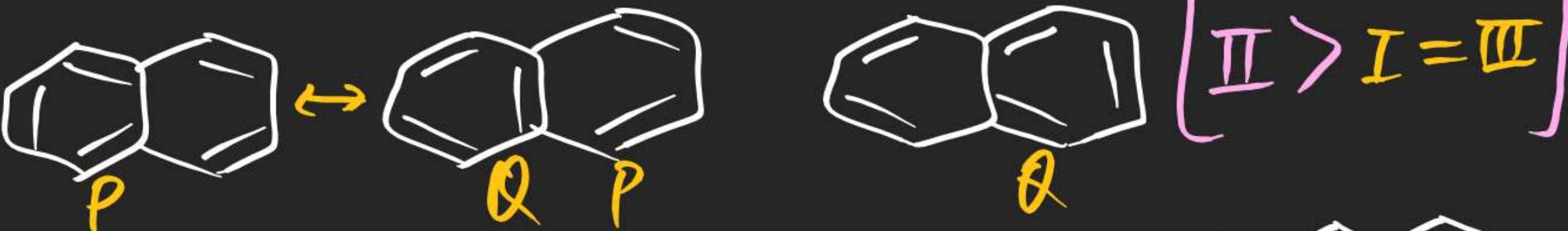


⑤ RS having higher no. of Benzenoid segment is more stable.

(9)



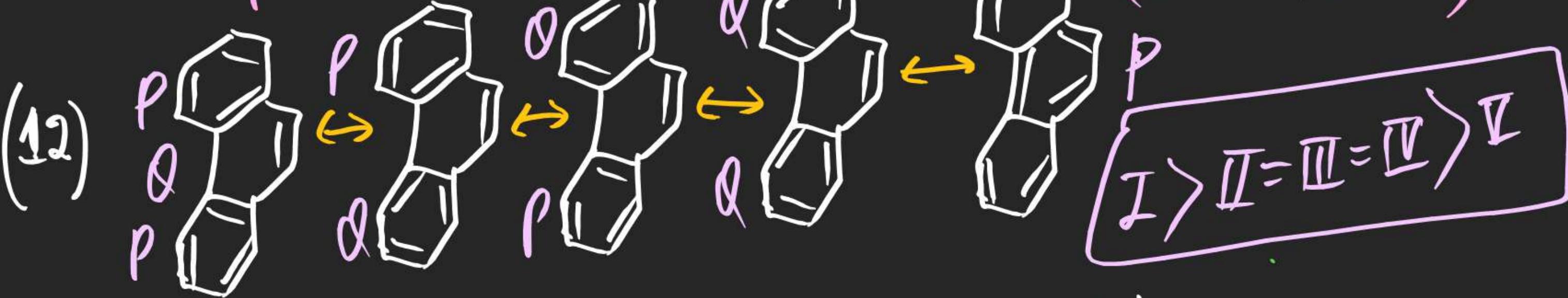
(10)



(11)

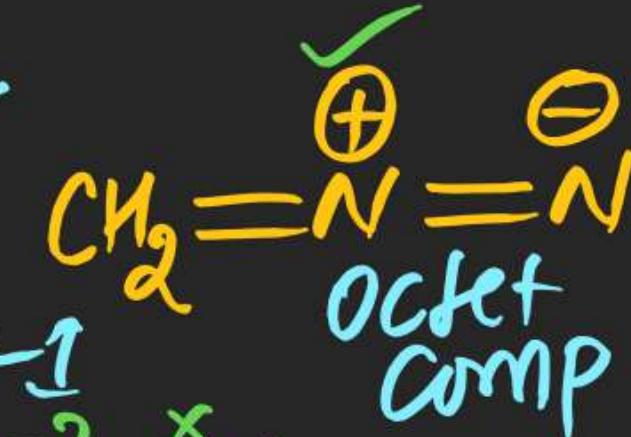


(12)



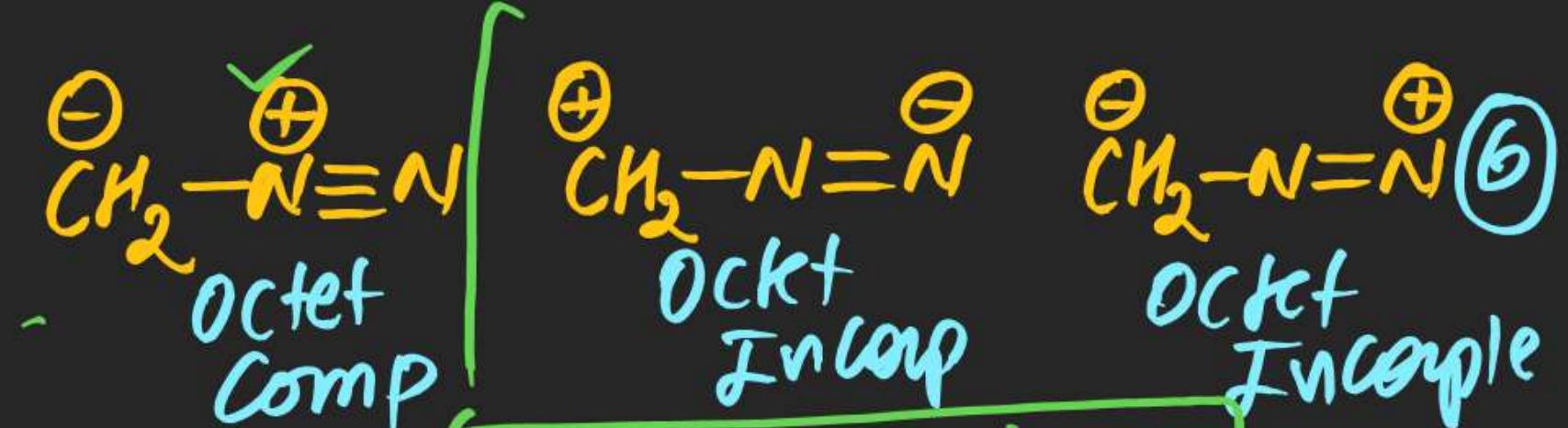
IIT Ad

(13)



