

# Periodic properties

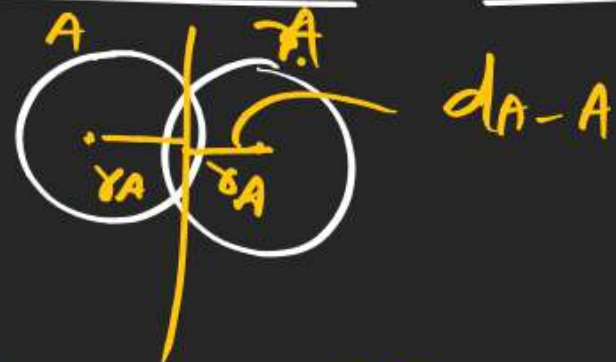
- ① Atomic radii → We can't measure  
because atom does not have  
certain boundary

- ② Covalent radii

Homonuclear diatomic molecule ( $H_2$ ,  $C_2$ ,  $Br_2$ )

ex find the  $r_c$ ,  
if  $d_{Cl-Cl}$  is  $1.98 \text{ \AA}$

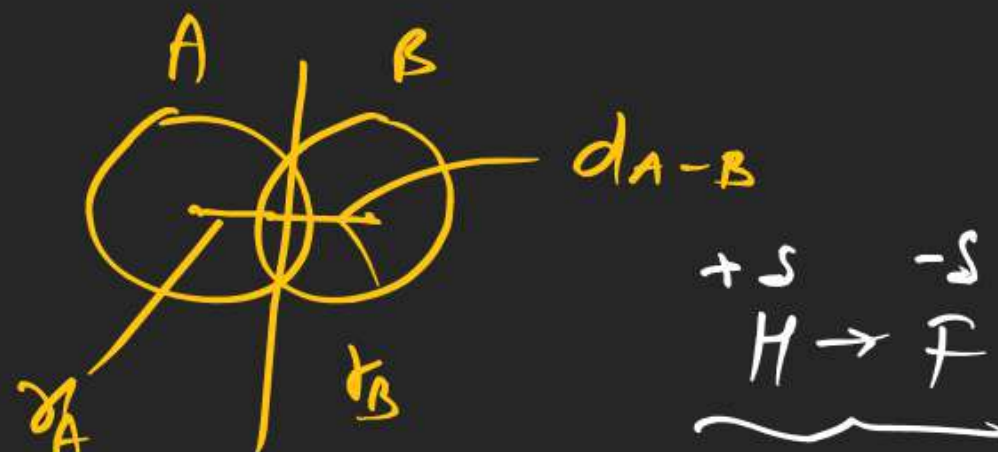
$$r_{Cl} = \frac{1.98}{2}$$



$$d_{A-A} = r_A + r_A$$

$$2r_A = d_{AA}$$

$$r_A = \frac{d_{A-A}}{2}$$

Heteronuclear diatomic molecule

Shomaker  
and Stevensen

$$d_{A-B} = r_A + r_B - 0.09 |\Delta\chi| \text{ \AA}$$

$\Delta\chi = \epsilon$  in diff

$r_A =$  radius of A in A'

$r_B =$  radius of B in A'

