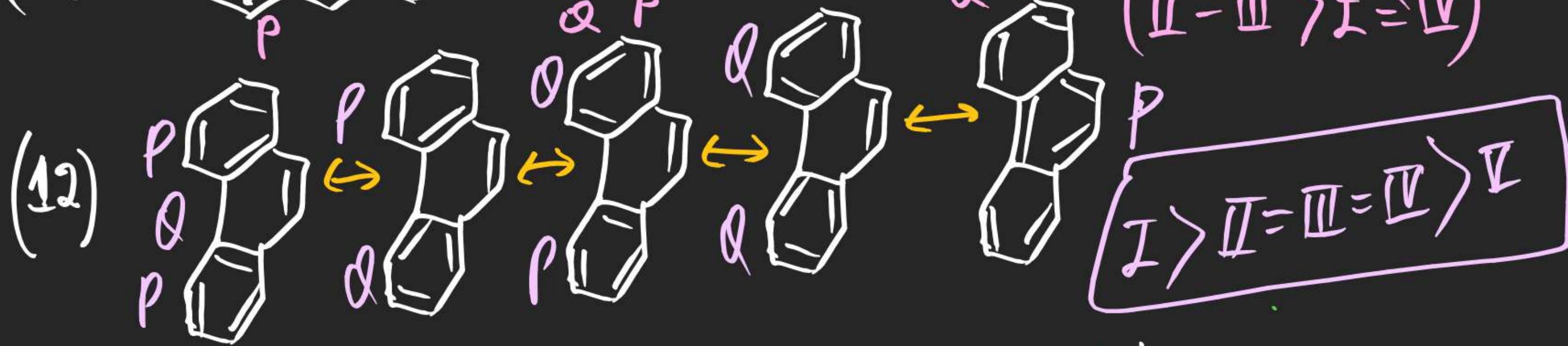
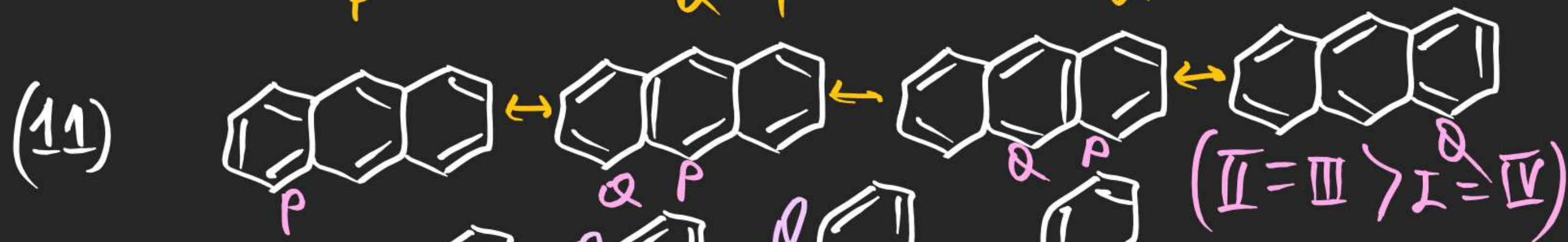
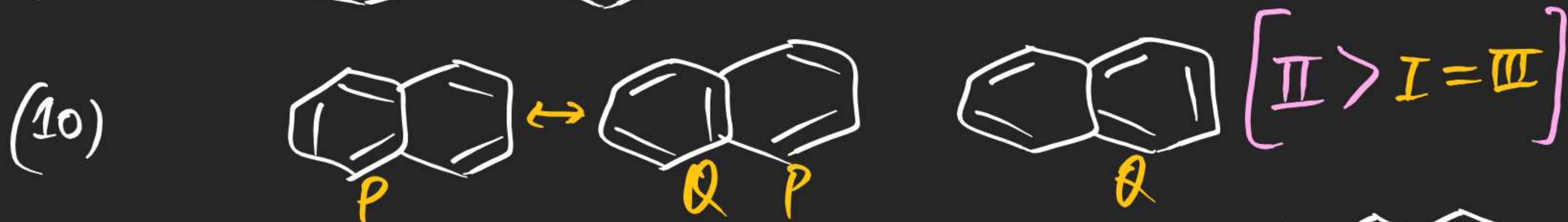


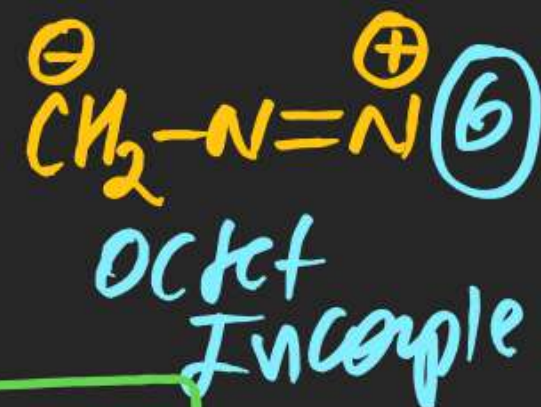
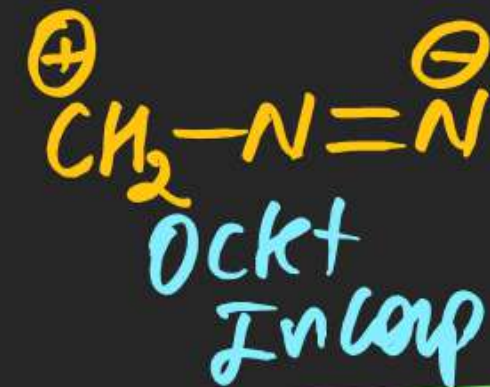
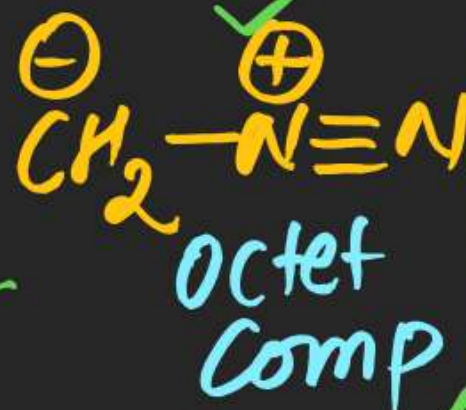
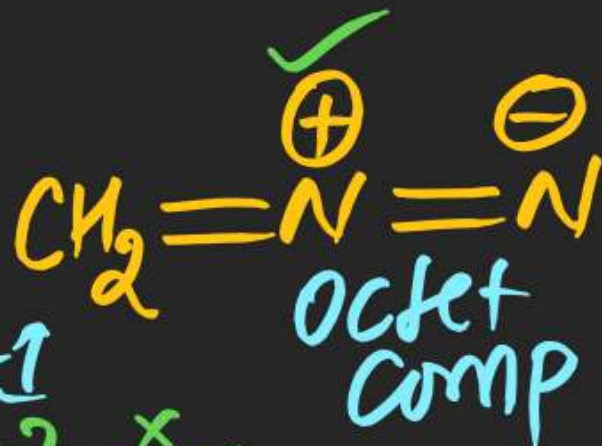
⑤ RS having higher no. of Benzenoid Segment is more stable.