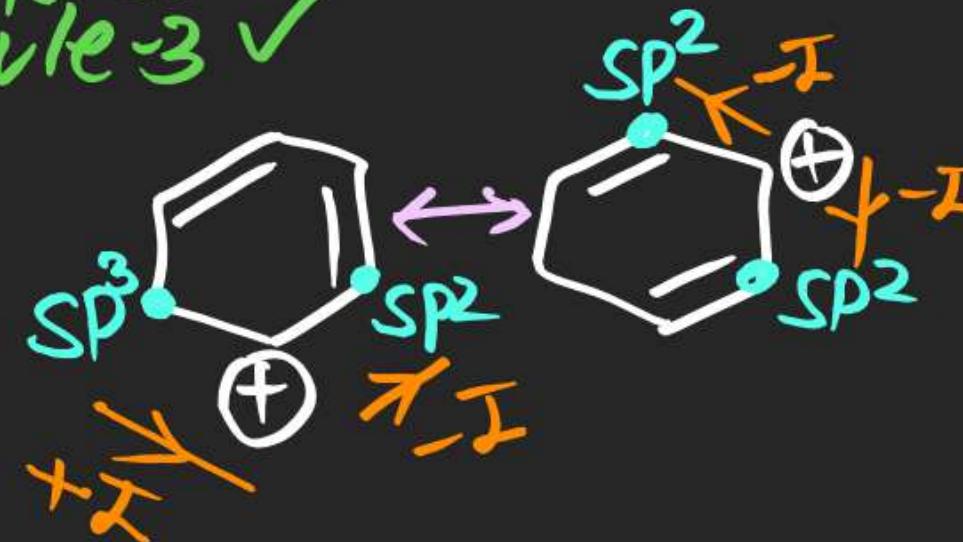
Rule-1

Rule-2 ✗
Rule-3 ✓

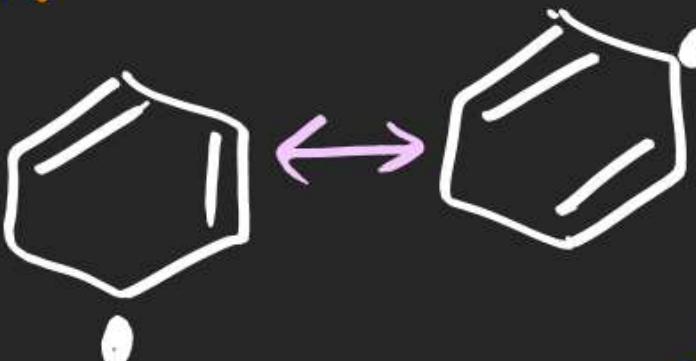


Rule-2 ✗
Rule-3 ✓

(14)

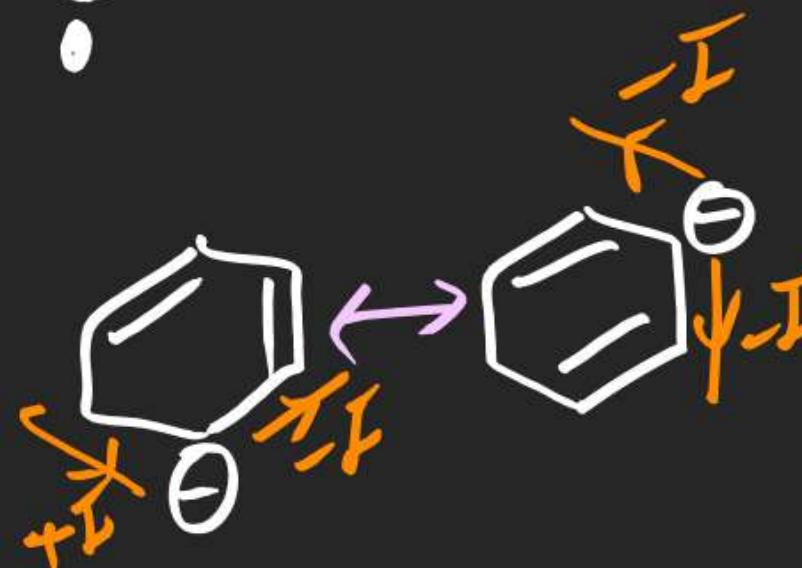


(15)