$$d_{A-B} = r_A + r_B - 9 |\Delta\chi| \text{ in pm}$$

S.B.C.R

D.B.C.R

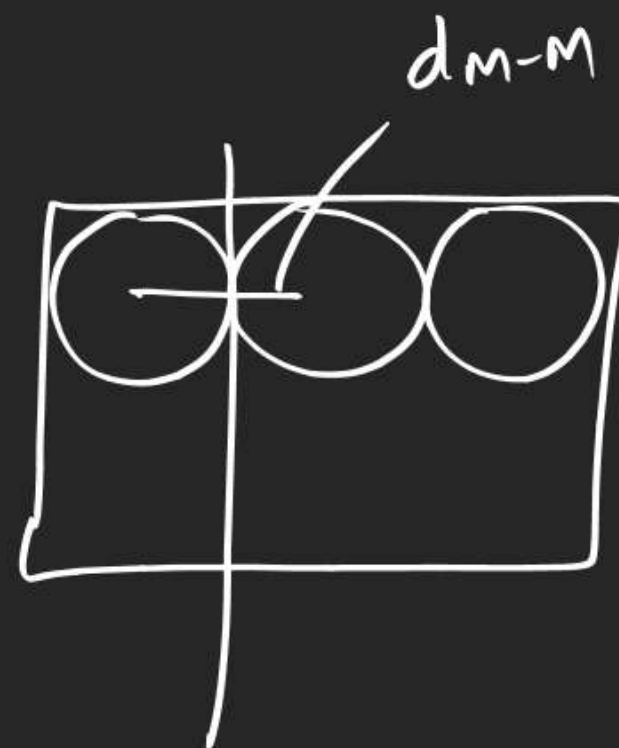
T.B.C.R

$C-C > C=C > C\equiv C$

Use Gm not measure

X     $H-H$      $O=O$

② Metallic



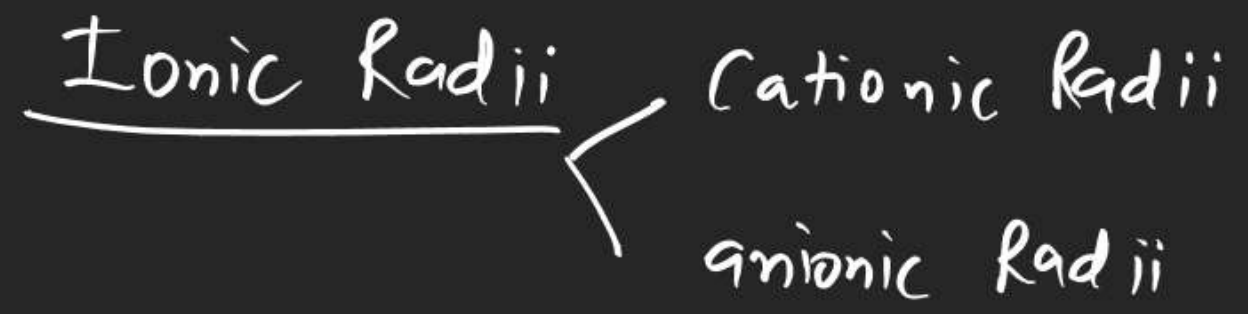
$$d_{M-M} = r_M + r_M$$

$$r_M = \frac{d_{M-M}}{2}$$

Q =  $d_{Cu-Cu}$  is 129 pm

