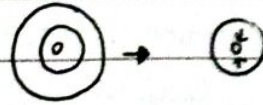


## ATOMIC & IONIC RADII -

→ Ionic radii < Atomic radii

↳ lose outermost shell of  $e^-$  ∴ one shell less

↳  $e^-$  more strongly held



→ Down group, atomic & ionic radii ↑

↳ no. of  $e^-$  shells ↑ ∴ distance of outer  $e^-$  from nucleus ↑

↳ no. of  $e^-$  shells ↑ ∴ shielding ↑

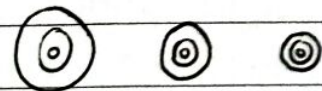
↳ nuclear attraction ↓ ∴



→ Across period, atomic & ionic radii ↓

↳ same no. of shells ∴ shielding effect same.

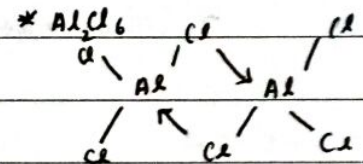
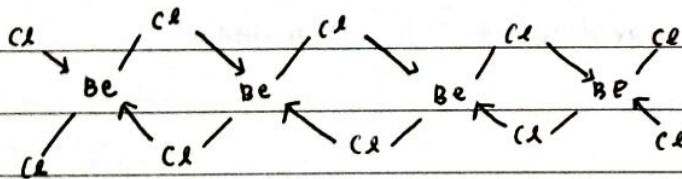
↳ but proton no. ↑ ∴ nuclear attraction ↑



## REACTION W/ CHLORINE -

↳ only Be reaction differs

↳ forms linear molecule w/ dative covalent bond (not dimer)



\* Be only has 2 outer  $e^-$

## REACTIONS W/ WATER (OXIDES) -

↳  $BeO$ ,  $MgO$  insoluble

↳ basicity ↑ down the group.

## TEST FOR SULPHATES -

↳ acidify solution w/  $HNO_3$

↳ add  $BaCl_2$

↳ (+ve) white ppt  $BaSO_4$

### → THERMAL STABILITY -

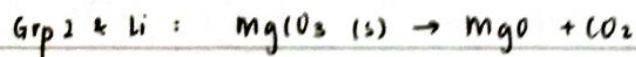
- ↳ polarisation  $\therefore$   $e^-$  cloud of anion distorted  $\therefore$  less stable on heating.
- ↳ oxides stable  $\therefore$  O small & hard to polarise.
- \* Li small  $\therefore$  charge density  $\uparrow$   $\therefore$  behave like Grp 2.

### → NITRATES \* $NO_3^-$

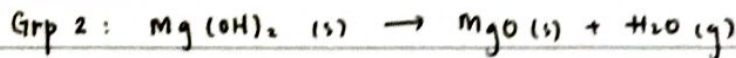


### → Carbonates

Grp 1: stable



### → Hydroxides



Grp 1: stable

- ↳ going down the group, thermal stability  $\uparrow$
- ↳ cation larger  $\therefore$  charge density  $\downarrow$   $\therefore$  less polarising.

### → USES -