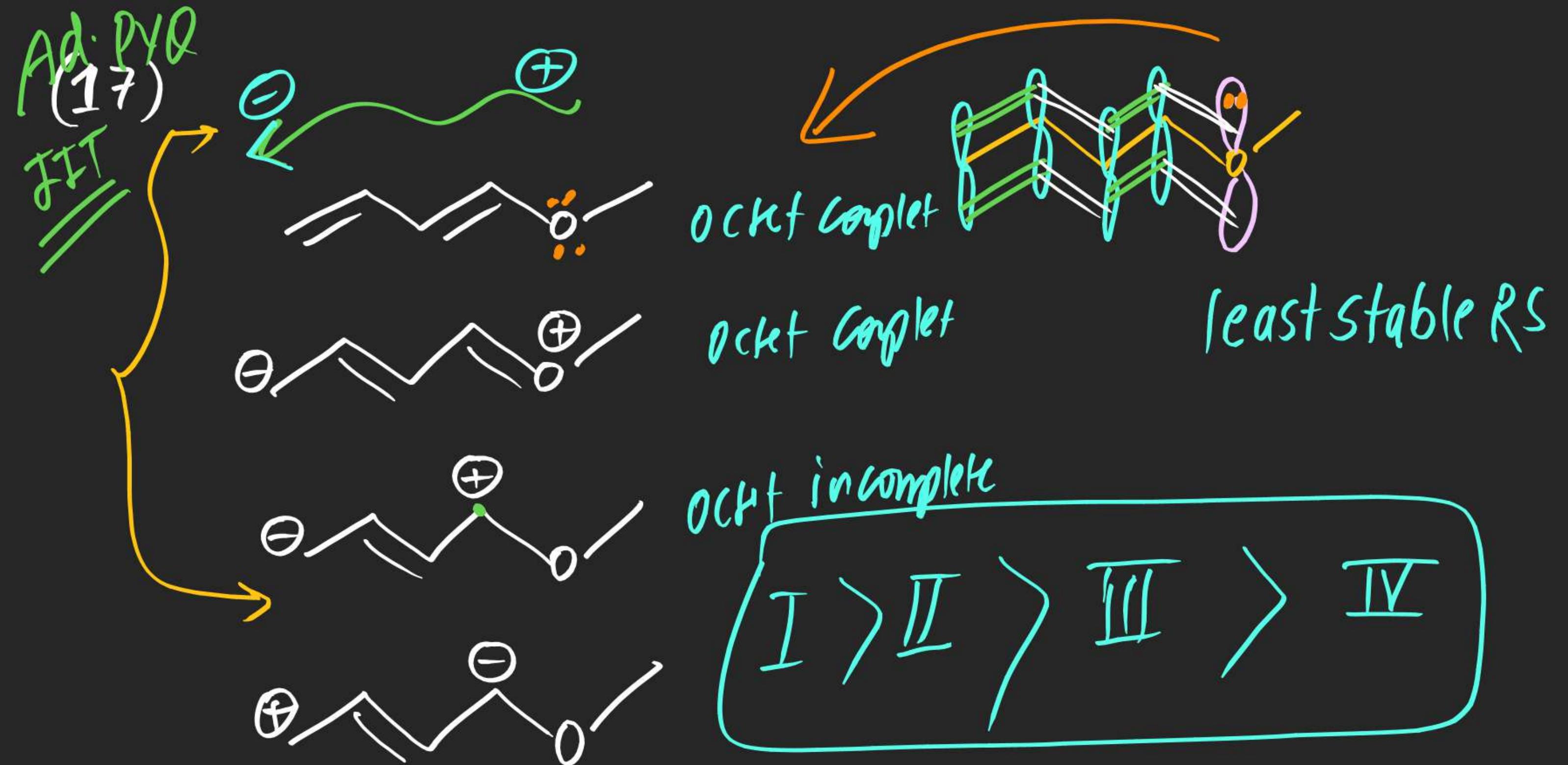


I > II

(16)

