Al 1.5	Si 1.8	P 2.1	S 2.5	Cl 3.0
K 0.8						Br 2.8
Rb 0.8						
Cs 0.7						I 2.5

$$\underline{H} = \underline{P} = \underline{2.1}$$

$$\underline{N} = \underline{Cl} = \underline{3.0}$$

$$\underline{C} = \underline{S} = \underline{I} = 2.5$$

B 2.0 Al 1.5 Ga 1.6 In 1.7 Tl 1.8

Table 3.8(a) Electronegativity Values (on Pauling scale) Across the Periods

Atom (Period II)	Li	Be	B	C	N	O	F
Electronegativity	1.0	1.5	2.0	2.5	3.0	3.5	4.0
Atom (Period III)	Na	Mg	Al	Si	P	S	Cl
Electronegativity	0.9	1.2	1.5	1.8	2.1	2.5	3.0

Table 3.8(b) Electronegativity Values (on Pauling scale) Down a Family

Atom (Group I)	Electronegativity Value	Atom (Group 17)	Electronegativity Value
Li	1.0	F	4.0
Na	0.9	Cl	3.0
K	0.8	Br	2.8
Rb	0.8	I	2.5
Cs	0.7	At	2.2

QUESTION



The electronegativity of the following elements increases in the order

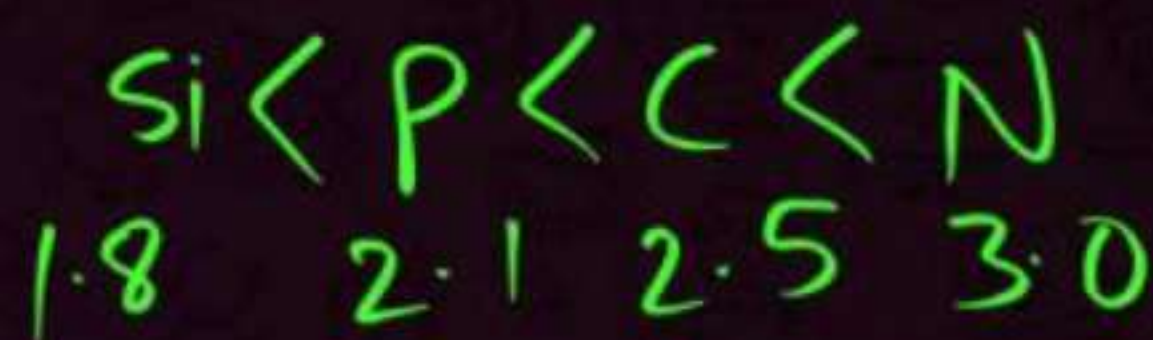
(AFMC 2010)

A C, N, Si, P

B N, Si, C, P

~~**C** Si, P, C, N~~

D P, Si, N, C



Ans. ()



FACTORS AFFECTING E.N.