then  $r_{Cu}$

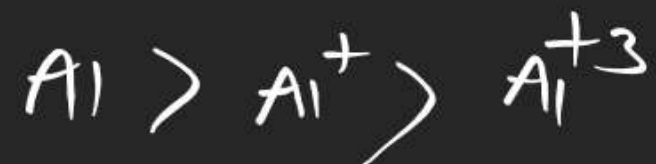
$$\frac{129}{2} = 64.5 \text{ pm}$$



Cationic Radii

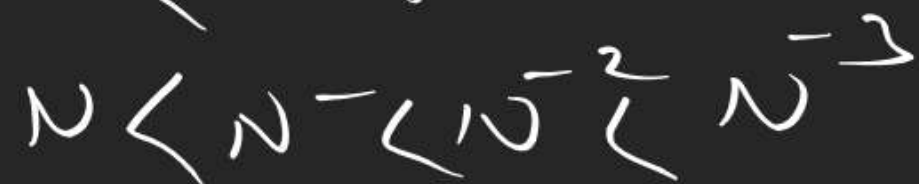


↑ positive charge    ↓ size



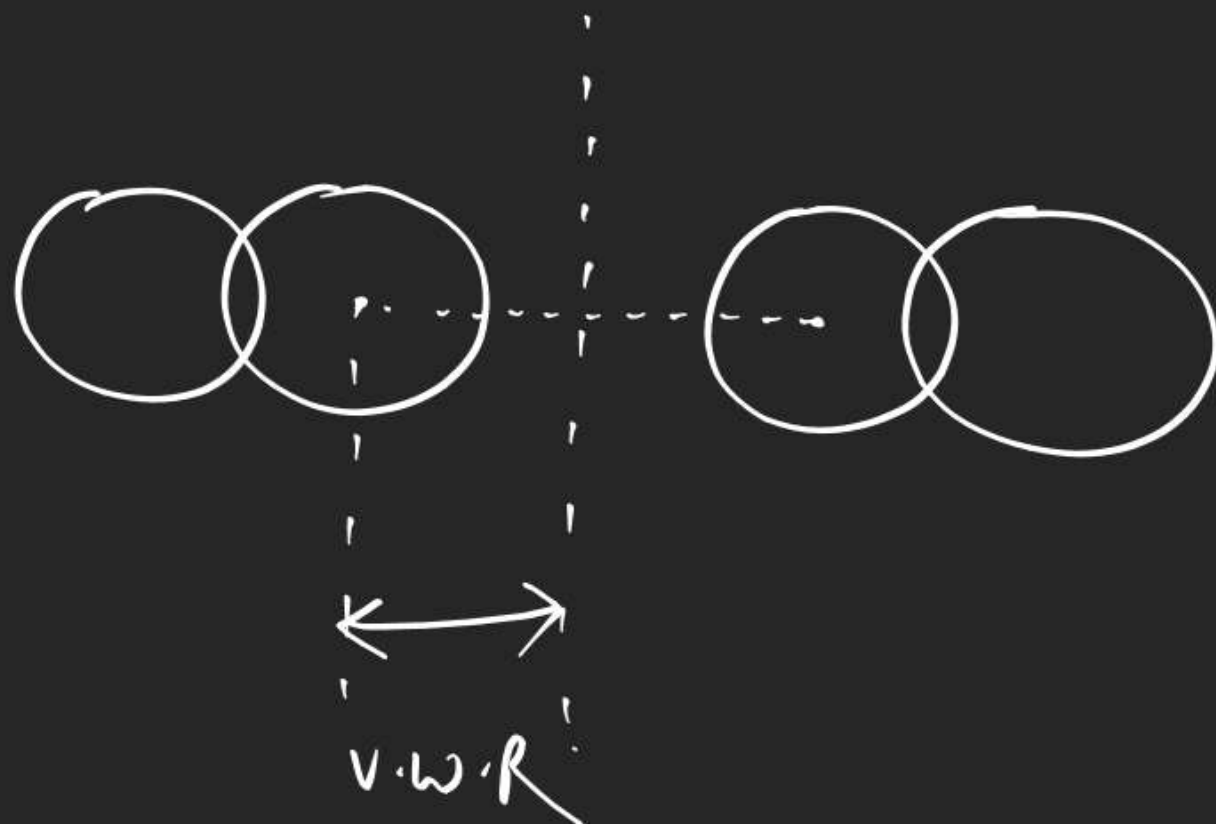
Anionic

-ive charge ↑ size ↑





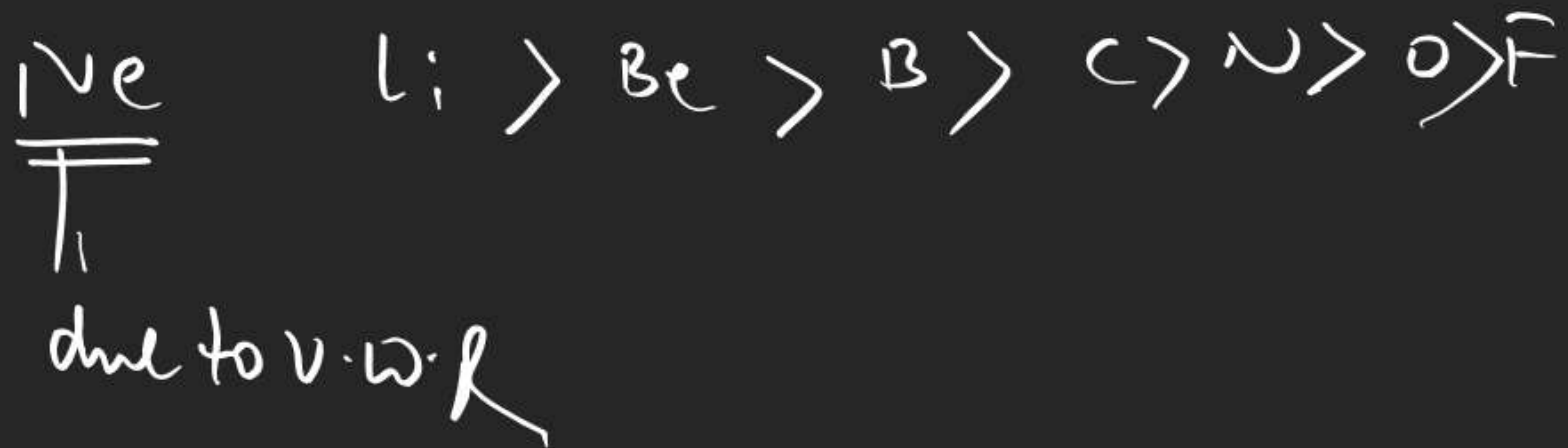
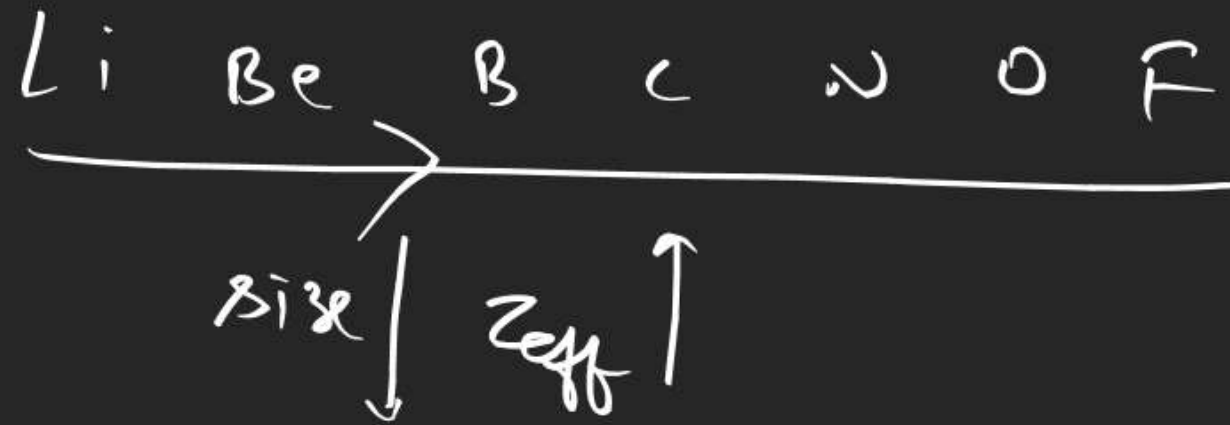
V.W.R



$$V.W.R > C.R$$

Noble gas and  $C_{70}$  etc

along the period





down the group

Li	Bc
Na	Mg
K	Ca
Rb	Sr
Cs	Ba

size ↑ because shell ↑

P-Block

$\left. \begin{array}{l} 143 \text{ pm B} \\ 135 \text{ pm Al} \end{array} \right\}$   
 $\left. \begin{array}{l} 135 \text{ pm Ga} \\ \text{In} \\ \text{Te} \end{array} \right\}$