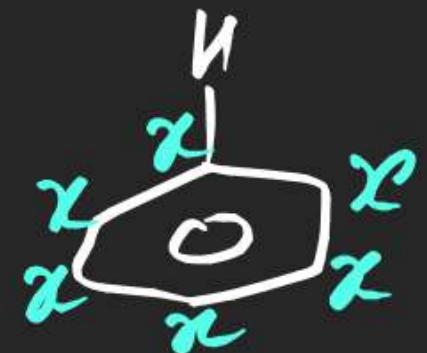


II < I



Resonance / mesomeric Effect:-

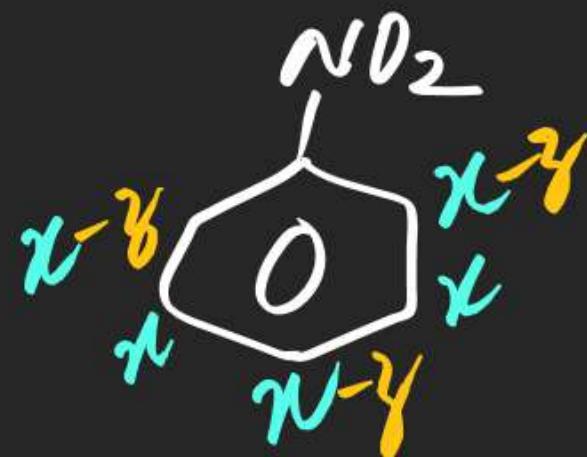
Resonance



Resonance effect

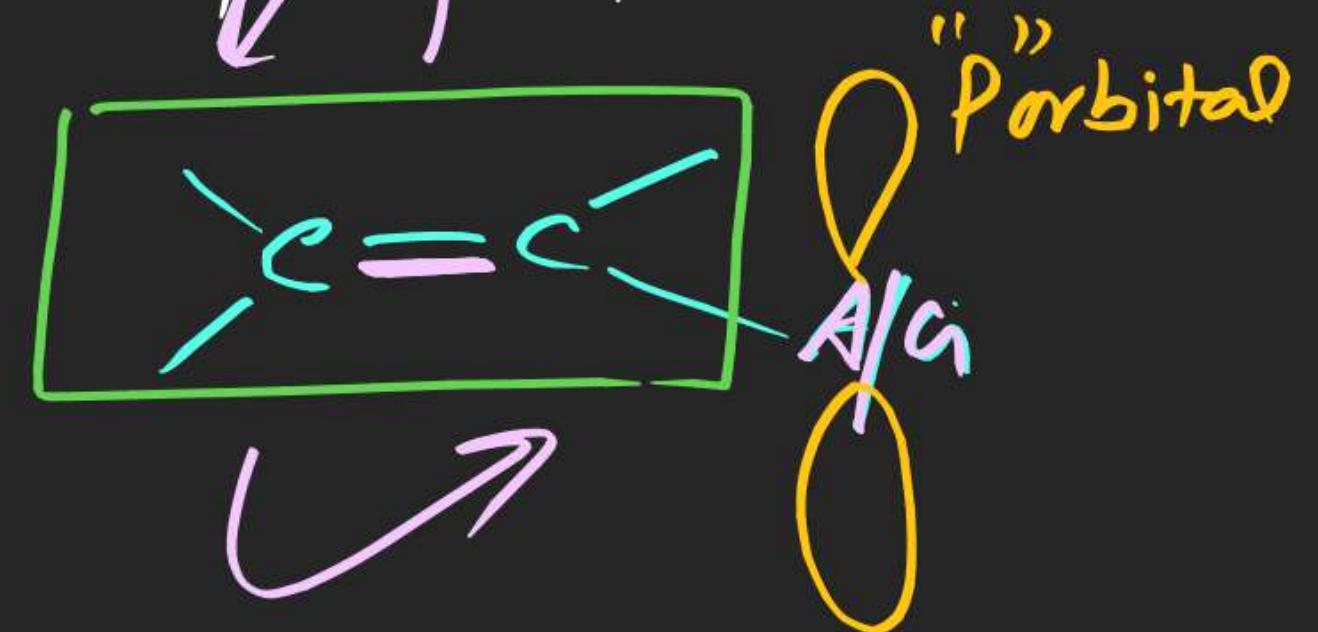


Resonance effect of NH_2
(Electron donating)



Resonance effect of NO_2
(Electron withdrawing)

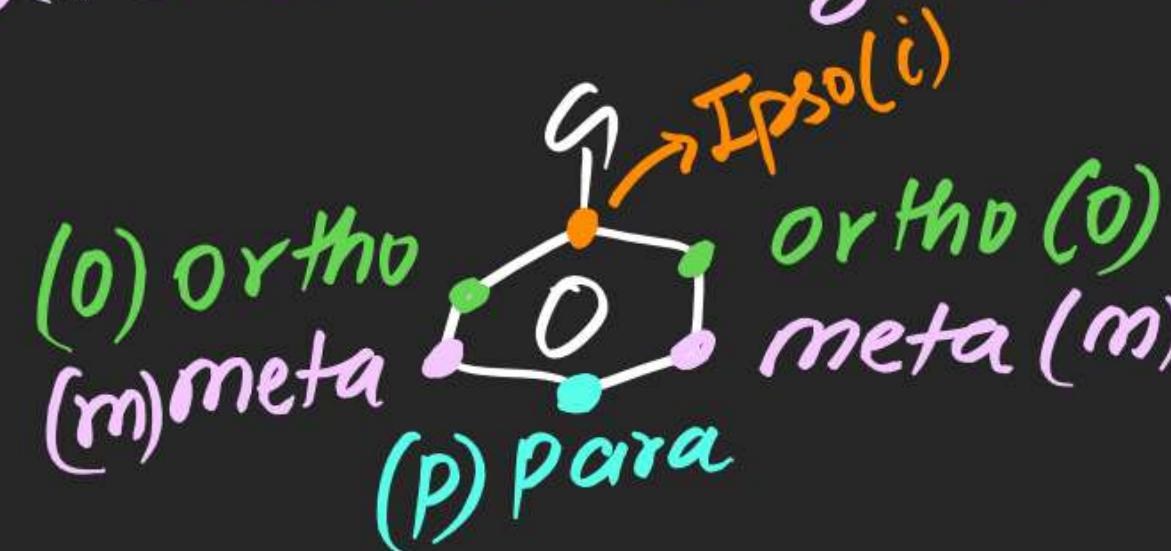
Resonance/Mesomeric Effect (R/m effect) Permanent displacement of πe^- density due to p orbital of directly attached A/G is known as R/m effect of that A/G on conjugated system.



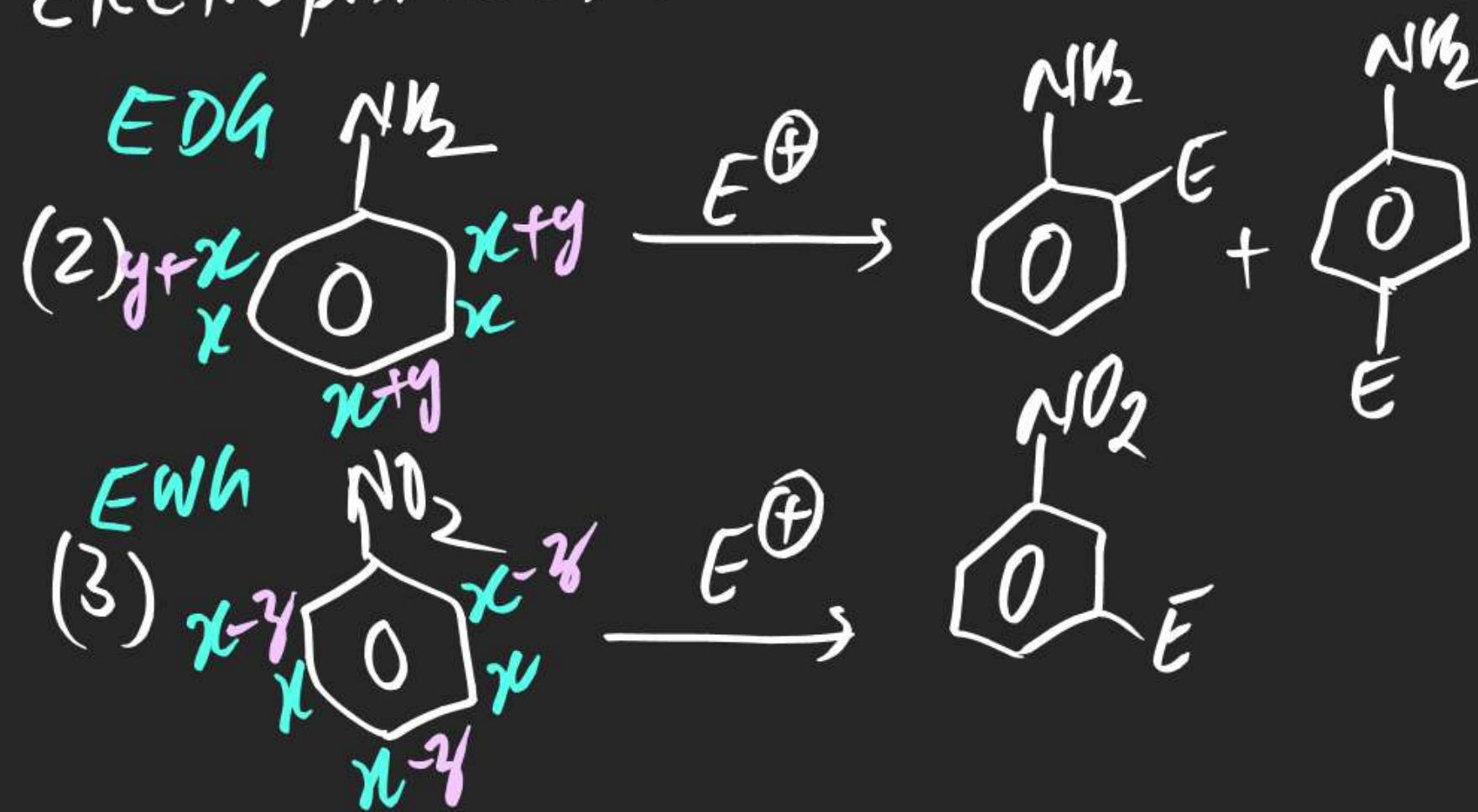
- Note
- (i) Permanent effect
 - (ii) Distance independent effect