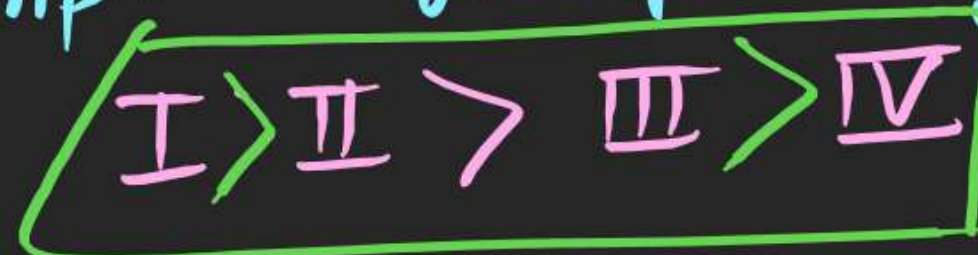
IIT Ad
(13)

Rule-1

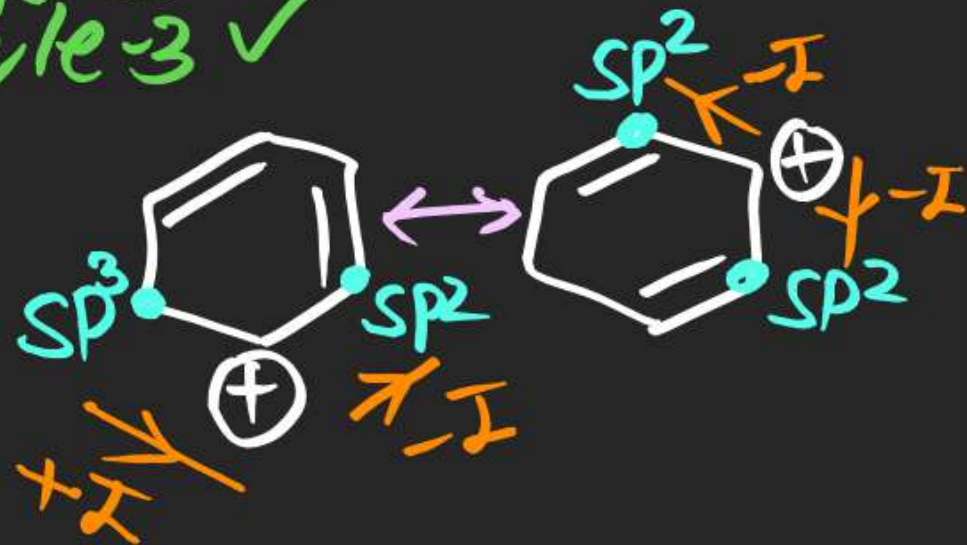
Rule-2 \times
Rule-3 \checkmark



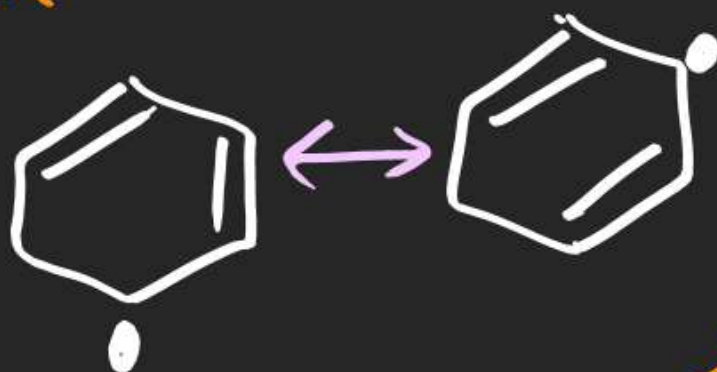
Rule-2 \times
Rule-3 \checkmark



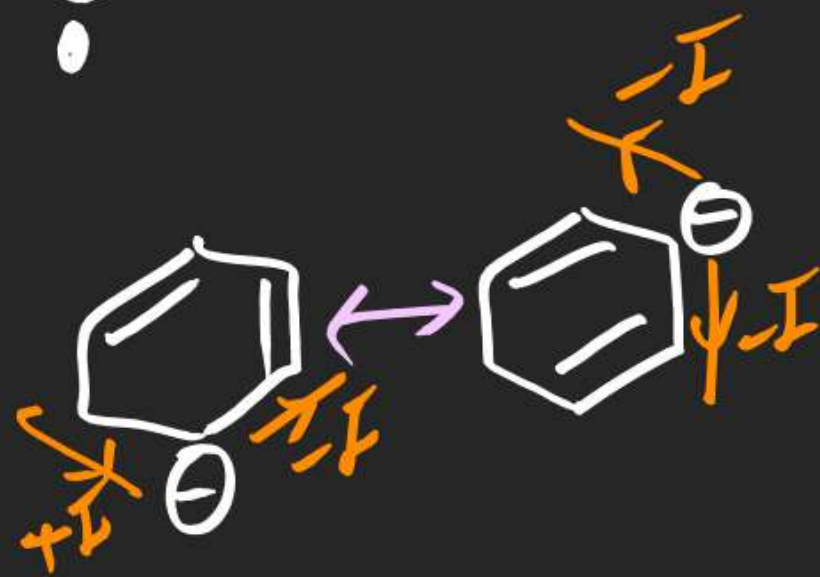
(14)



(15)



(16)