① Z_{eff} $\rightarrow$

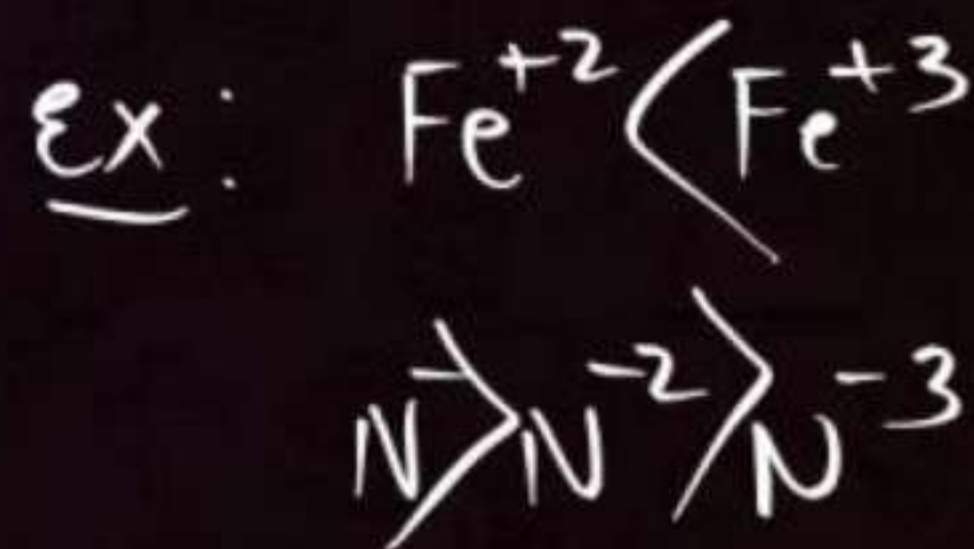
$$E.N. \propto Z_{eff}$$

② size $\rightarrow$

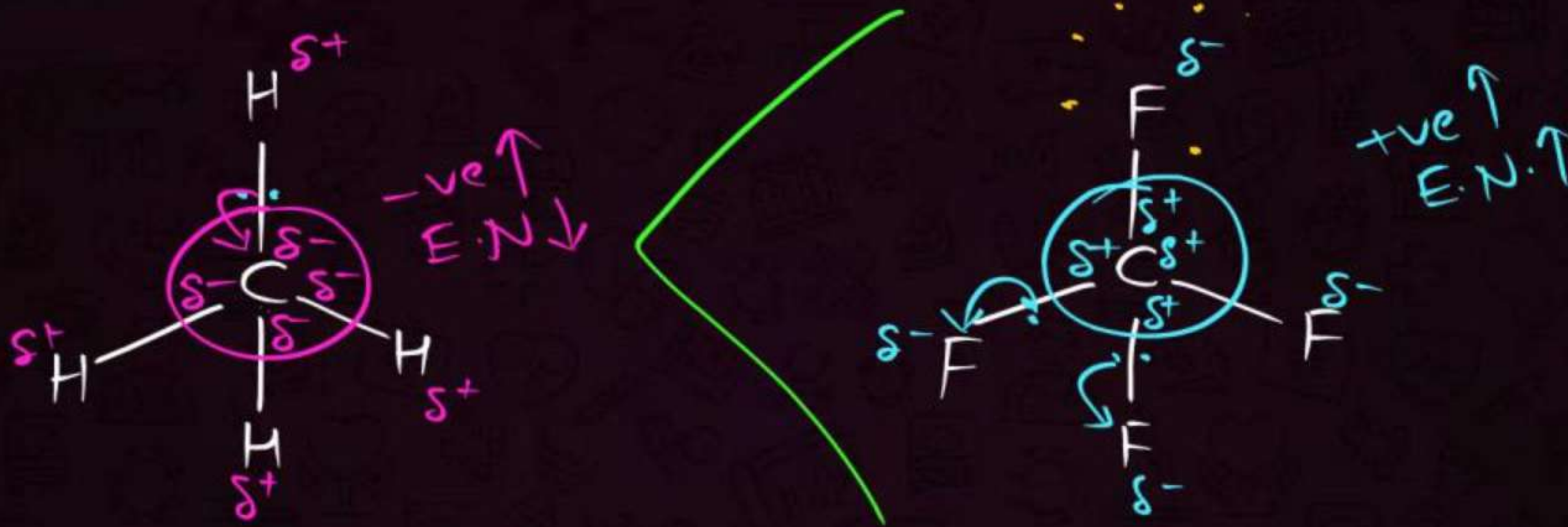
$$E.N. \propto \frac{1}{\text{size}}$$

③ +ve/-ve charge $\rightarrow$

$$E.N. \propto \frac{+ve}{-ve}$$



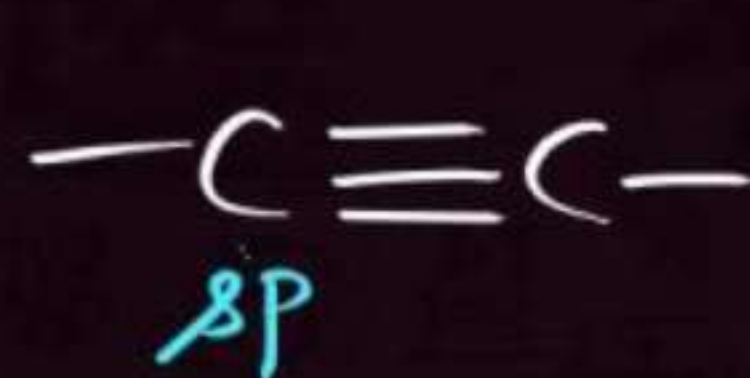
④ E.N. of Terminal atoms: $\rightarrow$



$\text{E.N.} \propto \text{E.N. of Terminal atom}$

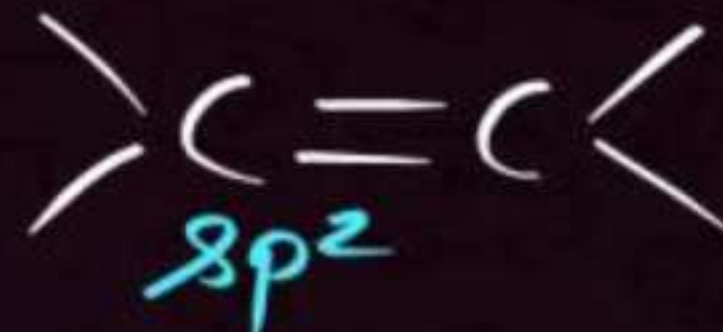
⑤ % s character : $\rightarrow$

E.N. $\propto$ % s char



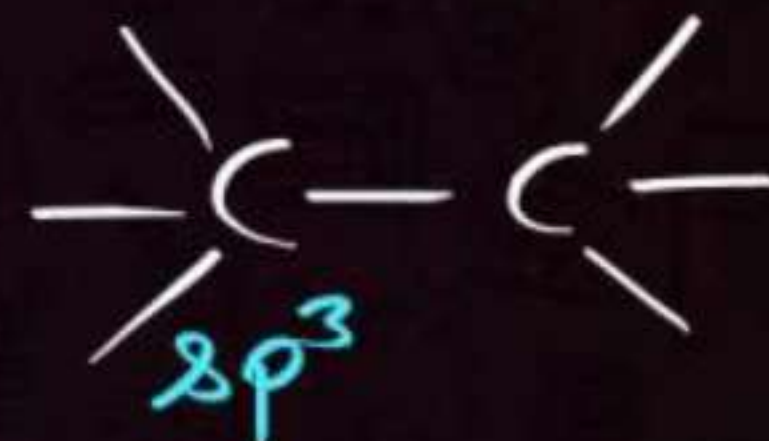
50% s

3.25



33.3% s

2.8



25% s

2.5



APPLICATIONS OF E.N.



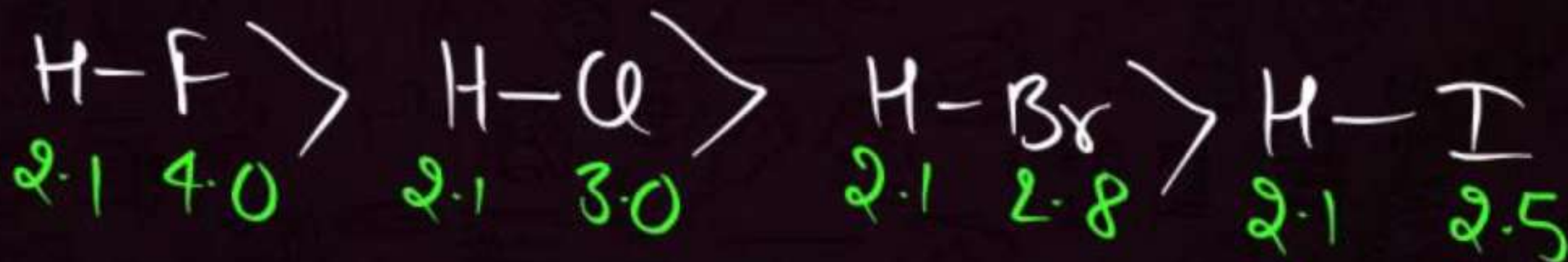
BOND POLARITY:



$\Delta EN = 0 \Rightarrow$ Non Polar Bond

$\Delta EN \neq 0 \Rightarrow$ Polar.

Ex: Polarity order



② METALLIC/NON-METALLIC NATURE :

$$\text{Metallic Nature} \propto \frac{1}{\text{E.N.}}$$

In a Group Metallic Nature ↑ se from Top to Bottom

$$\text{Non Metallic Nature} \propto \text{E.N.}$$

in a period Non Metallic Nature ↑ se.

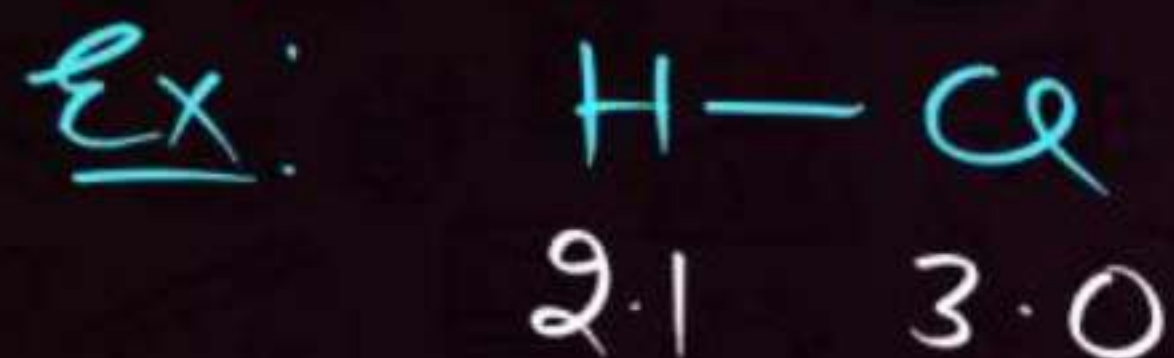
③ IUPAC NAMING OF INORGANIC COMPOUNDS:



oxygen difluoride ✓✓

④ % IONIC CHARACTER :

$$\% \text{I.C.} = 16 \times (\Delta \text{EN}) + 3.5 (\Delta \text{EN})^2$$

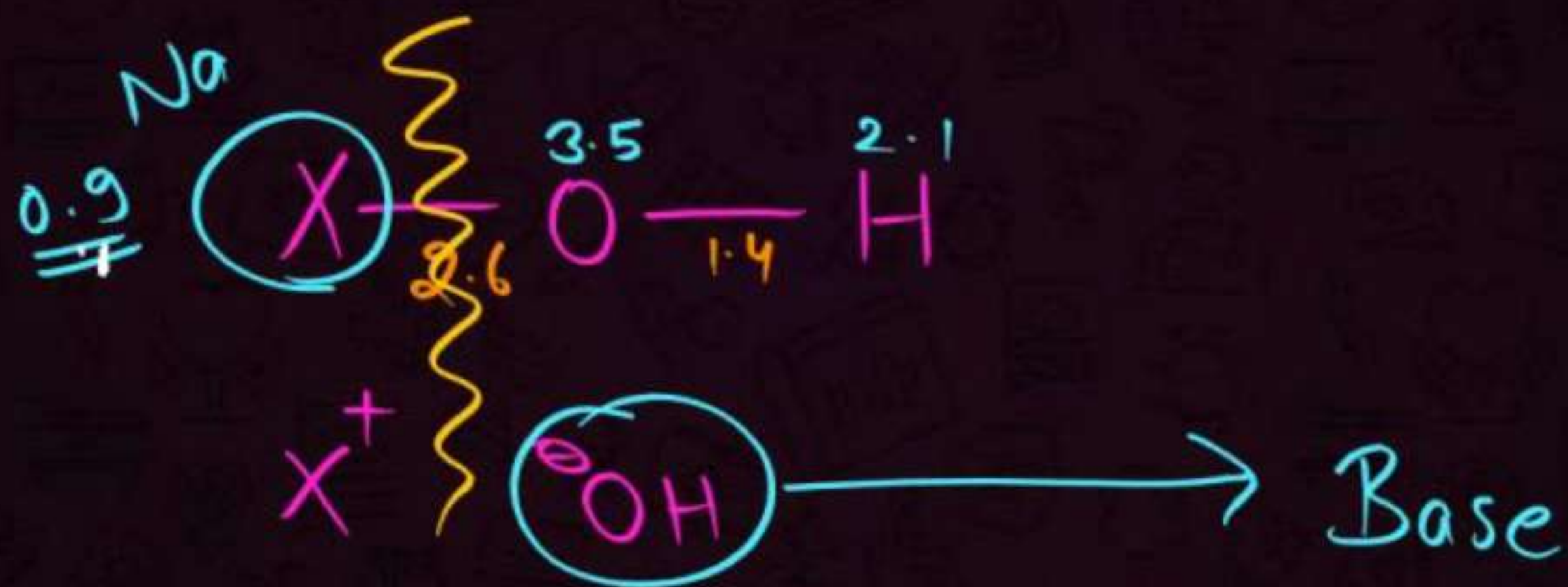
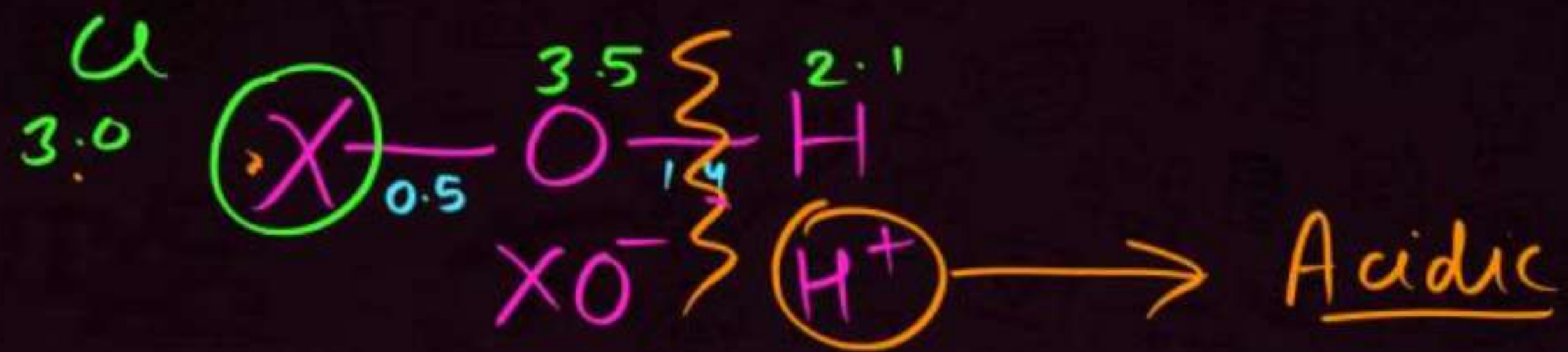


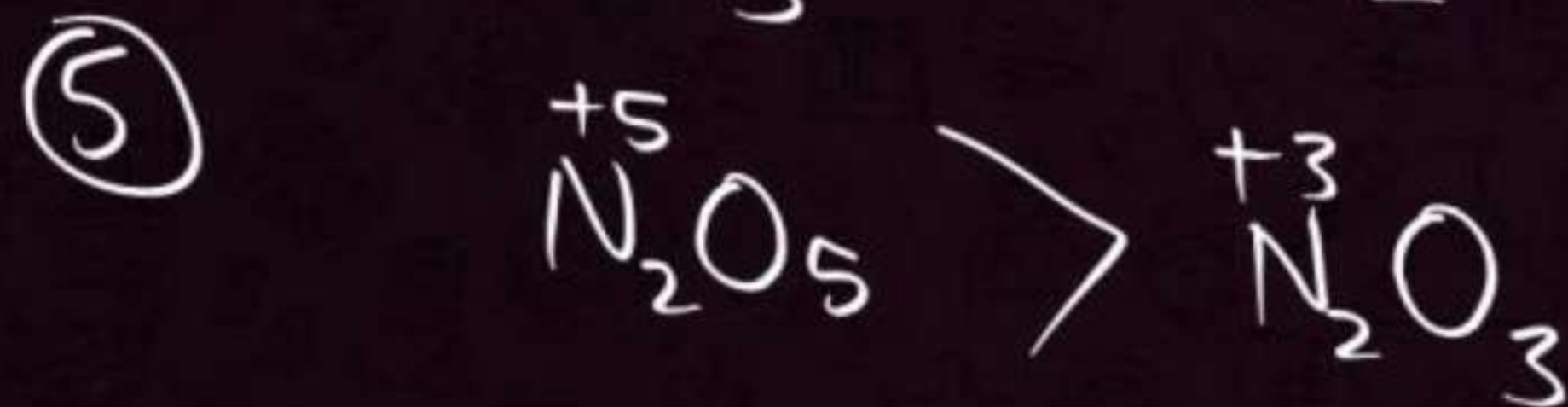
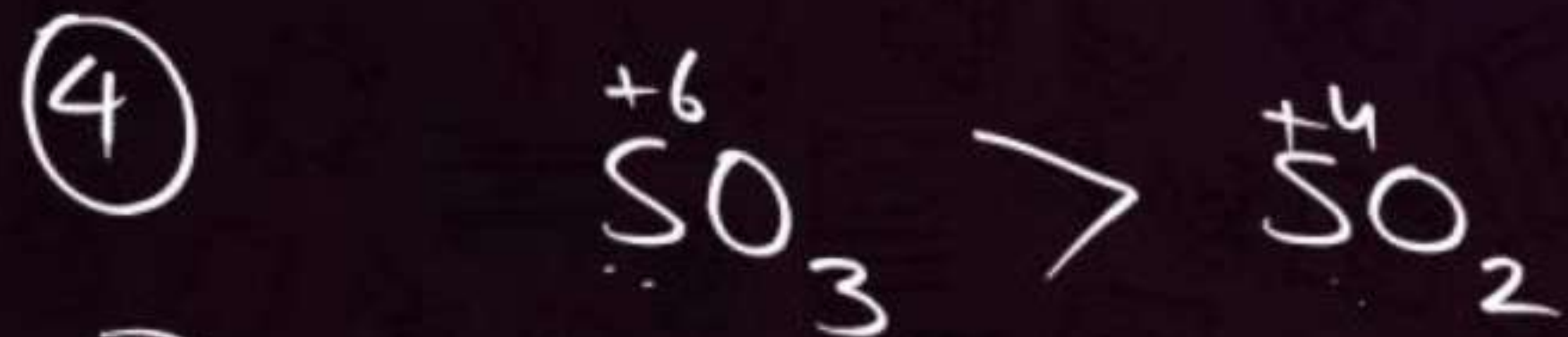
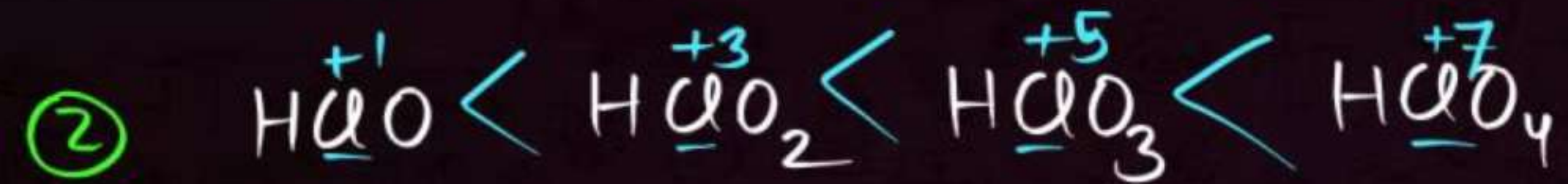
$$\Delta \text{EN} = 0.9$$

$$\text{Cov. char.} = 100 - \% \text{I.C.}$$

⑤ ^{*Imp*} NATURE OF OXIDES/HYDROXIDES :

Non Metal oxide/Hydroxide $\Rightarrow$ Acidic
Metal oxides/hydroxides $\Rightarrow$ Basic



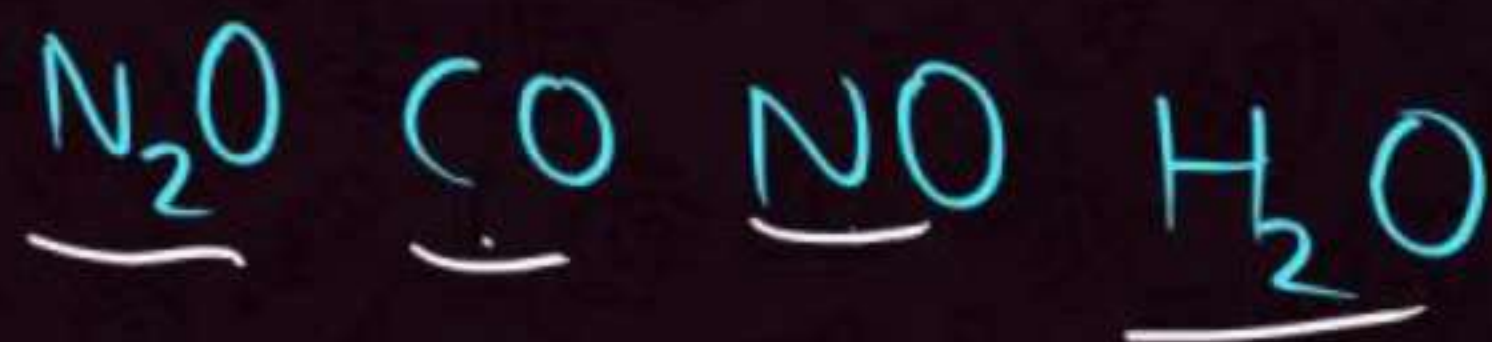


* Amphoteric Oxides $\rightarrow$ (can behave as Acid as well as base both)



Pb	Zn	Be	Al	Ga	Sn	Ce ⁺³	Mn ⁺⁴	Sb ⁺³	As ⁺³	V
PbO	ZnO	BeO	Al ₂ O ₃	Ga ₂ O ₃	SnO	Ce ₂ O ₃	MnO ₂	Sb ₂ O ₃	As ₂ O ₃	
PbO ₂					SnO ₂		Mn ₂ O ₃			
Pb ₃ O ₄							Mn ₃ O ₈			V ₂ O ₅

* Neutral Oxides $\rightarrow$



QUESTION



Which of the following is different from other three oxides?

- ~~A~~ MgO Basic
- B SnO
- C ZnO
- D PbO
- Amph.

Ans. ()

QUESTION



Which one of the following orders present the correct sequence of the increasing basic nature of the given oxides?

- ☒ **A** $\underline{\text{Al}_2\text{O}_3} < \underline{\text{MgO}} < \underline{\text{Na}_2\text{O}} < \underline{\text{K}_2\text{O}}$
- ☐ **B** $\text{MgO} < \text{K}_2\text{O} < \underline{\text{Al}_2\text{O}_3} < \text{Na}_2\text{O} \times$
- ☐ **C** $\text{Na}_2\text{O} < \text{K}_2\text{O} < \text{MgO} < \text{Al}_2\text{O}_3 \times$
- ☐ **D** $\text{K}_2\text{O} < \text{Na}_2\text{O} < \underline{\text{Al}_2\text{O}_3} < \text{MgO} \times$

Ans. ()

QUESTION

Incorrect order of basic character:

- A** Ag_2O > PbO ✓
- B** MgO > ZnO ✓
- C** Fe_2O_3 > Al_2O_3 ✓
- D** ~~Ag_2O > Cu_2O~~ ✓

Ans. ()

GENESIS OF PERIODIC TABLE

DOBEREINER'S TRIADS: → pair of 3 elements

Table 3.1 Dobereiner's Triads

Element	Atomic weight	Element	Atomic weight	Element	Atomic weight
Li	7	Ca	40	Cl	35.5
Na	23	Sr	88	Br	80
K	39	Ba	137	I	127

(K, Rb, Cs)

(Sc, Y, La)

(P, As, Sb)

(S, Se, Te) (H, F, Cl)

$$7 + 39$$

$$= 23.2$$

NEWLAND'S OCTAVE LAW :

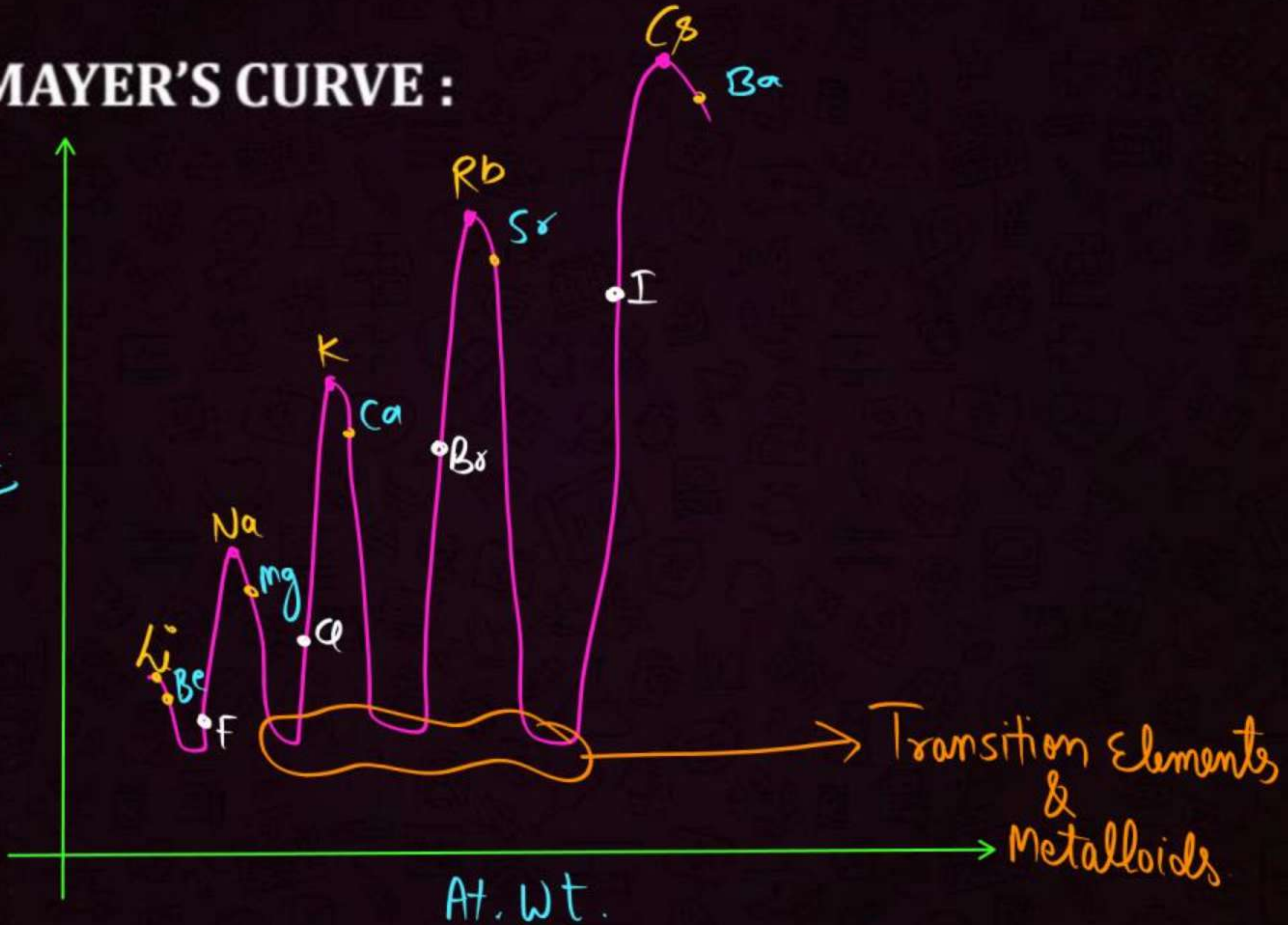
सा रे गा मा पा धा नी सा

Table 3.2 Newlands' Octaves

Element	Li	Be'	B	C	N	O	F
At. wt.	7	9	11	12	14	16	19
Element	Na	Mg	Al	Si	P	S	Cl
At. wt.	23	24	27	29	31	32	35.5
Element	K	Ca					
At. wt.	39	40					

LOTHER MAYER'S CURVE :

At. Vol





MENDELEEV'S PERIODIC TABLE :

→ 63 elements

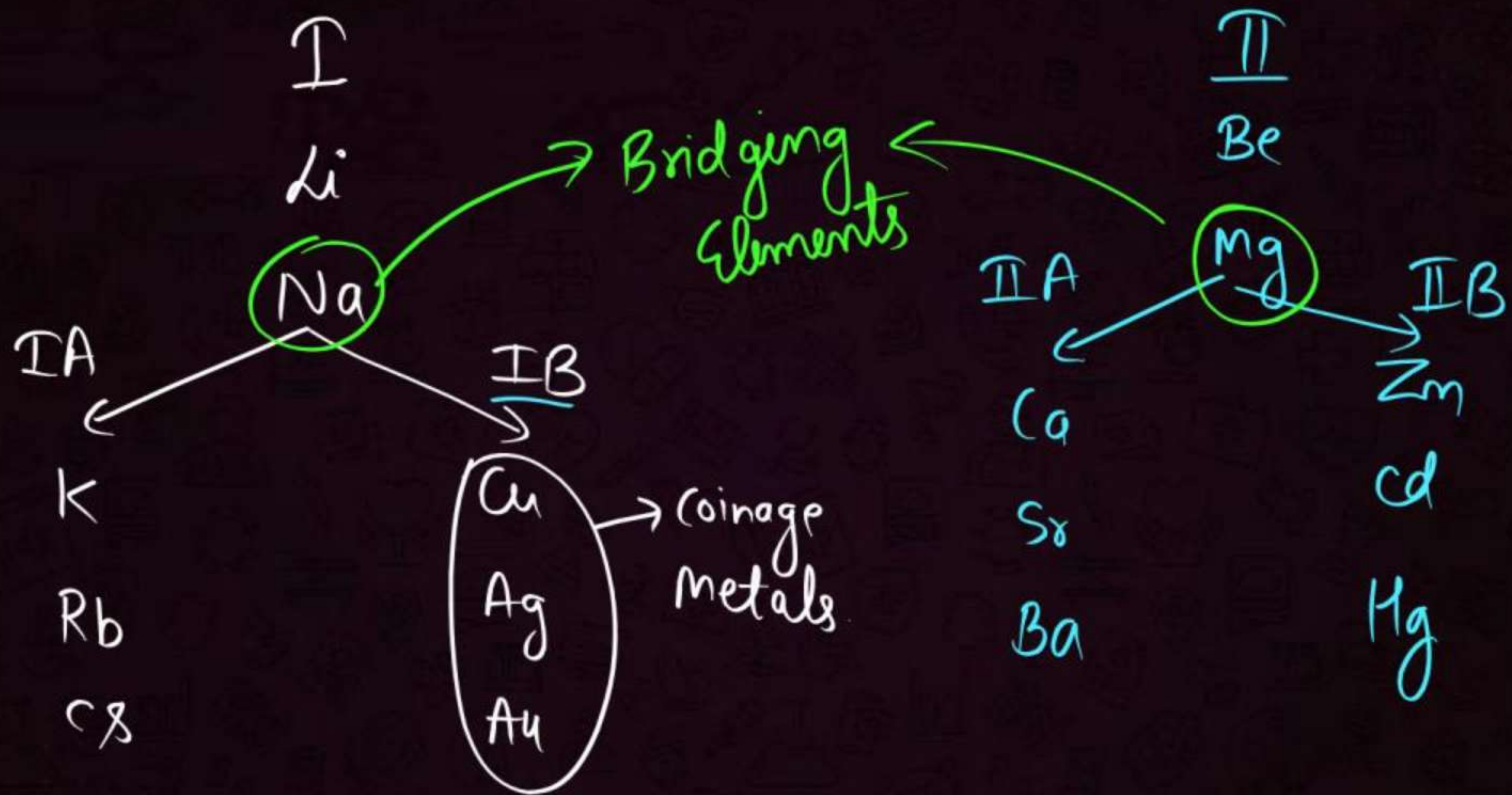
→ Based on Atomic Wt.

→ 8 vertical columns (Groups) & 12 horizontal rows (series)

7 period

→ 9th group was added after discovery of Noble gases & named as group zero.

→ Group 1 to 7 were divided into 2 subgroups.



VIII

Fe	Co	Ni
Ru	Rh	Pd
Os	Ir	Pt

गोला [Ga] → Eka Aluminium Al

Sc → Eka Boron B

जस्ता Zn → Eka Silicon Si

तम्र Cu → Eka Manganese Mn

Correction in At. Wt. → U Be Im Pt Au

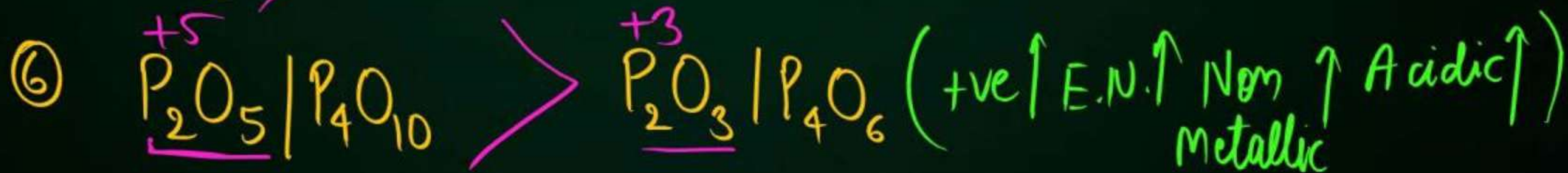
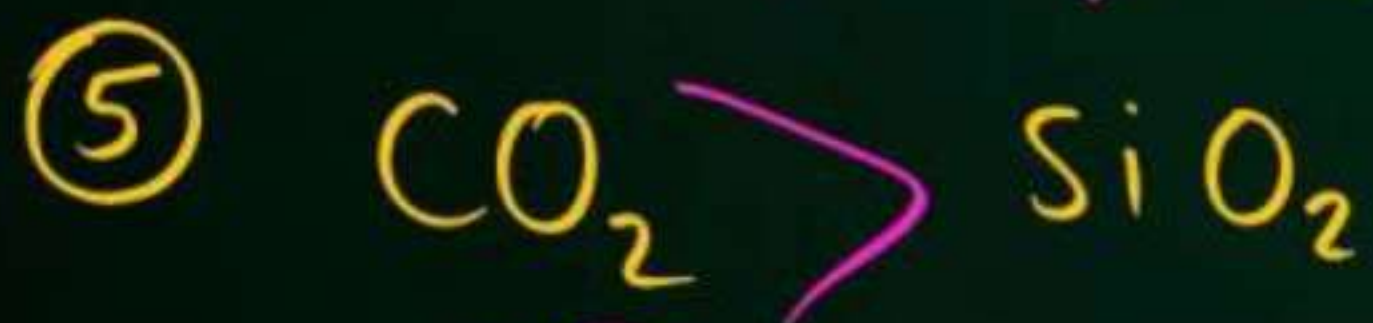
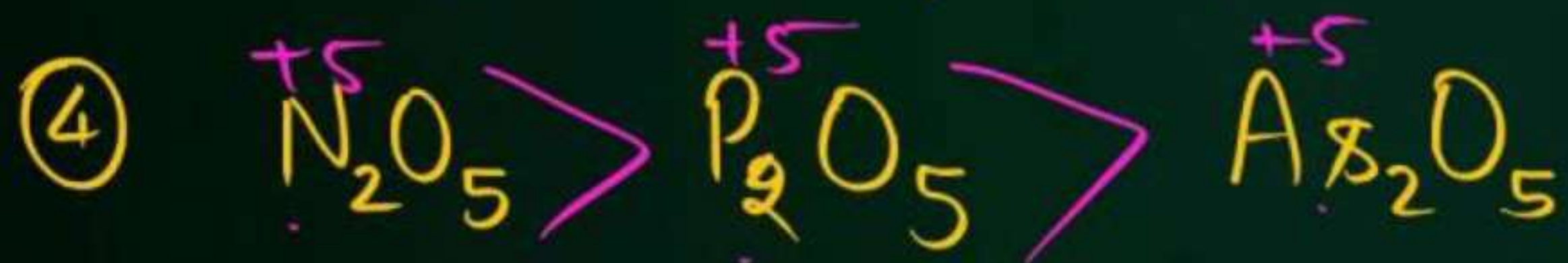
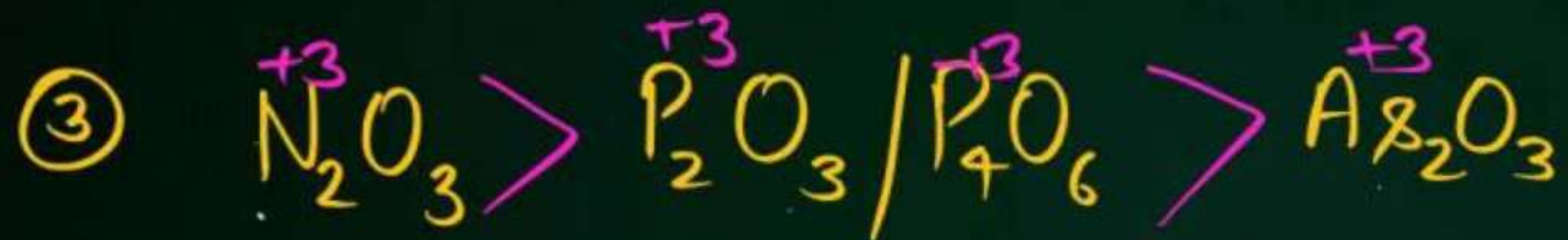
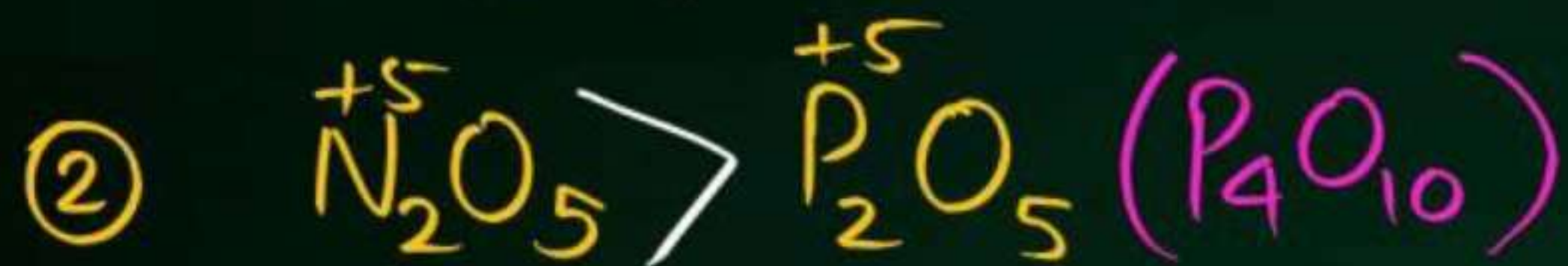
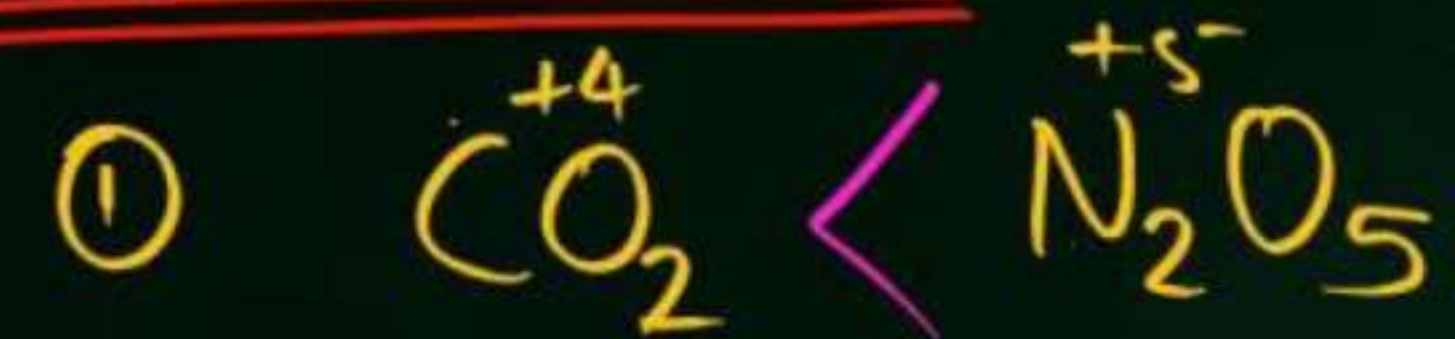


* Drawbacks



- ① Position of H
- ② Position of isotopes
- ③ Anomalous Pairs : (Ar, K) (Te, I)
39.9 39.1

Acidic Nature :-



E.N. $\uparrow$
Non metallic Nature $\uparrow$
Acidic $\uparrow$

Basic Nature :->

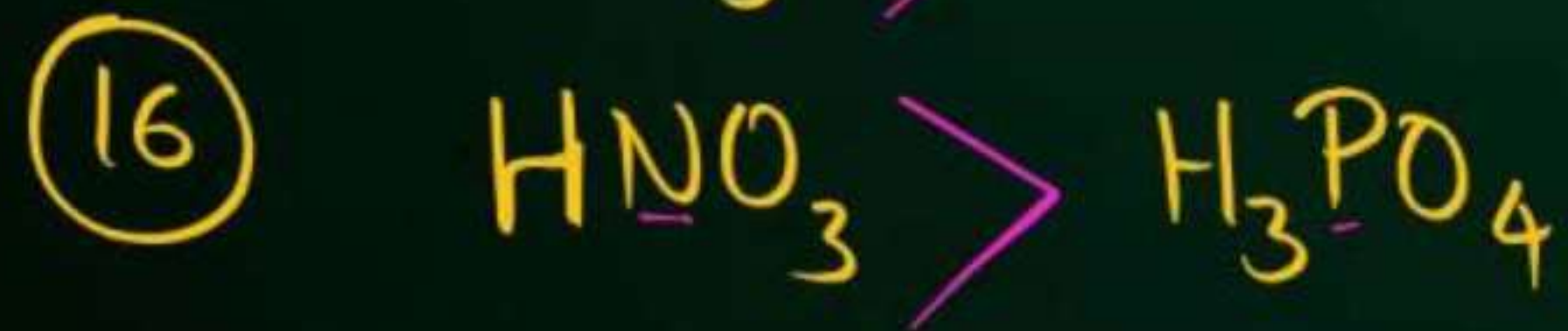
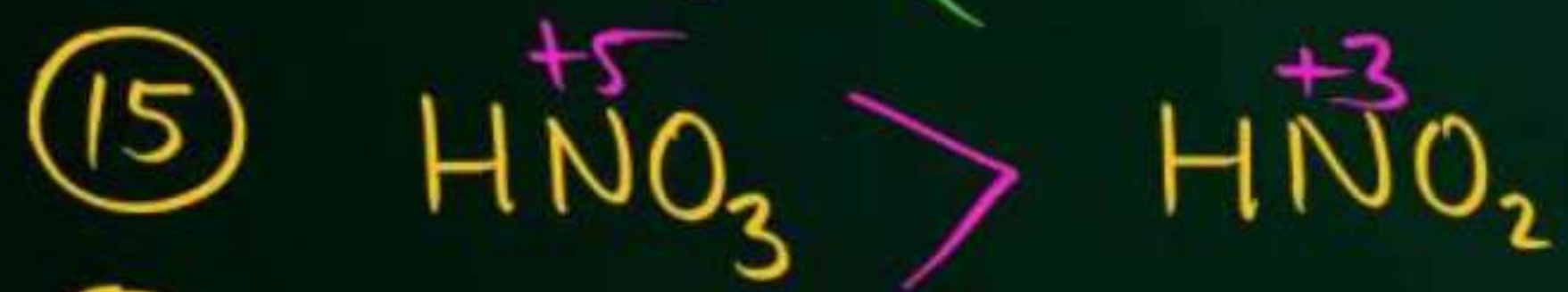
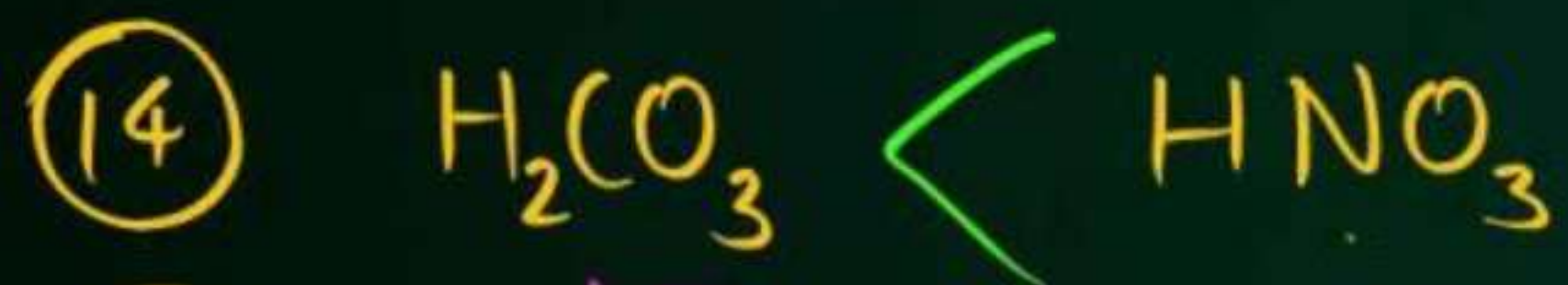


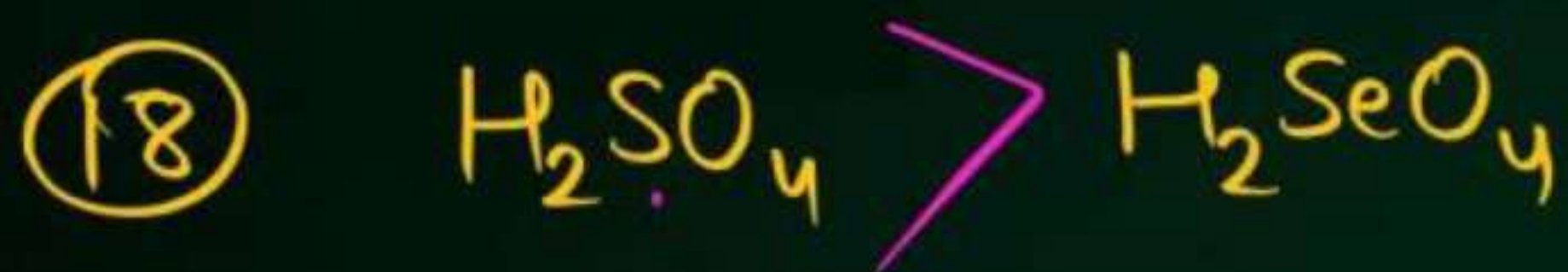
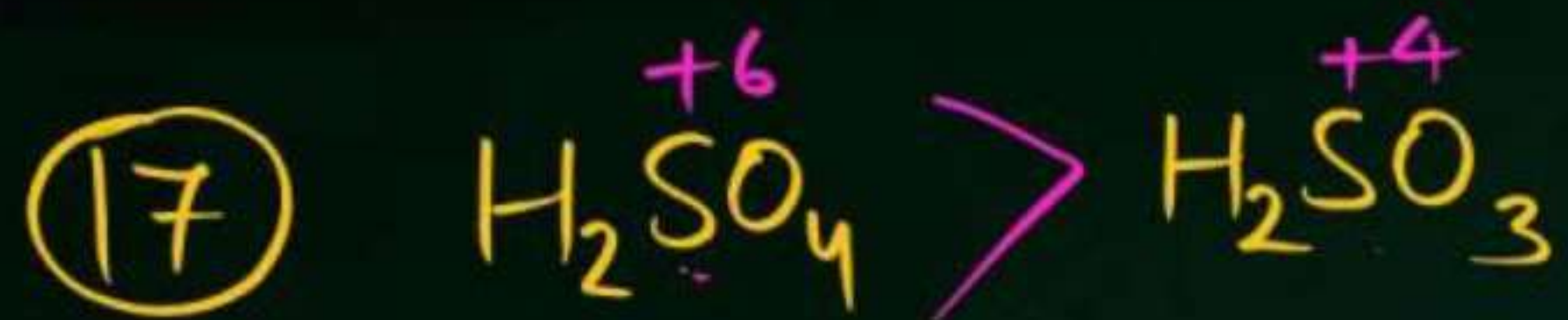
$\xrightarrow{\text{E.N.} \uparrow \text{Non Metallic} \uparrow \text{Metallic} \downarrow \text{Basic} \downarrow}$



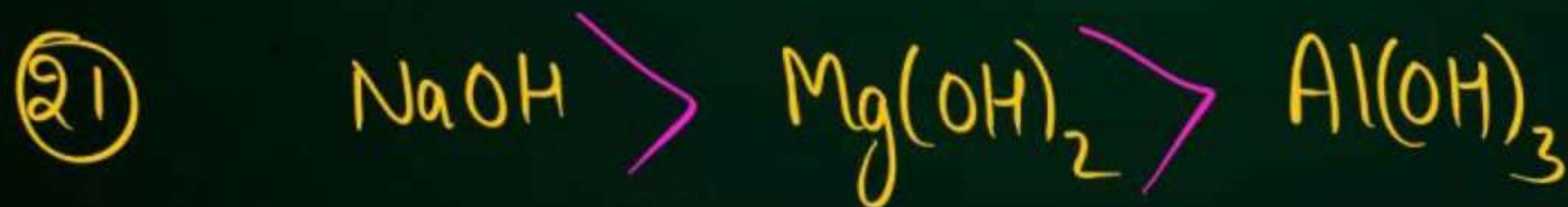
E.N. $\downarrow$
Metallic $\uparrow$
Basic $\uparrow$

Acidic Nature $\rightarrow$ (oxy Acids)





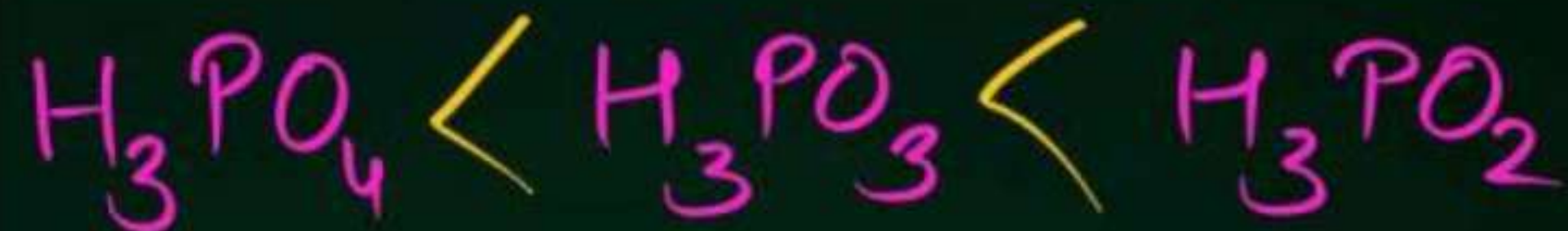
Basic Nature : $\rightarrow$



* Remember *



Acidic Nature :



Reducing Nature :

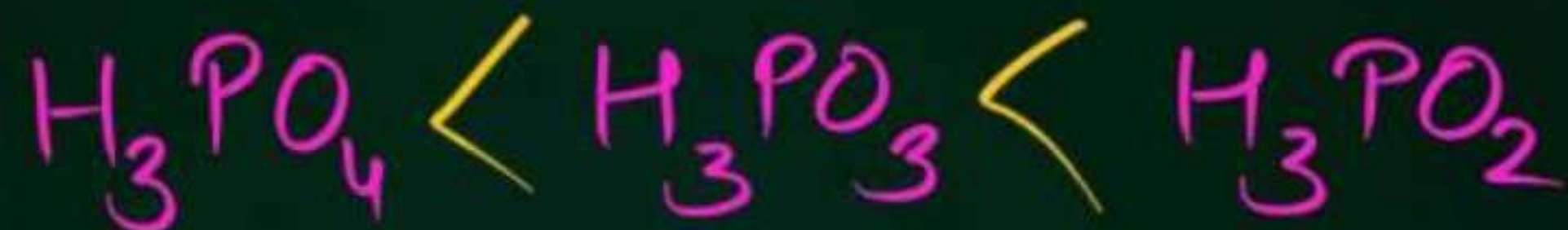


Table 2.6 Electronic Configurations of the Elements

Element Z	1s	2s	2p	3s	3p	3d	4s	4p	4d	4f	5s	5p	5d	5f	6s	6p	6d	7s
H 1	1																	
He 2	2																	
Li 3	2	1																
Be 4	2	2																
B 5	2	2	1															
C 6	2	2	2															
N 7	2	2	3															
O 8	2	2	4															
F 9	2	2	5															
Ne 10	2	2	6															
Na 11	2	2	6	1														
Mg 12	2	2	6	2														
Al 13	2	2	6	2	1													
Si 14	2	2	6	2	2													
P 15	2	2	6	2	3													
S 16	2	2	6	2	4													
Cl 17	2	2	6	2	5													
Ar 18	2	2	6	2	6													
K 19	2	2	6	2	6		1											
Ca 20	2	2	6	2	6		2											
Sc 21	2	2	6	2	6	1	2											
Ti 22	2	2	6	2	6	2	2											
V 23	2	2	6	2	6	3	2											
Cr* 24	2	2	6	2	6	5	1											
Mn 25	2	2	6	2	6	5	2											
Fe 26	2	2	6	2	6	6	2											
Co 27	2	2	6	2	6	7	2											
Ni 28	2	2	6	2	6	8	2											
Cu* 29	2	2	6	2	6	10	1											
Zn 30	2	2	6	2	6	10	2											

Ga	31	2	2	6	2	6	10	2	1			
Ge	32	2	2	6	2	6	10	2	2			
As	33	2	2	6	2	6	10	2	3			
Se	34	2	2	6	2	6	10	2	4			
Br	35	2	2	6	2	6	10	2	5			
Kr	36	2	2	6	2	6	10	2	6			
Rb	37	2	2	6	2	6	10	2	6	1		
Sr	38	2	2	6	2	6	10	2	6	2		
Y	39	2	2	6	2	6	10	2	6	1		
Zr	40	2	2	6	2	6	10	2	6	2		
Nb*	41	2	2	6	2	6	10	2	6	4		
Mo*	42	2	2	6	2	6	10	2	6	5		
Tc	43	2	2	6	2	6	10	2	6	5		
Ru*	44	2	2	6	2	6	10	2	6	7		
Rh*	45	2	2	6	2	6	10	2	6	8		
Pd*	46	2	2	6	2	6	10	2	6	10		
Ag*	47	2	2	6	2	6	10	2	6	10	1	
Cd	48	2	2	6	2	6	10	2	6	10	2	
In	49	2	2	6	2	6	10	2	6	10	2	1
Sn	50	2	2	6	2	6	10	2	6	10	2	2
Sb	51	2	2	6	2	6	10	2	6	10	2	3
Te	52	2	2	6	2	6	10	2	6	10	2	4
I	53	2	2	6	2	6	10	2	6	10	2	5
Xe	54	2	2	6	2	6	10	2	6	10	2	6

* Elements with exceptional electronic configurations

Element Z		1s	2s	2p	3s	3p	3d	4s	4p	4d	4f	5s	5p	5d	5f	6s	6p	6d	7s
Cs	55	2	2	6	2	6	10	2	6	10		2	6			1			
Ba	56	2	2	6	2	6	10	2	6	10		2	6			2			
La*	57	2	2	6	2	6	10	2	6	10		2	6	1		2			
Ce*	58	2	2	6	2	6	10	2	6	10	2	2	6			2			
Pr	59	2	2	6	2	6	10	2	6	10	3	2	6			2			
Nd	60	2	2	6	2	6	10	2	6	10	4	2	6			2			
Pm	61	2	2	6	2	6	10	2	6	10	5	2	6			2			
Sm	62	2	2	6	2	6	10	2	6	10	6	2	6			2			
Eu	63	2	2	6	2	6	10	2	6	10	7	2	6			2			
Gd*	64	2	2	6	2	6	10	2	6	10	7	2	6	1		2			
Tb	65	2	2	6	2	6	10	2	6	10	9	2	6			2			
Dy	66	2	2	6	2	6	10	2	6	10	10	2	6			2			
Ho	67	2	2	6	2	6	10	2	6	10	11	2	6			2			
Er	68	2	2	6	2	6	10	2	6	10	12	2	6			2			
Tm	69	2	2	6	2	6	10	2	6	10	13	2	6			2			
Yb	70	2	2	6	2	6	10	2	6	10	14	2	6			2			
Lu	71	2	2	6	2	6	10	2	6	10	14	2	6	1		2			
Hf	72	2	2	6	2	6	10	2	6	10	14	2	6	2		2			
Ta	73	2	2	6	2	6	10	2	6	10	14	2	6	3		2			
W	74	2	2	6	2	6	10	2	6	10	14	2	6	4		2			
Re	75	2	2	6	2	6	10	2	6	10	14	2	6	5		2			
Os	76	2	2	6	2	6	10	2	6	10	14	2	6	6		2			
Ir	77	2	2	6	2	6	10	2	6	10	14	2	6	7		2			
Pt*	78	2	2	6	2	6	10	2	6	10	14	2	6	9		1			
Au*	79	2	2	6	2	6	10	2	6	10	14	2	6	10		1			
Hg	80	2	2	6	2	6	10	2	6	10	14	2	6	10		2			
Tl	81	2	2	6	2	6	10	2	6	10	14	2	6	10		2	1		
Pb	82	2	2	6	2	6	10	2	6	10	14	2	6	10		2	2		
Bi	83	2	2	6	2	6	10	2	6	10	14	2	6	10		2	3		
Po	84	2	2	6	2	6	10	2	6	10	14	2	6	10		2	4		
At	85	2	2	6	2	6	10	2	6	10	14	2	6	10		2	5		
Rn	86	2	2	6	2	6	10	2	6	10	14	2	6	10		2	6		





Fr	87	2	2	6	2	6	10	2	6	10	14	2	6	10	2	6	1
Ra	88	2	2	6	2	6	10	2	6	10	14	2	6	10	2	6	2
Ac	89	2	2	6	2	6	10	2	6	10	14	2	6	10	2	6	2
Th	90	2	2	6	2	6	10	2	6	10	14	2	6	10	2	6	2
Pa	91	2	2	6	2	6	10	2	6	10	14	2	6	10	2	6	2
U	92	2	2	6	2	6	10	2	6	10	14	2	6	10	2	6	2
Np	93	2	2	6	2	6	10	2	6	10	14	2	6	10	2	6	2
Pu	94	2	2	6	2	6	10	2	6	10	14	2	6	10	2	6	2
Am	95	2	2	6	2	6	10	2	6	10	14	2	6	10	2	6	2
Cm	96	2	2	6	2	6	10	2	6	10	14	2	6	10	2	6	2
Bk	97	2	2	6	2	6	10	2	6	10	14	2	6	10	2	6	2
Cf	98	2	2	6	2	6	10	2	6	10	14	2	6	10	2	6	2
Es	99	2	2	6	2	6	10	2	6	10	14	2	6	10	2	6	2
Fm	100	2	2	6	2	6	10	2	6	10	14	2	6	10	2	6	2
Md	101	2	2	6	2	6	10	2	6	10	14	2	6	10	2	6	2
No	102	2	2	6	2	6	10	2	6	10	14	2	6	10	2	6	2
Lr	103	2	2	6	2	6	10	2	6	10	14	2	6	10	2	6	2
Rf	104	2	2	6	2	6	10	2	6	10	14	2	6	10	2	6	2
Db	105	2	2	6	2	6	10	2	6	10	14	2	6	10	2	6	2
Sg	106	2	2	6	2	6	10	2	6	10	14	2	6	10	2	6	2
Bh	107	2	2	6	2	6	10	2	6	10	14	2	6	10	2	6	2
Hs	108	2	2	6	2	6	10	2	6	10	14	2	6	10	2	6	2
Mt	109	2	2	6	2	6	10	2	6	10	14	2	6	10	2	6	2
Ds	110	2	2	6	2	6	10	2	6	10	14	2	6	10	2	6	2
Rg**	111	2	2	6	2	6	10	2	6	10	14	2	6	10	2	6	1

** Elements with atomic number 112 and above have been reported but not yet fully authenticated and named.



Homework



Practice Sheet

———— **FOR NOTES & DPP CHECK DESCRIPTION** ————

THANK YOU