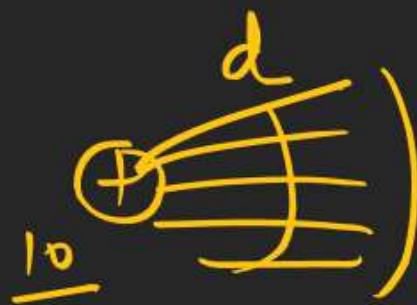
$\overset{167 \text{ pm}}{\text{B}} < \overset{170 \text{ pm}}{\text{Ga}} < \text{Al} < \text{In} < \text{Te}$

due to poor s-f of  
3d sub shell

Scandide contraction

$$Al = 1s^2 \cancel{2s^2} 2p^6 3s^2 3p^1$$

$$Ga = 1s^2 \cancel{2s^2} 2p^6 3s^2 3p^6 \underline{4s^2} \underline{3d^{10}} 4p^1$$



S.E

s > p > d > f

$$\begin{matrix} C \\ \nearrow Si \\ \searrow Ge \end{matrix} = \underline{up^2}$$

$Sn$   
 $Pb$

order of size

$$C < Si < Ge < Sn < Pb$$

d-Block

Sc	Ti	V	Cr	Mn	Fe	Co	Ni	Cu	Zn
$Z_{eff} > S.E$ $r \downarrow$					$Z_{eff} \approx S.E$ $r \approx \text{Const.}$			$Z_{eff} < S.E$ $r \uparrow$	

down the group

$Ti < Zr \approx Hf$   
150 pm 159 pm

3d   IT-S	Sc	Ti	.....	Cu	Zn
4d   IT-S	Y	Zr	.....	Ag	Cd
5d   III T-S	57 La	72 Hf	.....	Au	Hg
	89 Ac	104 Rf			



$Sc < Y < La < Ac$

89

Ac

La

104  
Rf

58  
Ce

71  
La

4f series / Lanthanide

90  
Th

103  
La

5f series / Actinide

3d series  
element

4d series  
element

$\approx$  5d series  
element

ans order of size



↑  
+ive charge | size |



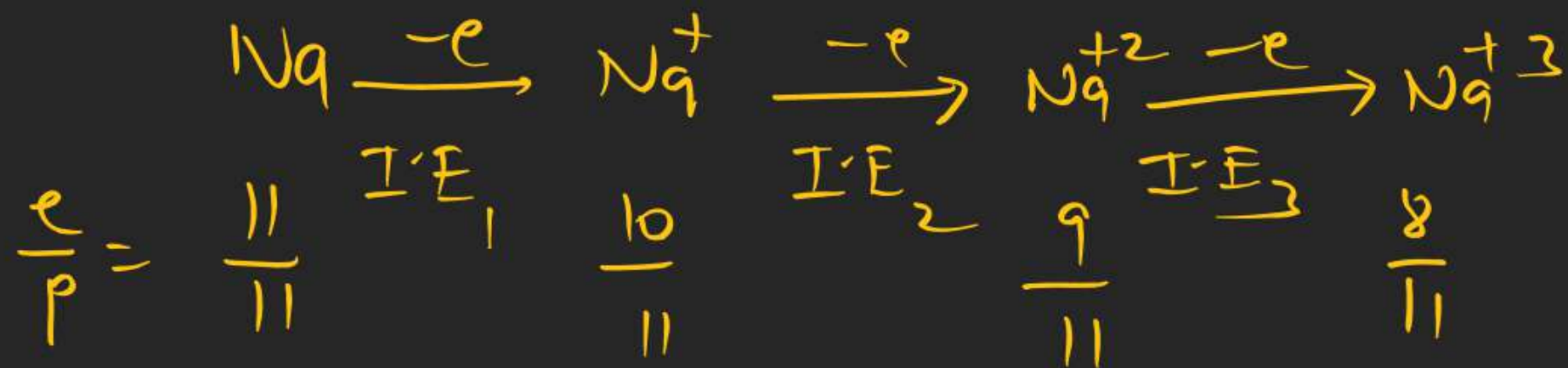




# I.E



## Successive



$$Z_{\text{eff of III}} > Z_{\text{eff of II}} > Z_{\text{eff of I.E}}$$

\*  $\boxed{\text{I.E}_3 > \text{I.E}_2 > \text{I.E}_1}$

$$\overset{\text{always}}{\left(\underline{I \cdot E}\right)_n} > \left(I \cdot E\right)_{n-1}$$

Unit

endothermic

$$1\text{eV/atom} = 96.4 \text{ kJ/mole}$$

$$1\text{eV/atom} = 23.1 \text{ kcal/mole}$$

fac.