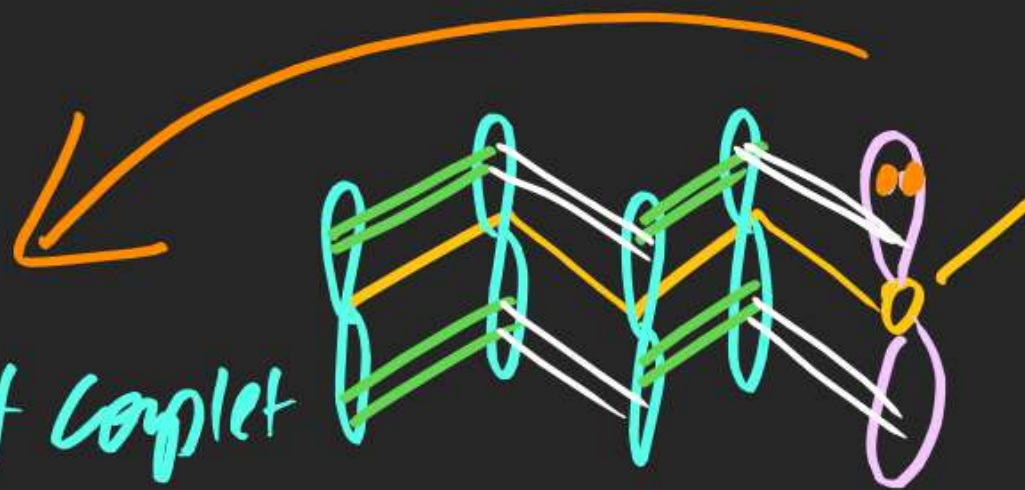
Ad. PYQ
(17)
JIT



octet complete

octet complete

octet incomplete

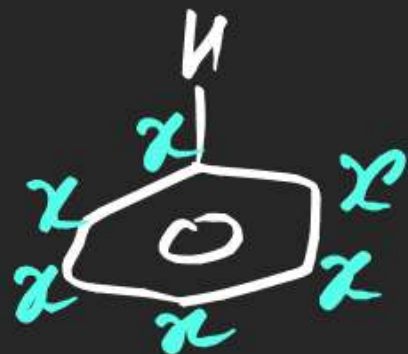


least stable RS

Resonance / Mesomeric Effect:-

Resonance

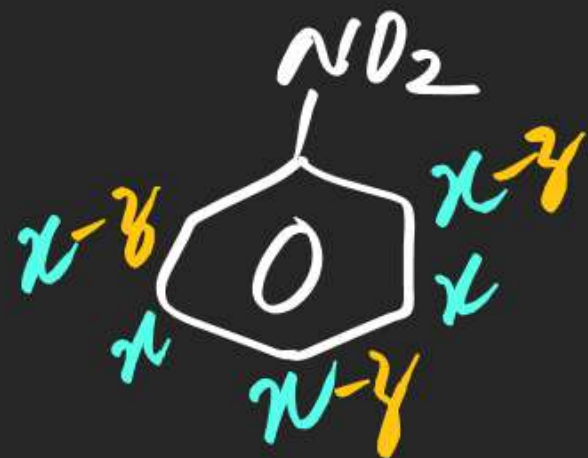
Resonance effect



✓



✓

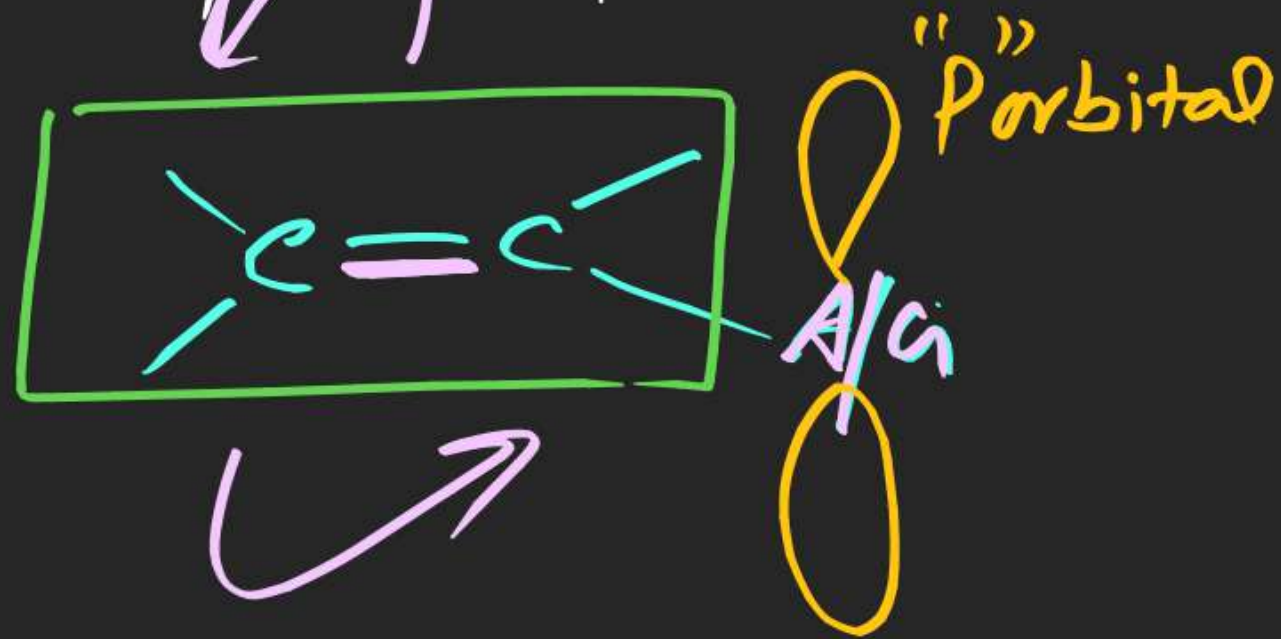


✓

Resonance effect of NH₂
(Electron donating)

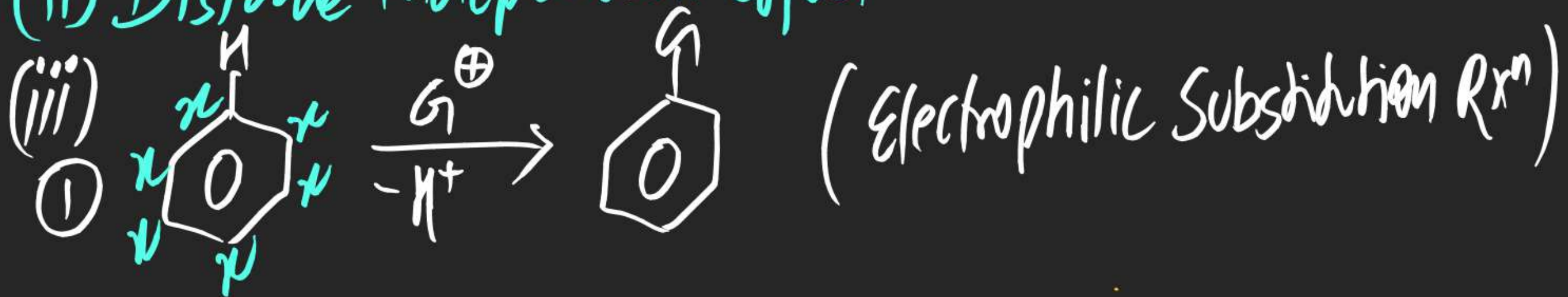
Resonance effect of NO₂
(Electron withdrawing)

Resonance/Mesomeric Effect (R/M effect) Permanent displacement of π e^- density due to p orbital of directly attached A/G is known as R/M effect of that A/G on conjugated system.