(iv) In Substituted Benzene



(v) Electrophilic Aromatic Substitution



दूसरे वाला देता है तो
 ortho & para पर देता है.
 hence ortho & para dimly

चौथे वाला देता है तो
 ortho & para में ही देता है.
 hence meta dimly.

Type of Resonance effect:

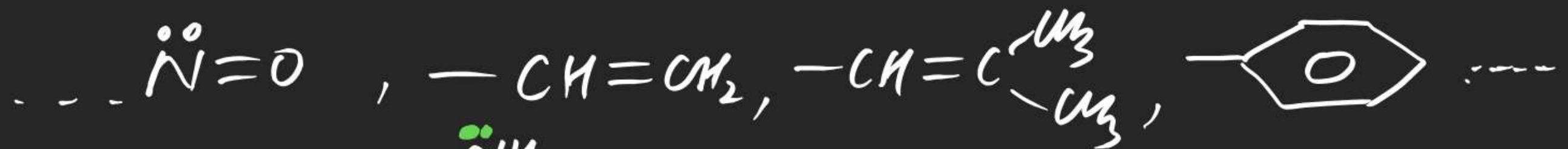
⇒ There are two type of Resonance effect.

(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 of A/G
 "P" orbital

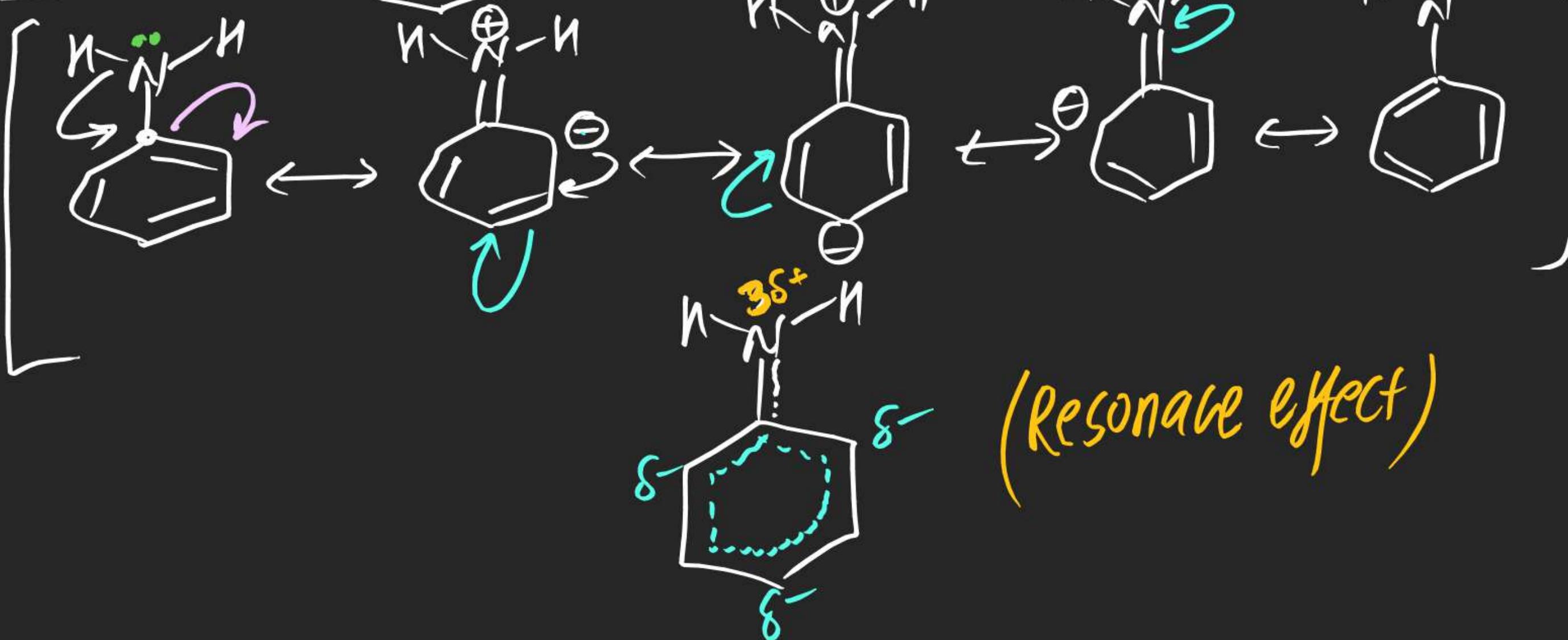
Note (i) +R/+m effect is Electron donating effect