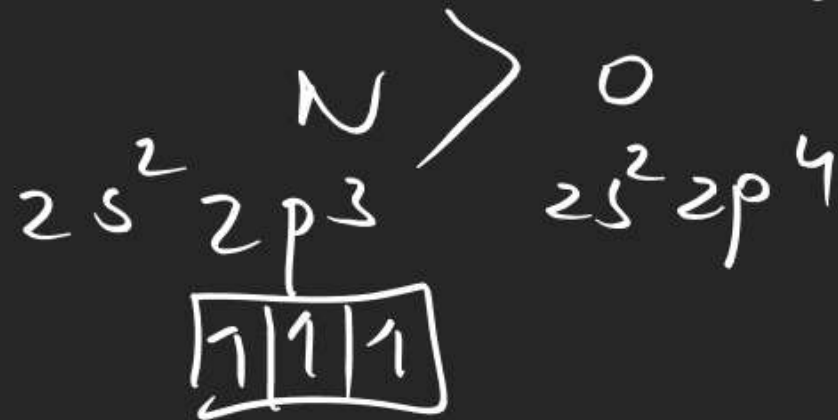
$$① \quad z \uparrow \quad \pm \cdot E \uparrow$$

$$② \quad z_{eff} \uparrow \quad \pm \cdot E \uparrow$$

$$③ \quad n \uparrow \quad \pm \cdot E \downarrow$$

$$④ \quad \sigma (s \cdot E) \uparrow \quad \pm \cdot E \downarrow$$

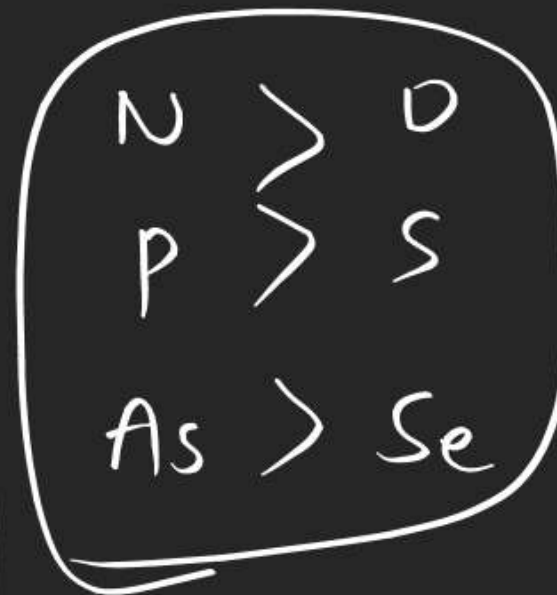
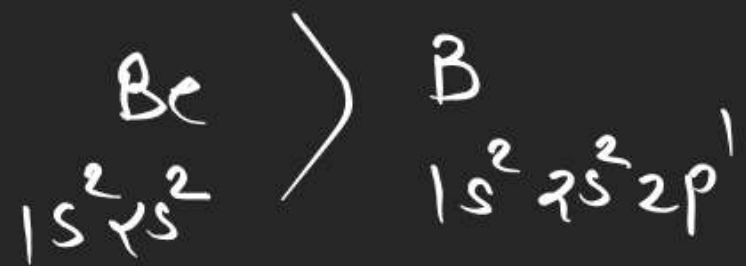
⑤ Half filled and fully filled



## Penetration effect

(Closeness toward nucleus)

$$s > p > d > f$$



along the period



$z_{\text{eff}} \uparrow$  size  $\downarrow$   $I-E \uparrow$



down the group