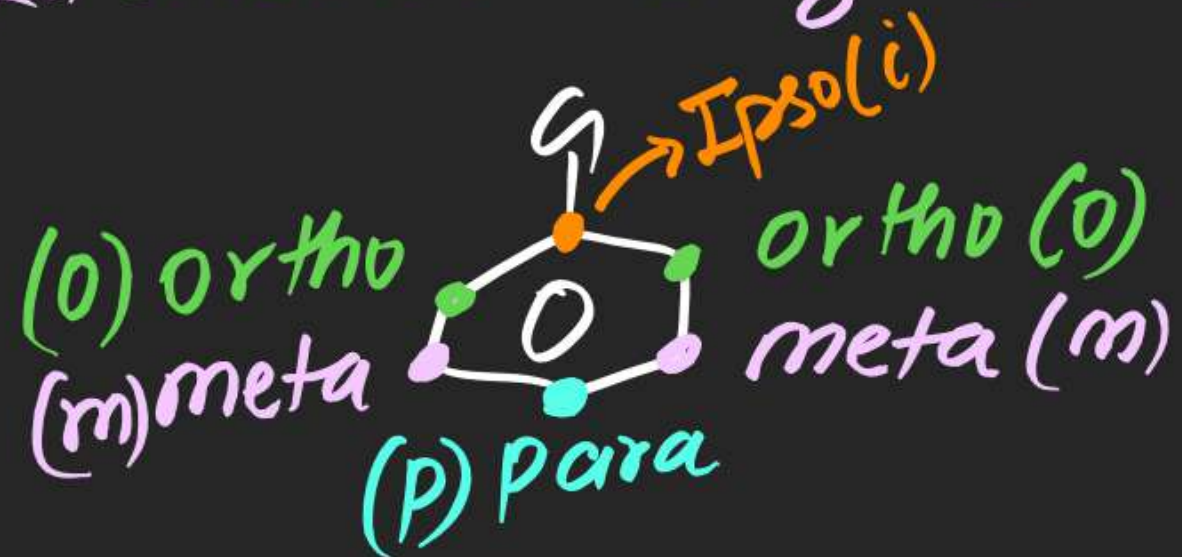


Note

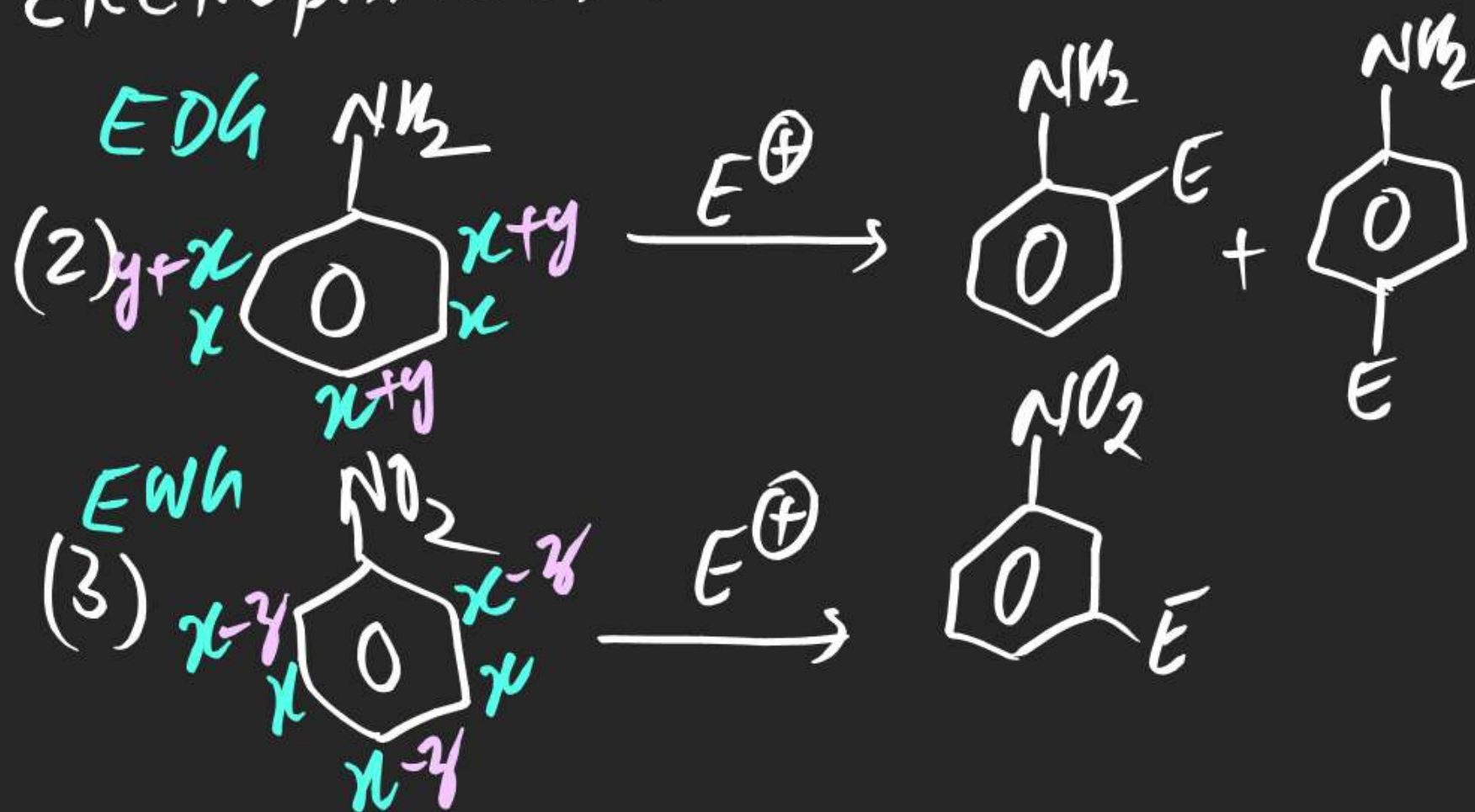
- (i) Permanent effect
- (ii) Distance independent effect



(iv) In Substituted Benzene



(v) Electrophilic Aromatic Substitution



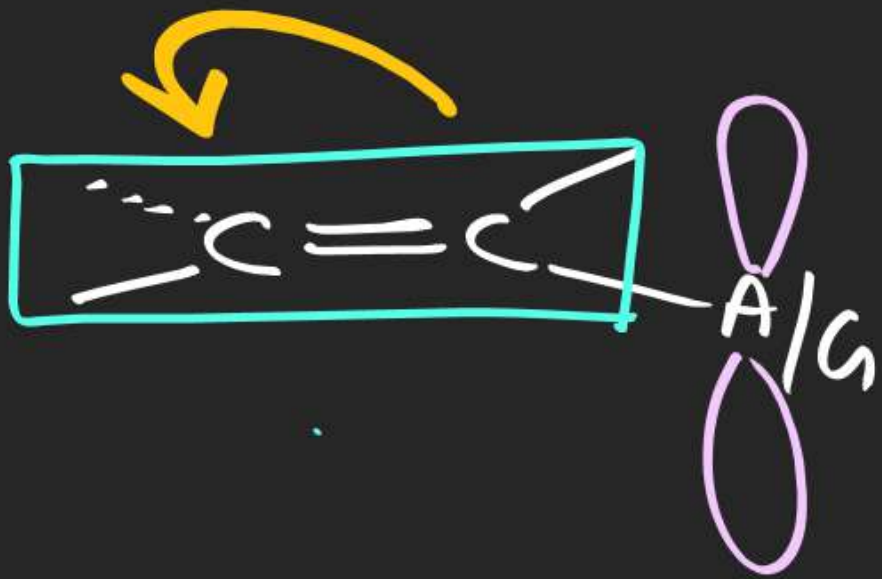
देने वाला देता है तो
ortho & para पर देता है.
hence ortho & para directing

लेने वाला लेता है तो
ortho & para से ही लेता है
hence meta directing.

Type of Resonance effect:

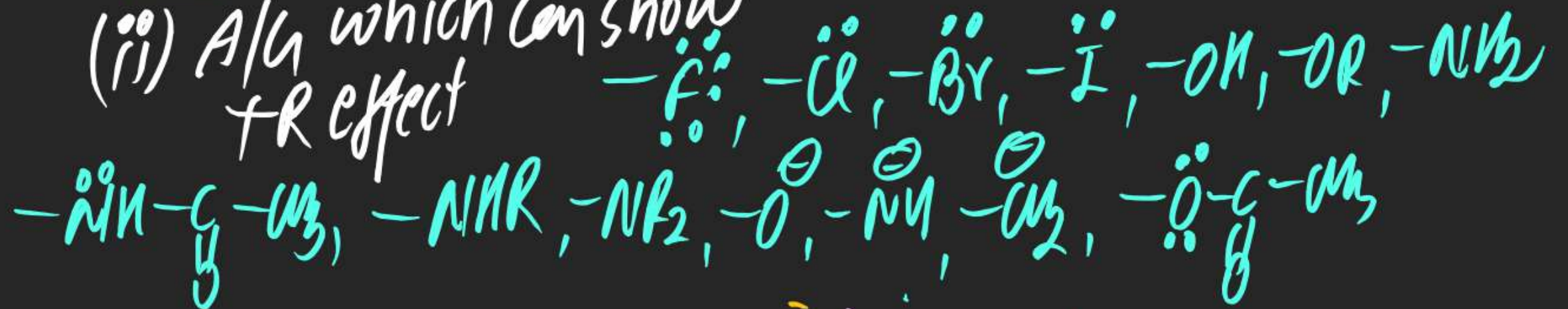
⇒ There are two types of Resonance effect.

(i) +Resonance/+mesomeric effect : Permanent polarisation of πe^- density due to "p" orbital of attached A/G away to itself is known as (+R/+m) effect +R/+m effect of A/G "p" orbital



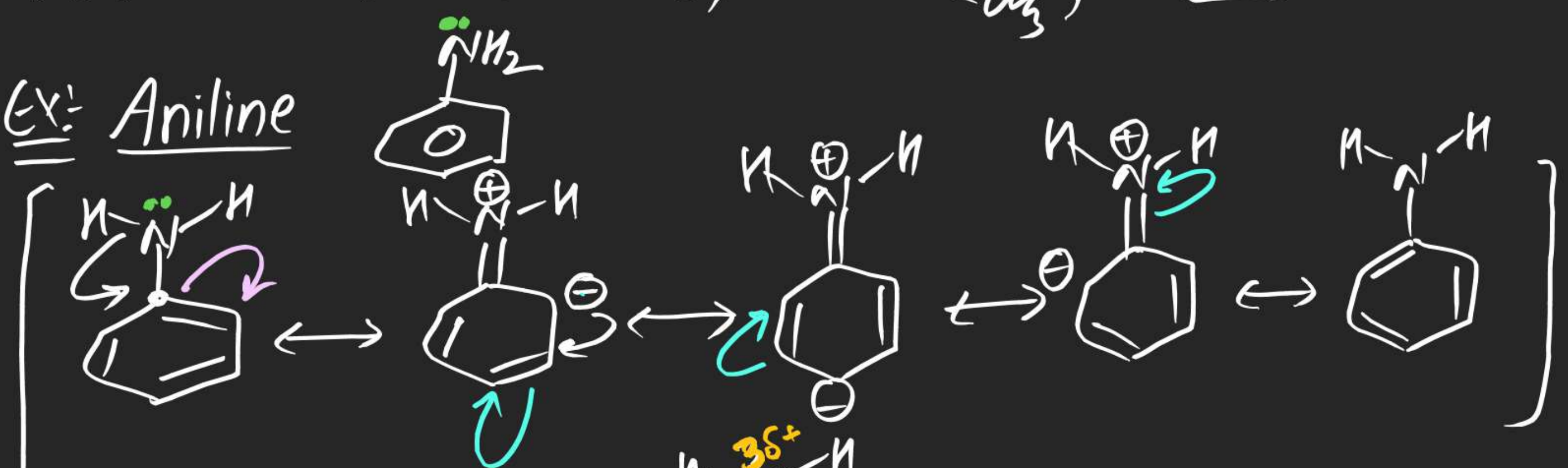
Note (i) +R/+m effect is Electron donating effect

(ii) A/G which can show +R effect

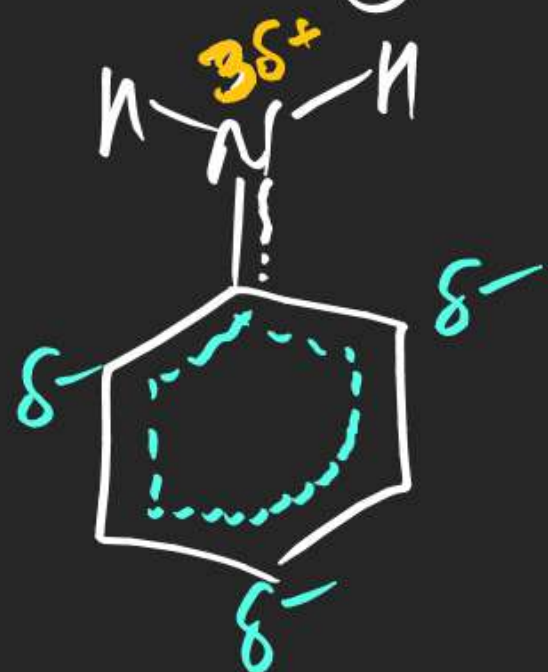




Ex: Aniline

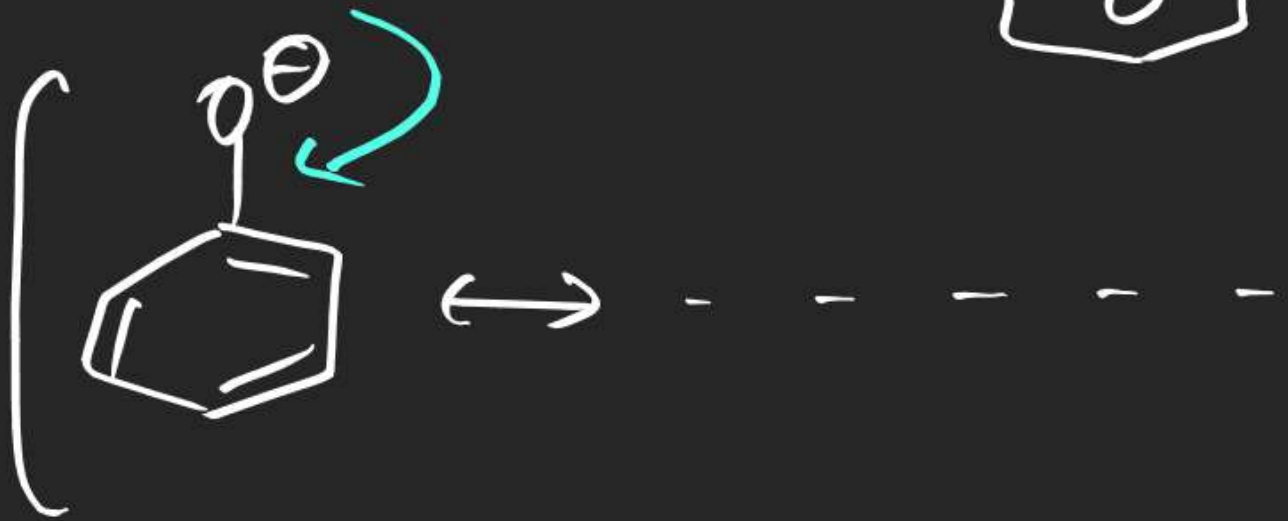


Total "5"
(RS)



(Resonance effect)

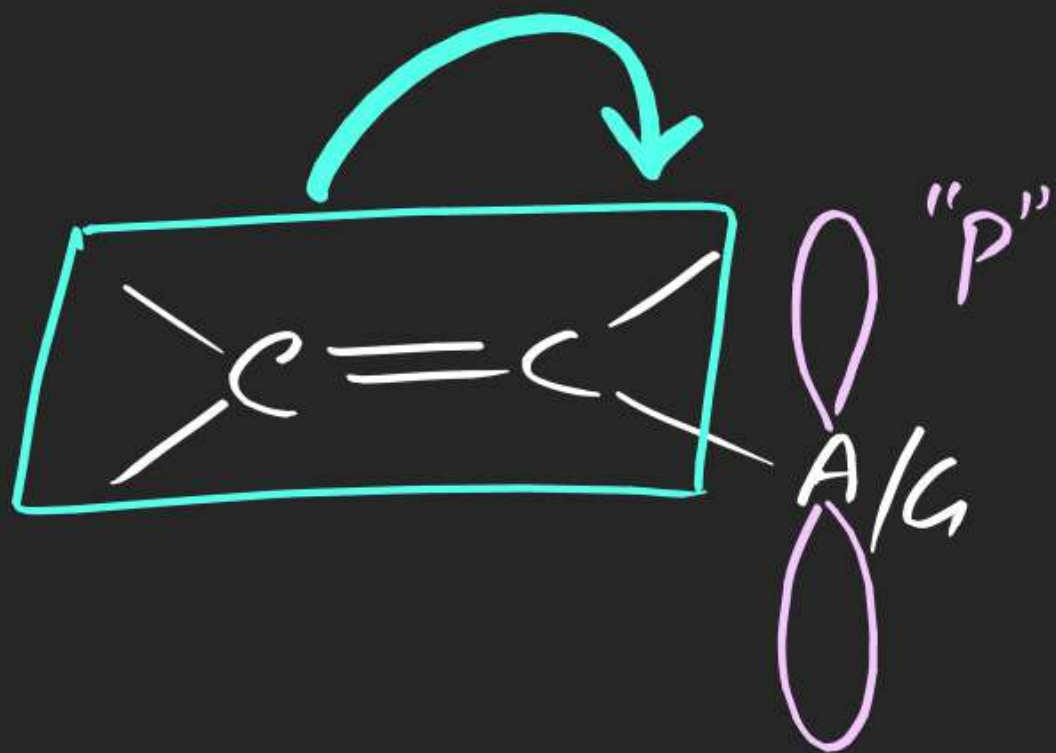
Ex! Phenoxide Ion



Note

+R series

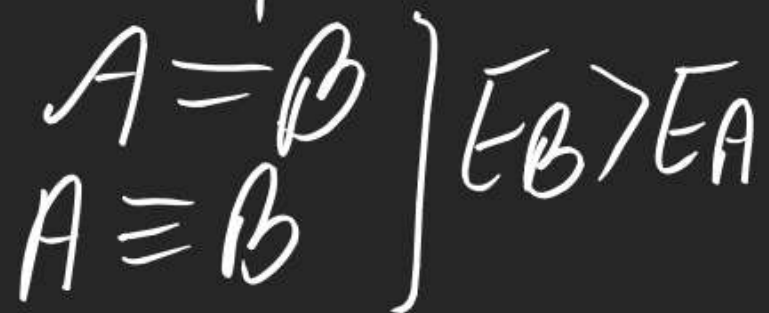
(#) - Resonance / - mesomeric effect (-R/-M effect):