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

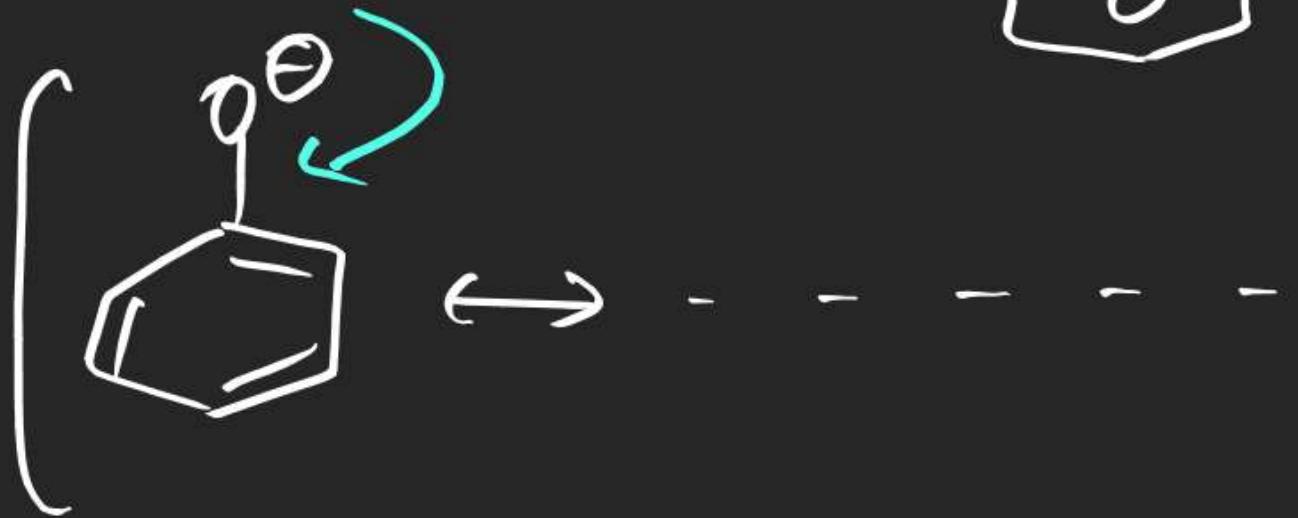




Ex: Aniline



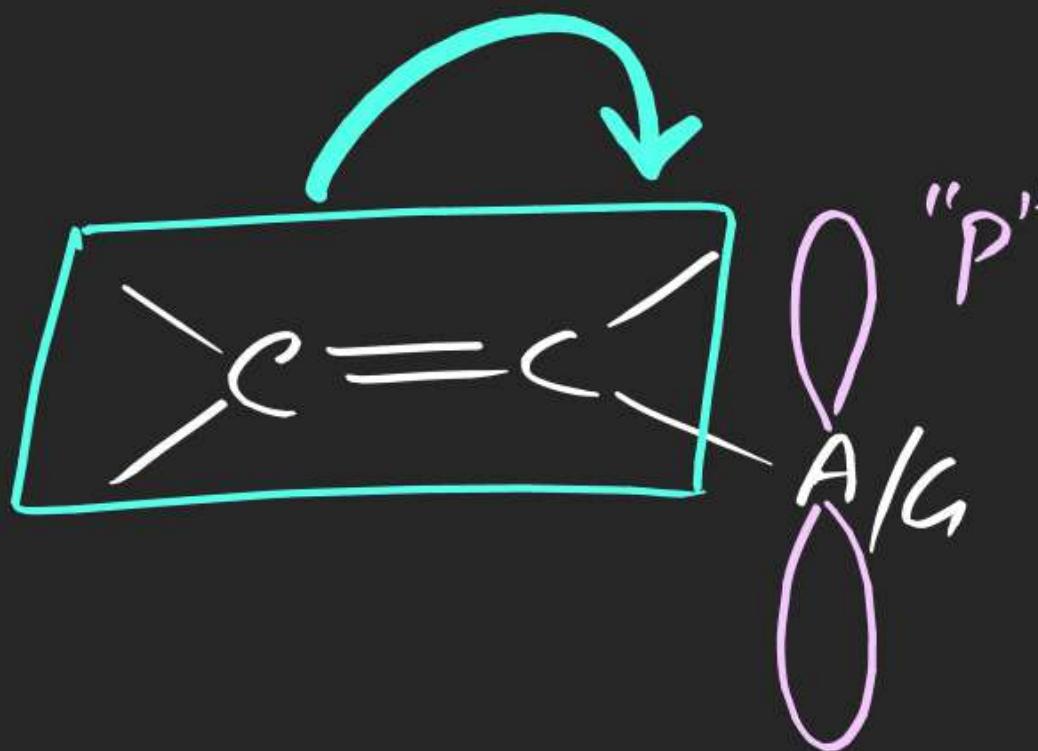
Ex: Phenoxide Ion



Note

+ R Series

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

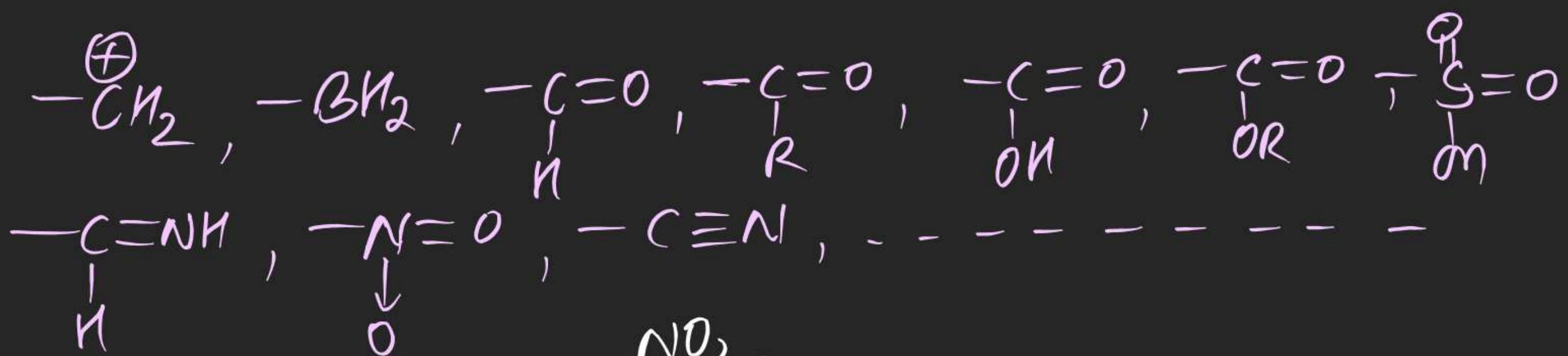


- movement of π e^- density toward attached A/G
- due to its "P" orbital are known as $-R/-M$
- effect of A/G

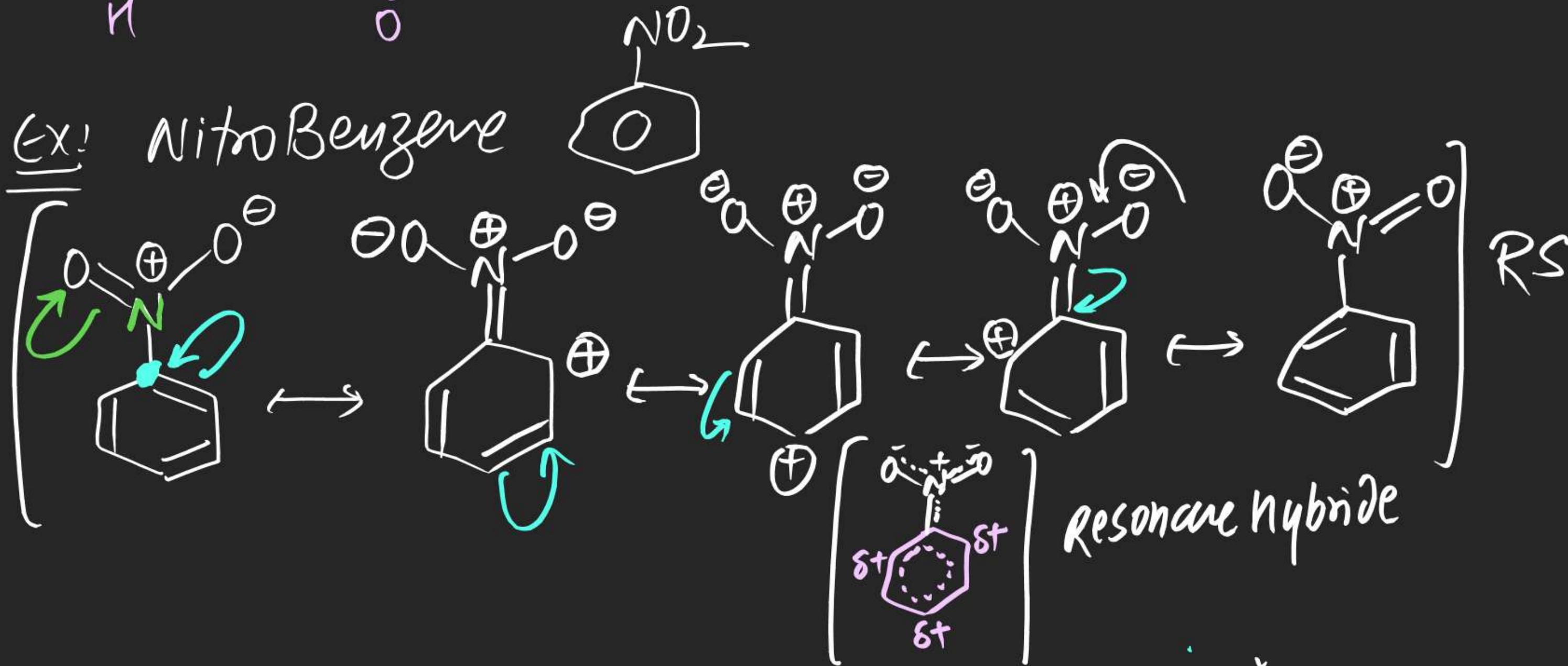
Note (i) $-R/-M$ effect is e^- withdrawing effect

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

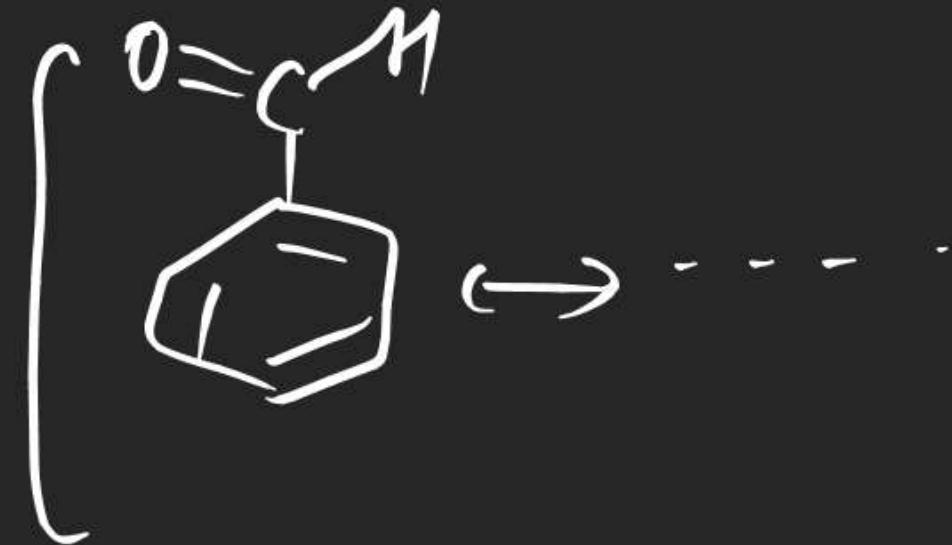
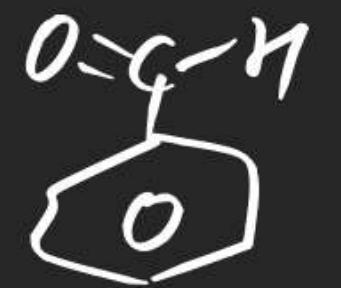
$$\begin{array}{l} A=\beta \\ A=B \end{array} \left. \right\} E_B > E_A$$



Ex: NitroBenzene



Ex: Benzaldehyde

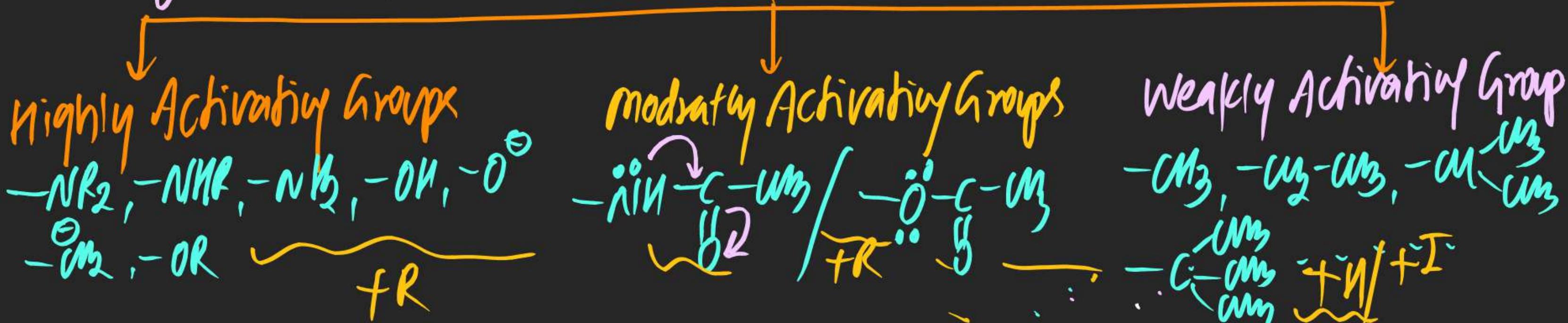


-R *benes*:

⇒ 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.



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

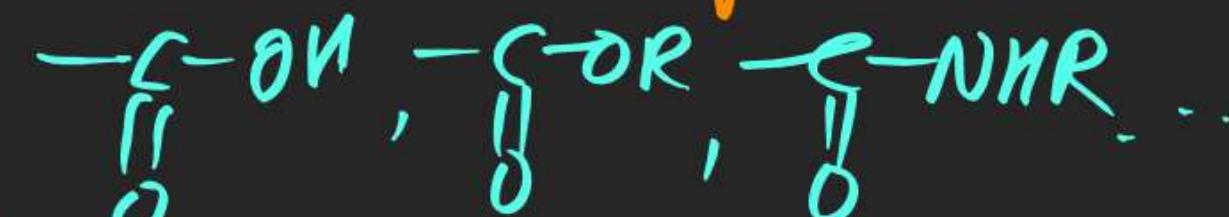
weakly deactivating Groups

halogens



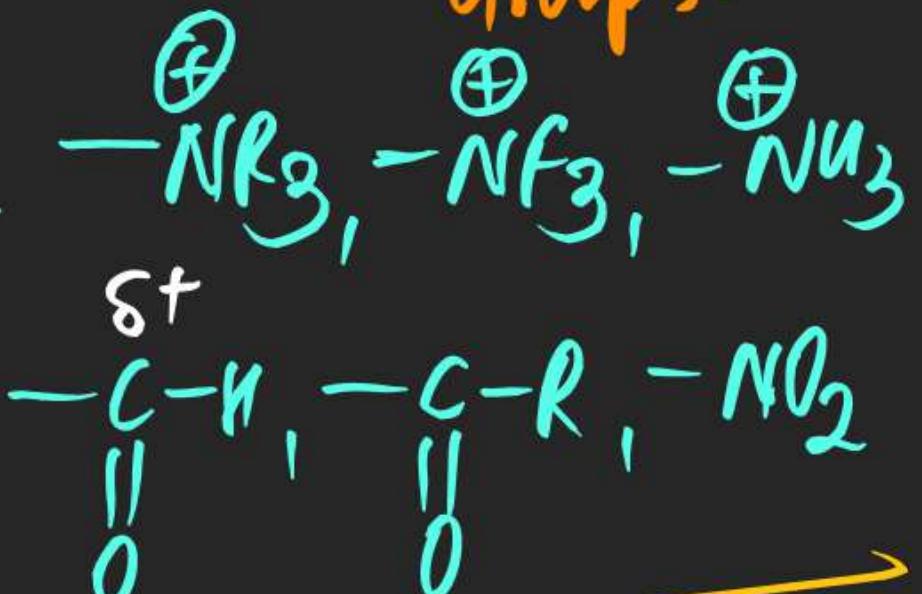
+R

moderately deactivating Groups



-R

Strongly deactivating Groups.



Note: