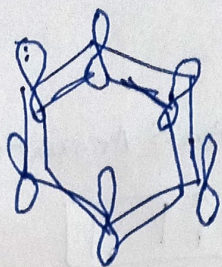
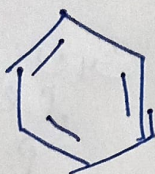
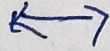
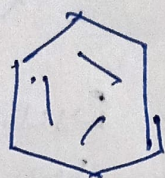


~~Aromaticity~~ Benzene



The ~~orbitals~~ p-orbitals are perpendicular to plane, all bonds lie in same plane.

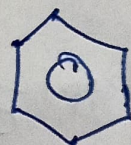


(ERS) Equivalent Resonating Structure.

No. of Resonating Structure of Benzene: (2)

No. of σ bond = $6 + 6 = 12$ (6 from C-C, 6 from C-H)
No. of π bond = 3

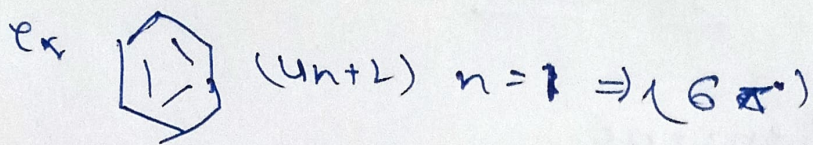
Due to resonance all bonds in C-C have Equivalent Bond length. ERS




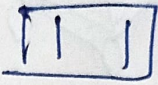

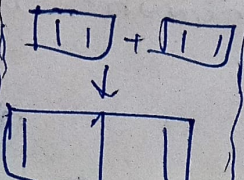
Aromaticity: $(4n+2)e^-$ Huckle Rule

Conditions for Aromaticity:

- ↳ Follow Huckle Rule $(4n+2)e^-$
- ↳ All bond should be in same plane for reso.
- ↳ (All carbon should be sp^2)
- ↳ Compound should be cyclic.
- ↳ Quasi Aromatic are also aromatic.



Non-Aromatic:

Characteristics	Aromatic	Anti Aromatic?	Non-Aromatic
Structure	 Cyclic, planar sp^2	 Cyclic, planar sp^2	 Cyclic/acyclic sp^2, sp, sp^2
No. of πe^-	$(4n+2) \pi e^-$	$4n \pi e^-$	Any πe^-
MOT	Upaired e^- in bonding MO.	Some πe^- in non-bonding MO orbital.	πe^- in bonding & Non-bonding M.O.
Overlapping	Favourable overlapping of p-orbital	Unfavourable overlapping of p-orbital	Simple overlapping like alkenes.
Resonance Energy	Very high	Zero.	4 ± 8 kcal/mol
Stability	High stability	Unstable	Normal stability
Characteristic Redn	Electrophilic Aromatic Subst.	Dimerization to gain stability 	Electrophilic Aromatic redn like normal alkenes.