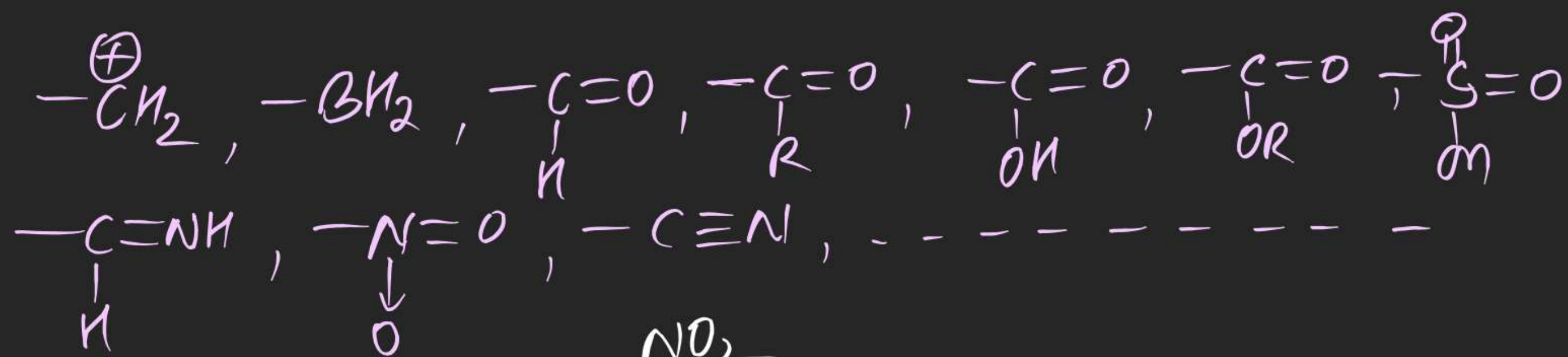


Permanent displacement of π e^- density toward attached Alk due to its "p" orbital are known as -R/-M effect of Alk

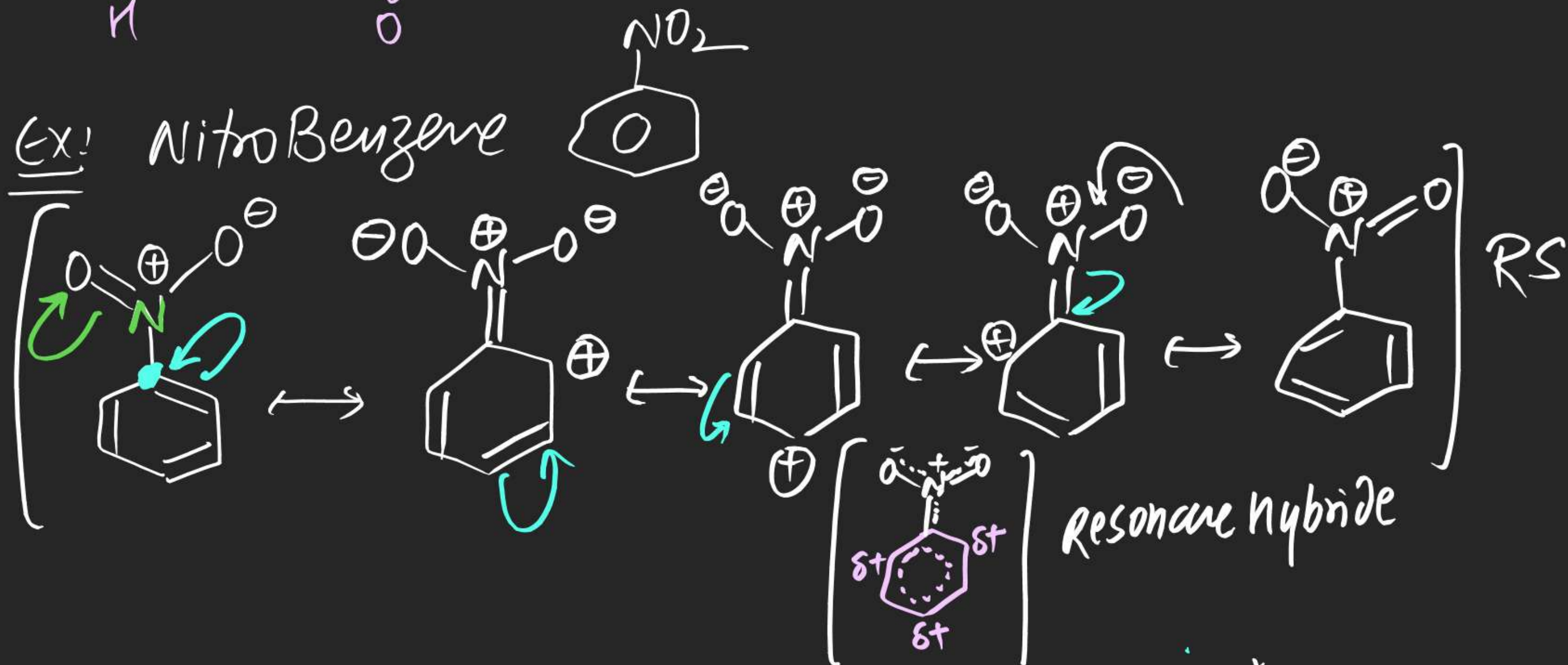
Note (i) -R/-M effect is e^- withdrawing effect

(ii) -R/-M effect is shown by groups having either exactly vacant orbital or may be vacant (polarised bonds)

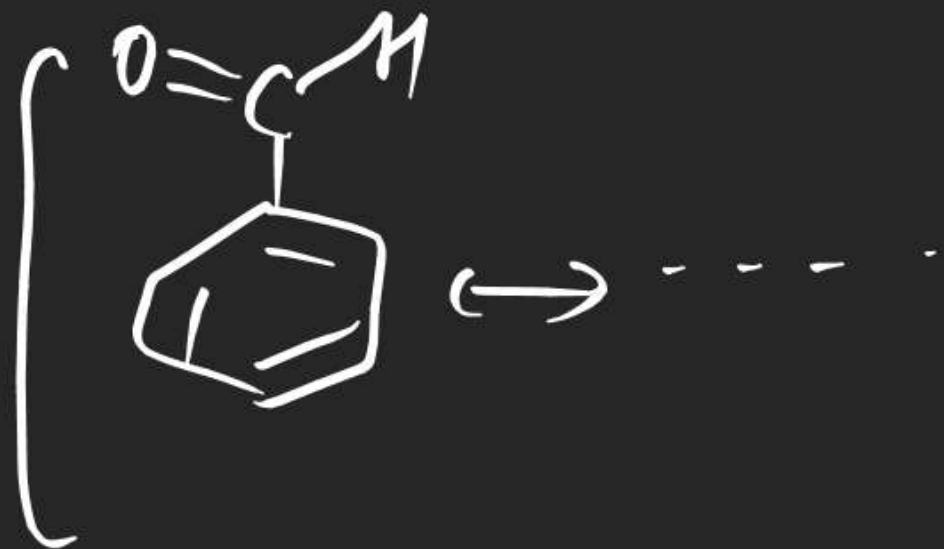
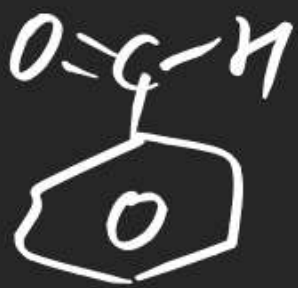




Ex! Nitrobenzene



Ex! Benzaldehyde



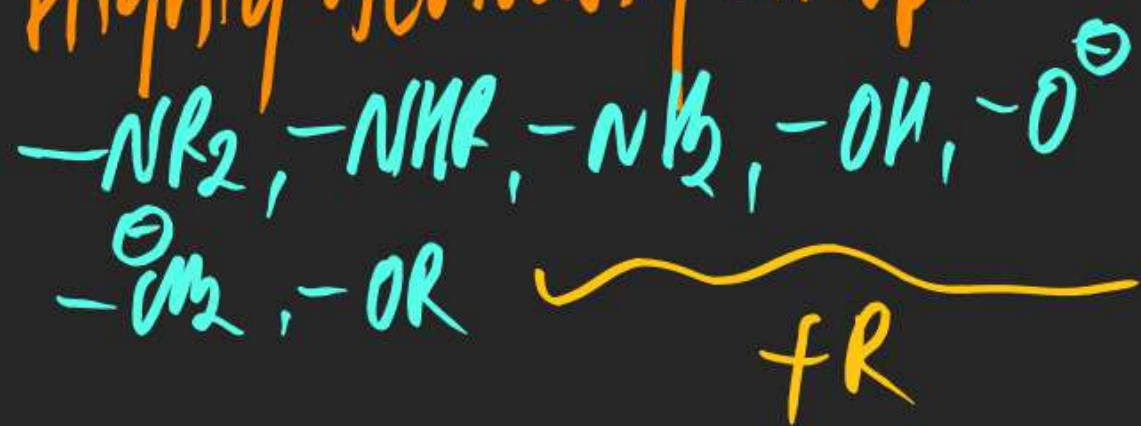
- R series :-

\Rightarrow order of Rate of electrophilic attack.



Activating Compounds: All compounds which show higher rate of electrophilic substitution than rate of electrophilic substitution of Benzene, are known as Activating Compounds.

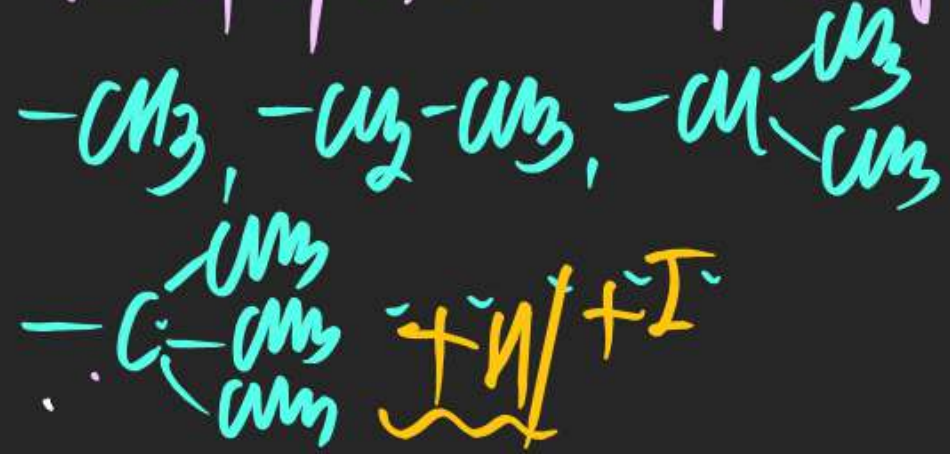
Highly Activating Groups



moderately Activating Groups



Weakly Activating Group



Deactivating Compounds All compounds which show slower rate of electrophilic substitution than Benzene, known as Deactivating Compounds.

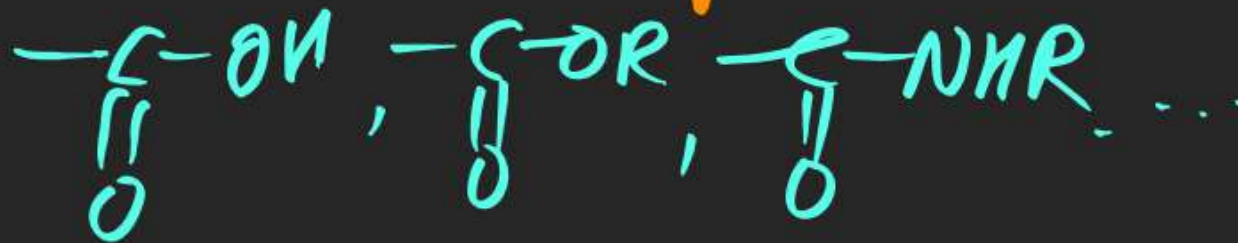
Weakly deactivating Groups

halogens



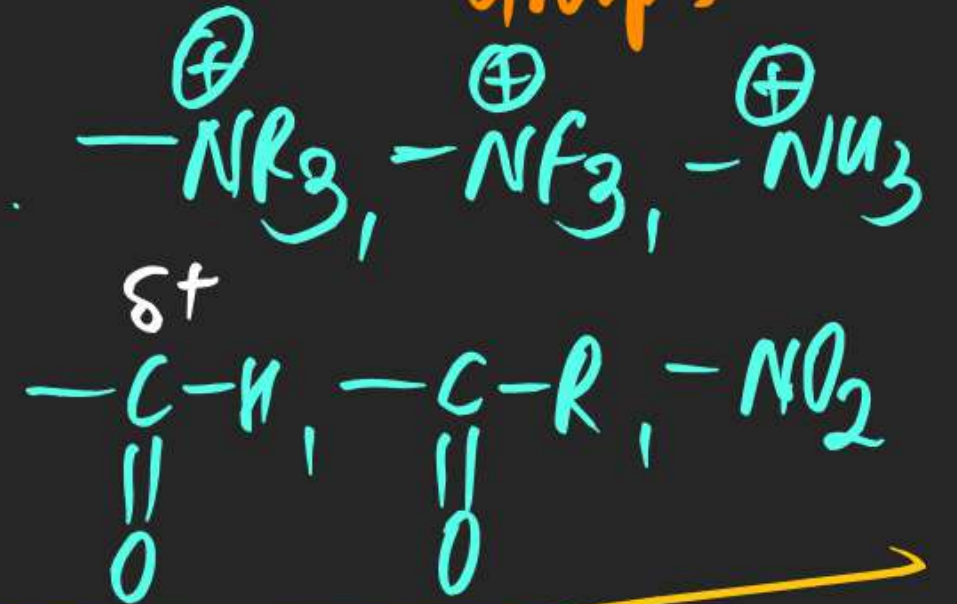
+R

moderately deactivating Groups



-R

Strongly deactivating Groups.



Note: