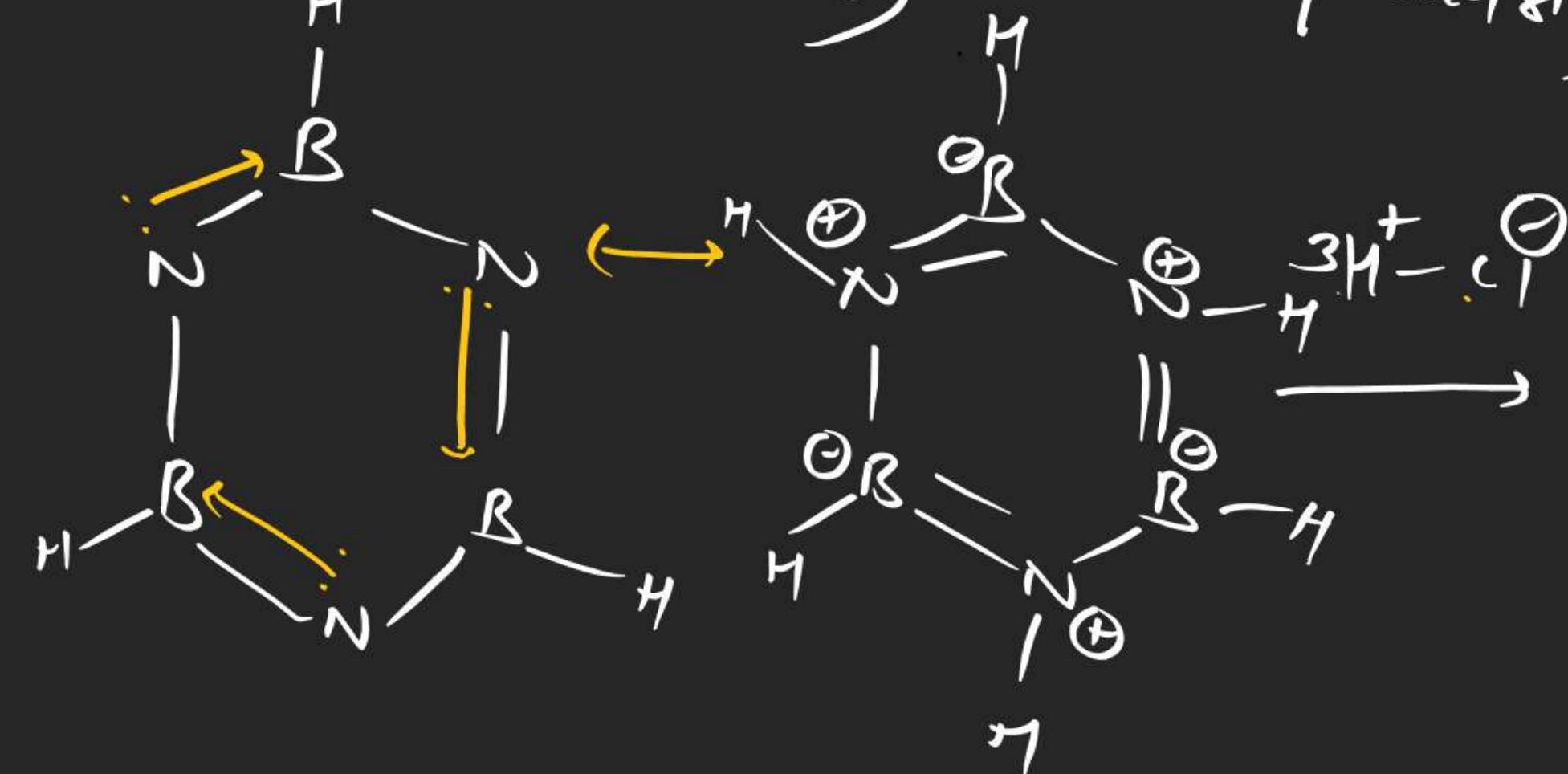
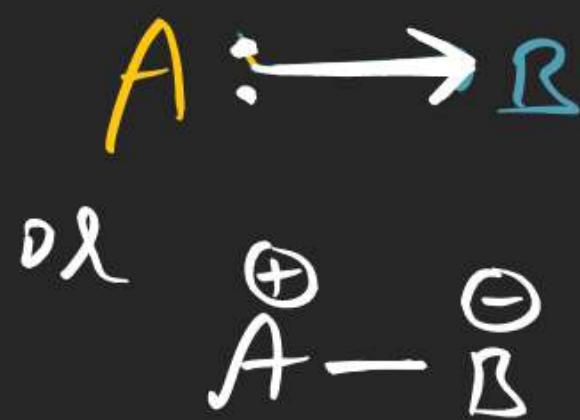
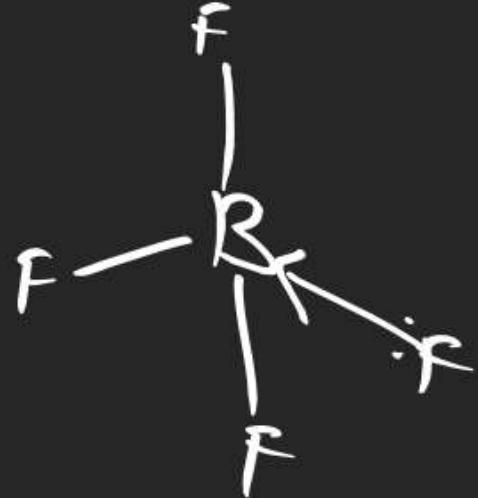


CHEMICAL BONDING

Condition of Back bonding

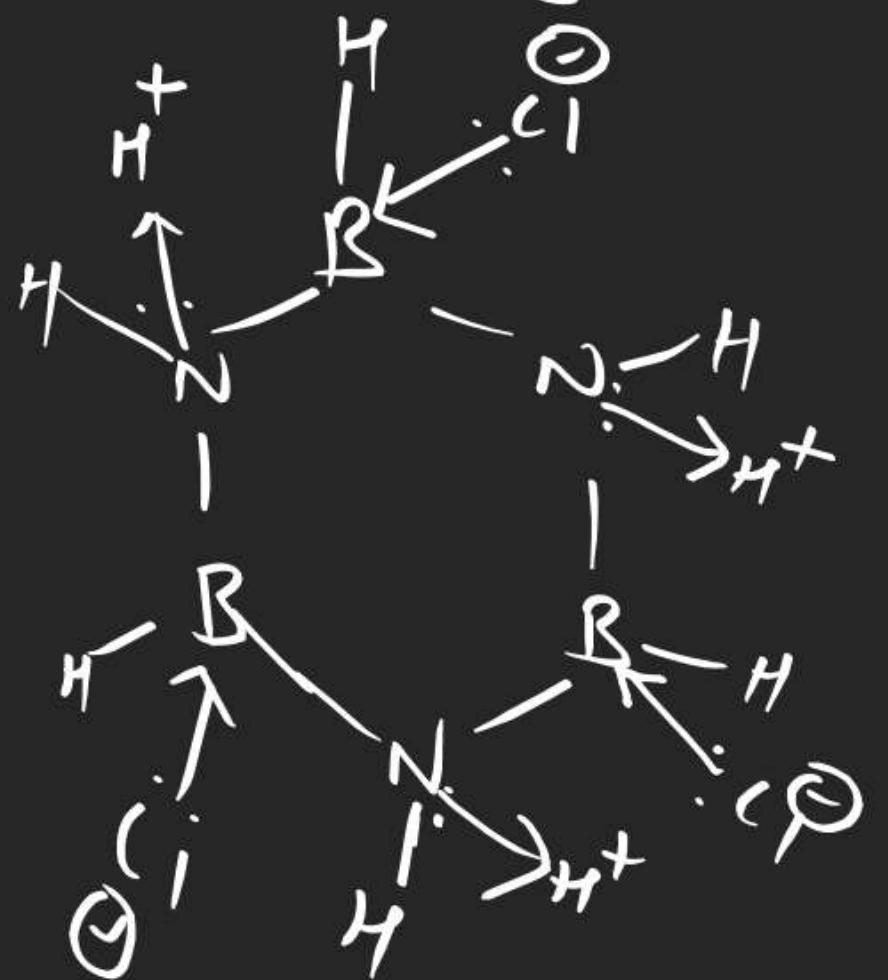
Inorganic benzene ($B_3N_3H_6$) (boronate | borazine)

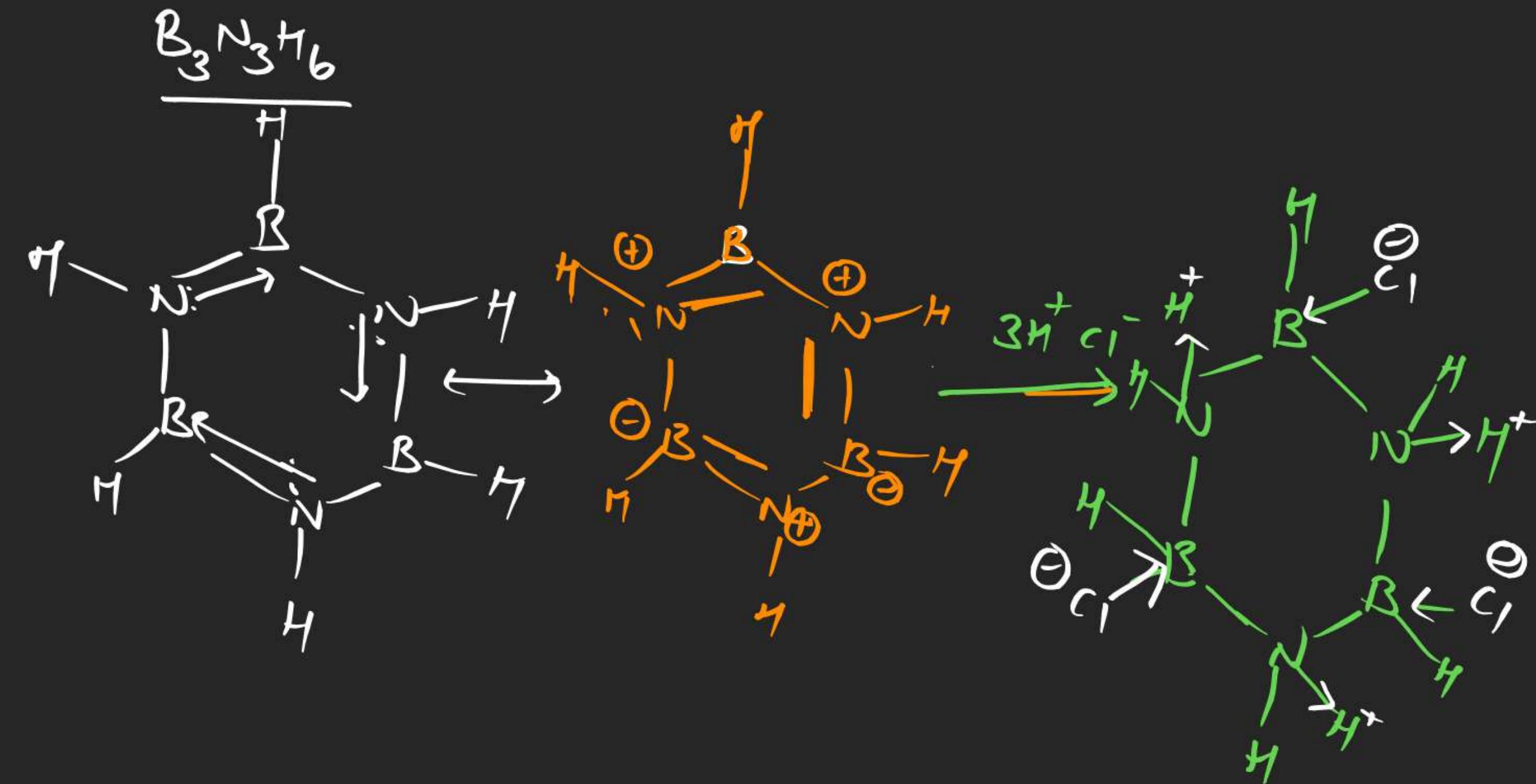




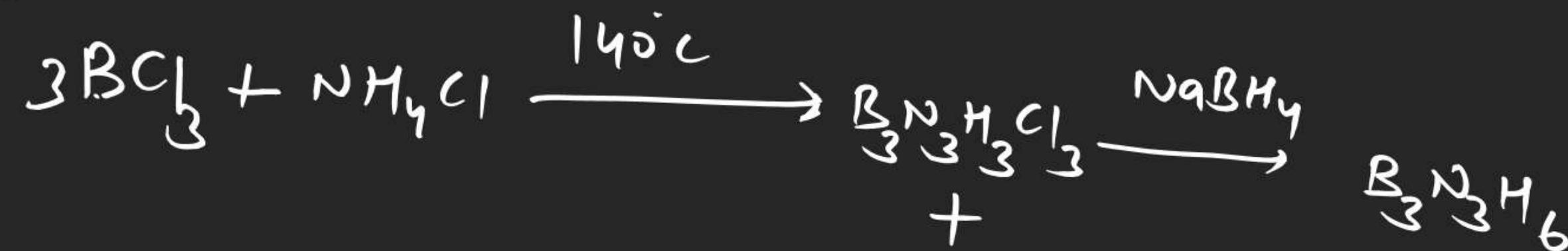


Draw the structure of $\text{B}_3\text{N}_3\text{H}_9\text{Cl}_3$





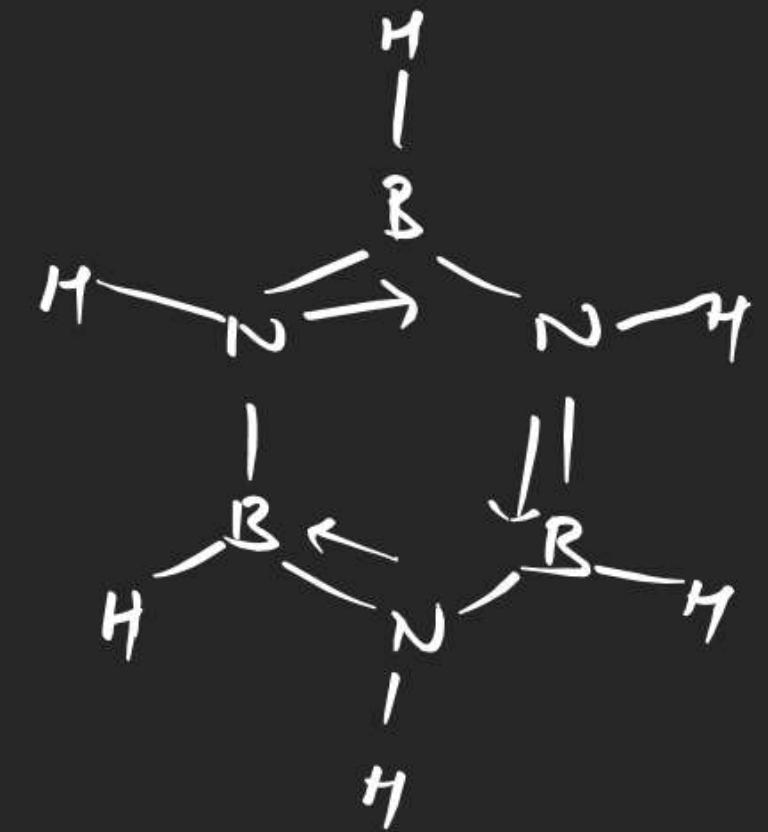
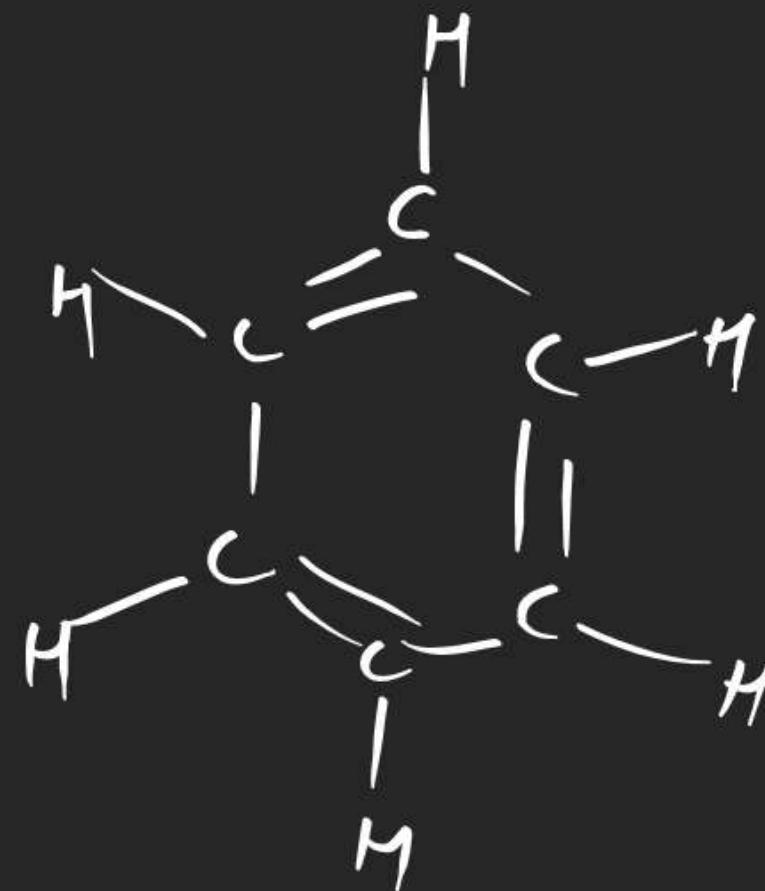
mp



Me-MgBr

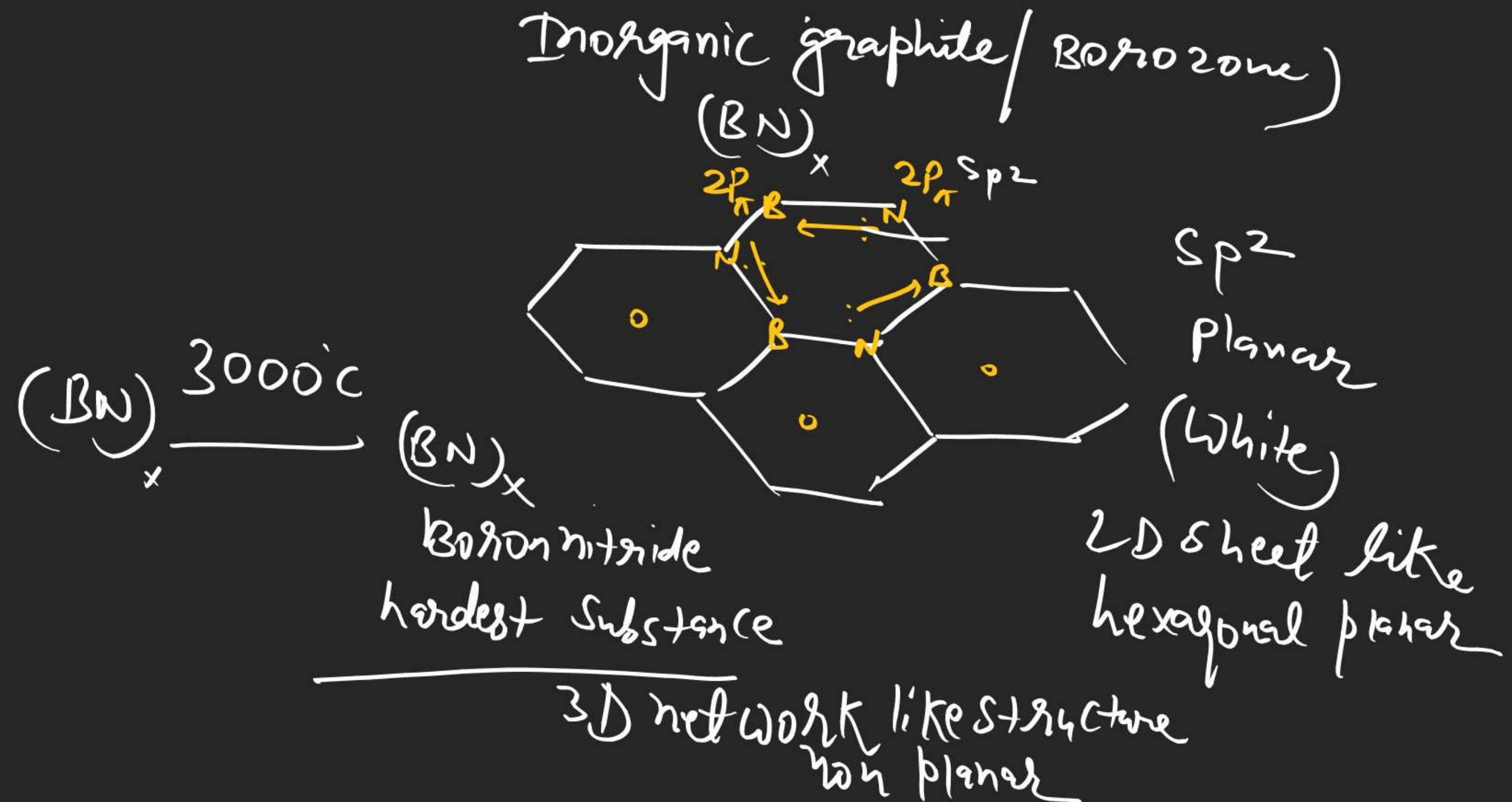


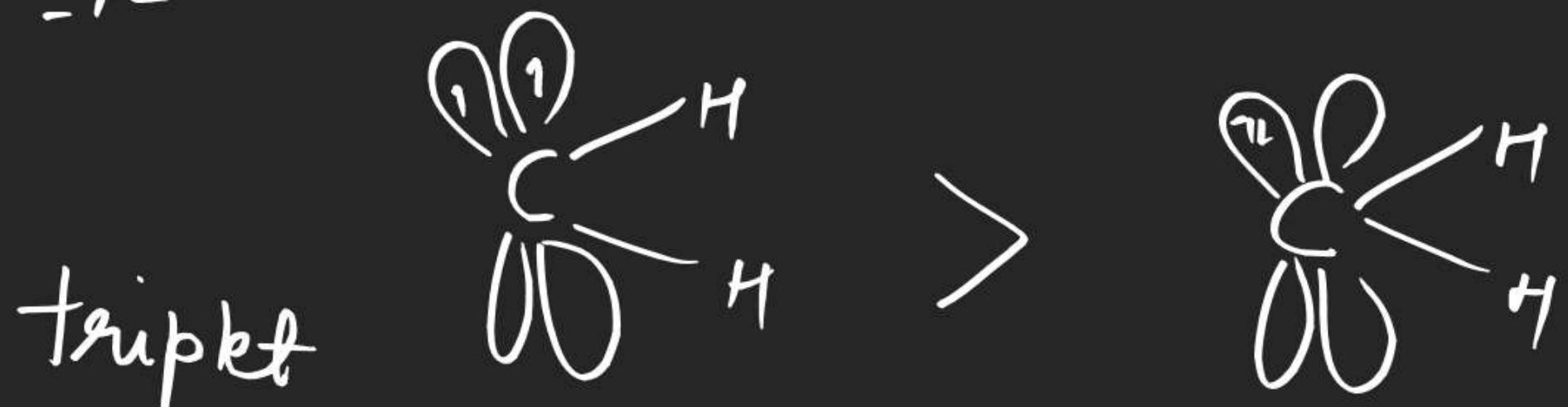
$\text{B}_3\text{N}_3\text{H}_3(\text{Me})_3$



Organic benzene and Inorganic benzene are
ISOstructural.

Note → Inorganic benzene is more reactive than organic benzene due to bond polarity





multiplicity = $2|S| + 1$

S = total Spin

$$\begin{array}{l} 2 \times 1 + 1 \\ 3 \end{array}$$

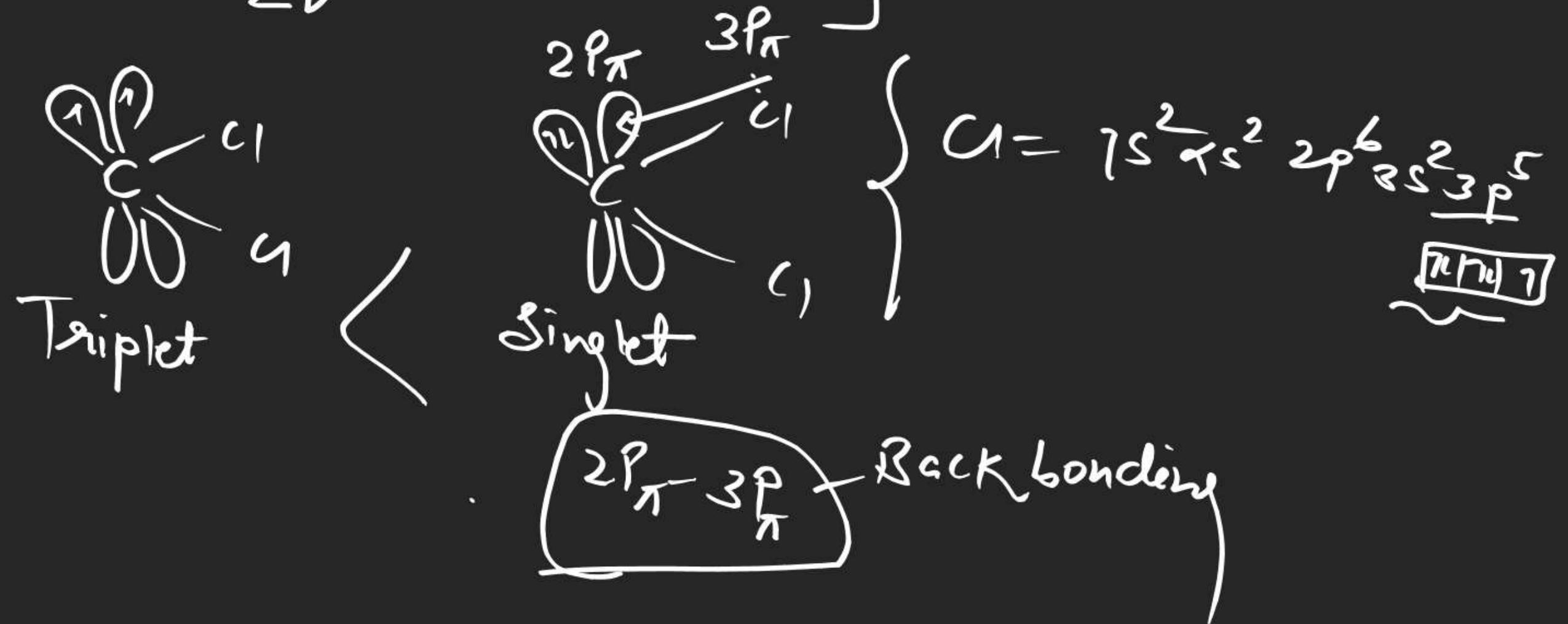
$2 \times 0 + 1$
1
Singlet

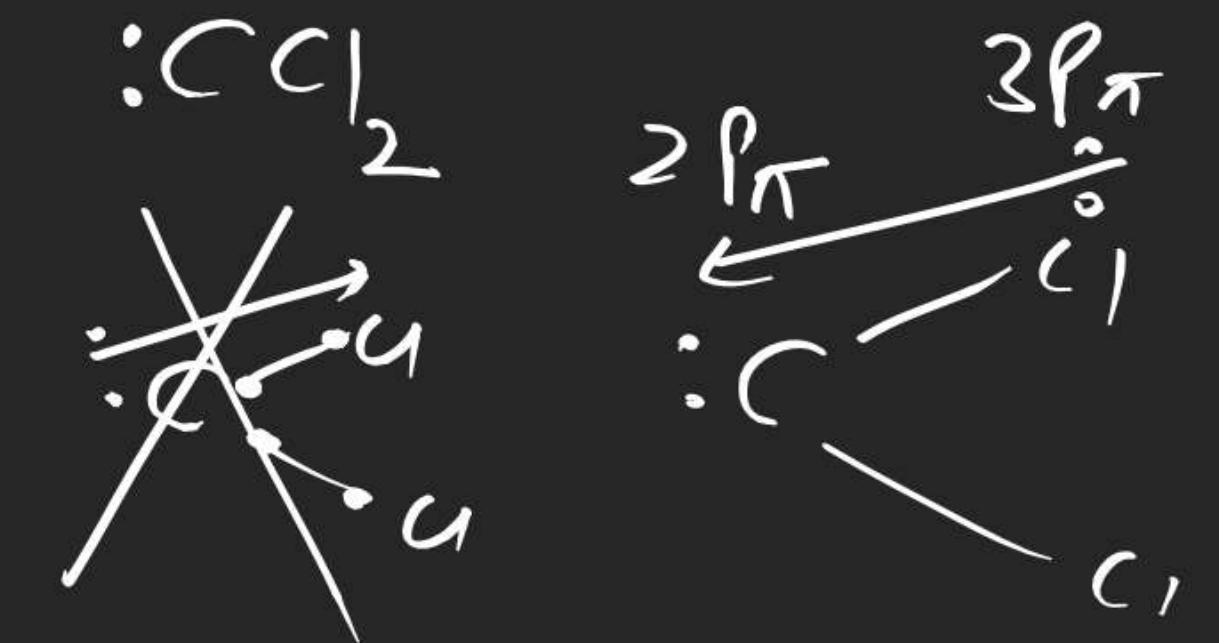
Note

Triplet Carbene is
more stable than

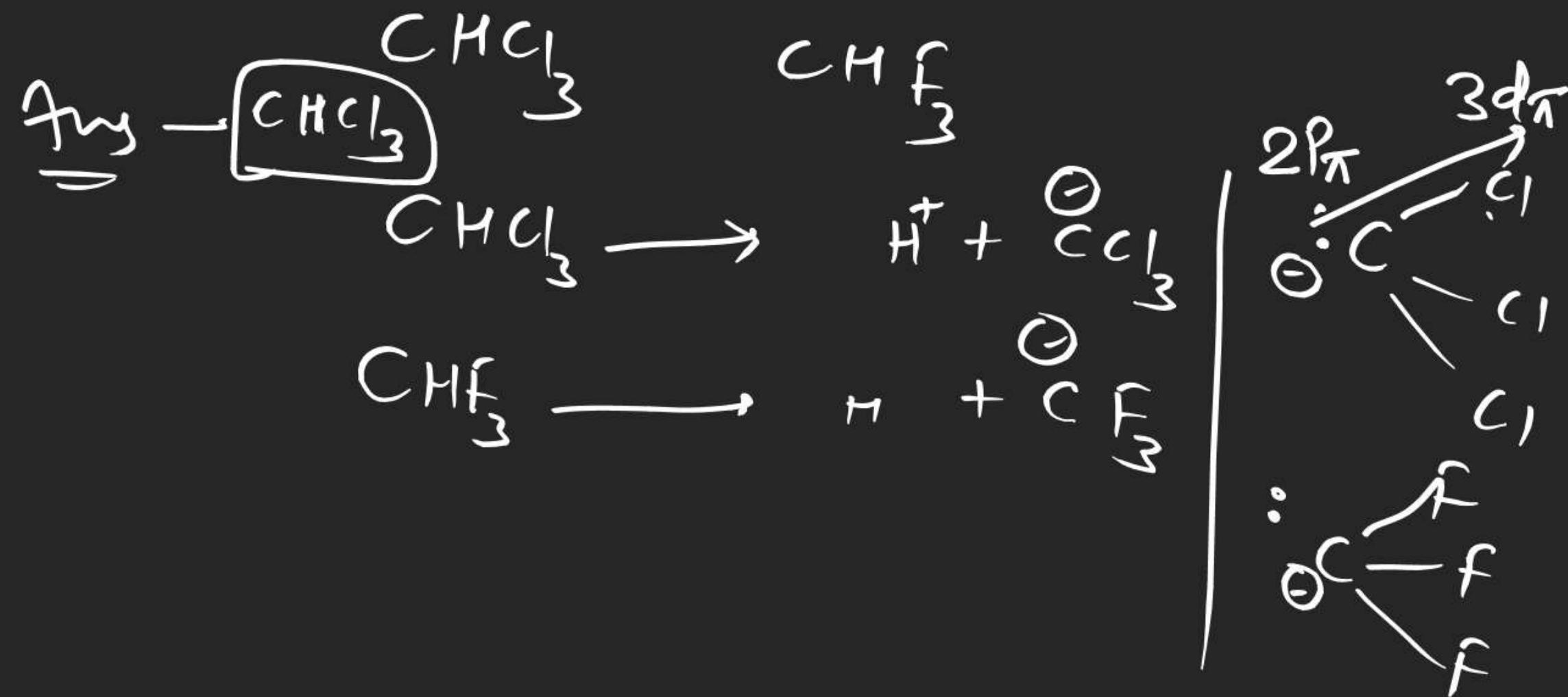
Singlet Carbene
due to Hund's Rule

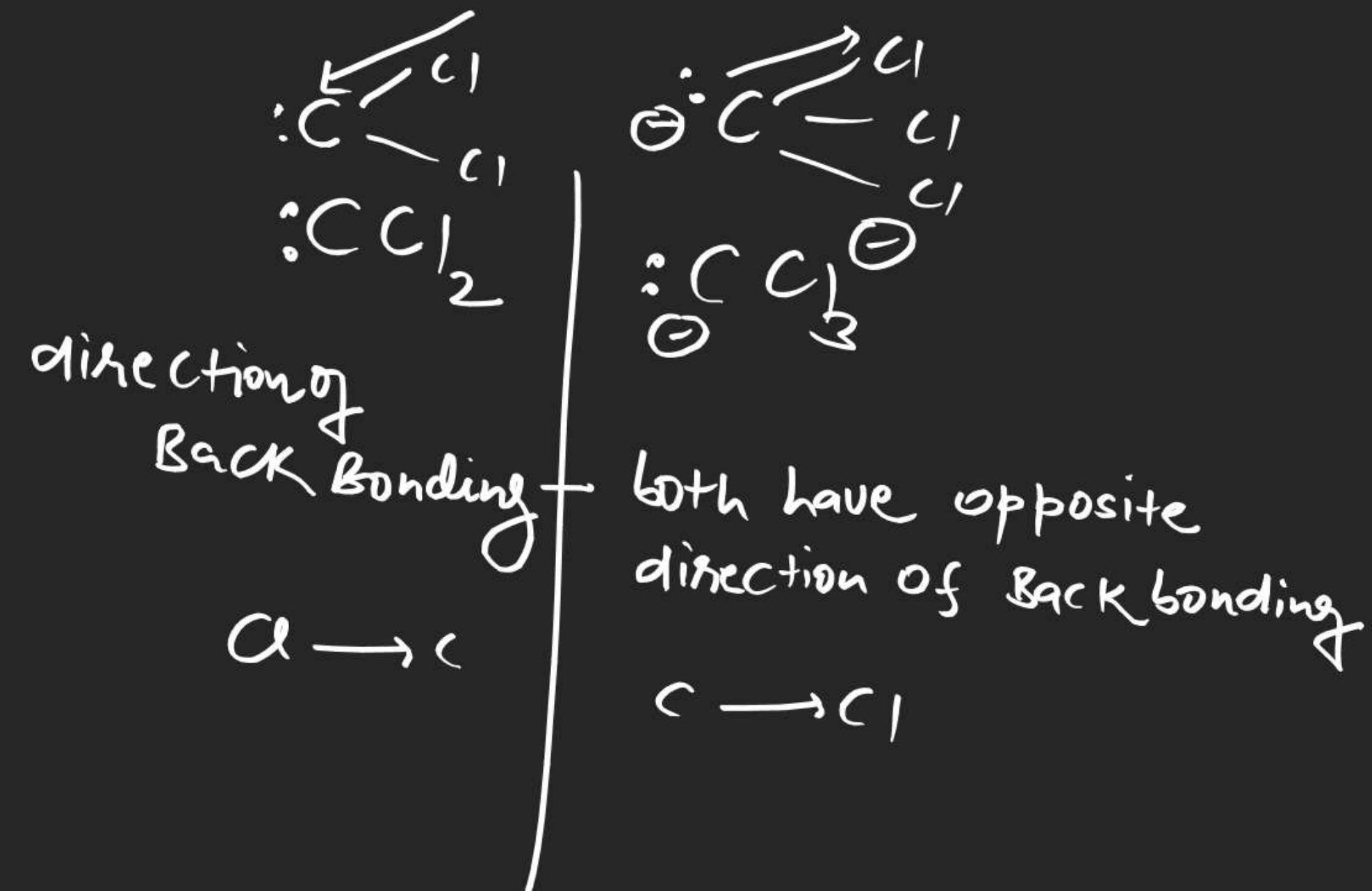
$$:\text{C}\text{Cl}_2 \left[\text{dichloro carbene} \right]$$

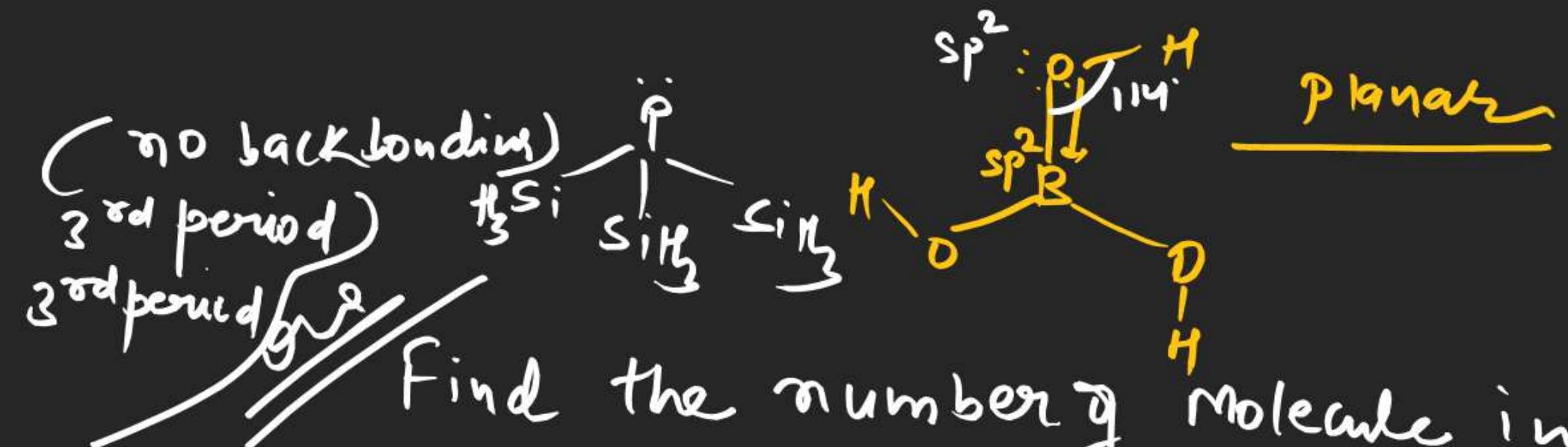




one which is better proton donor acid

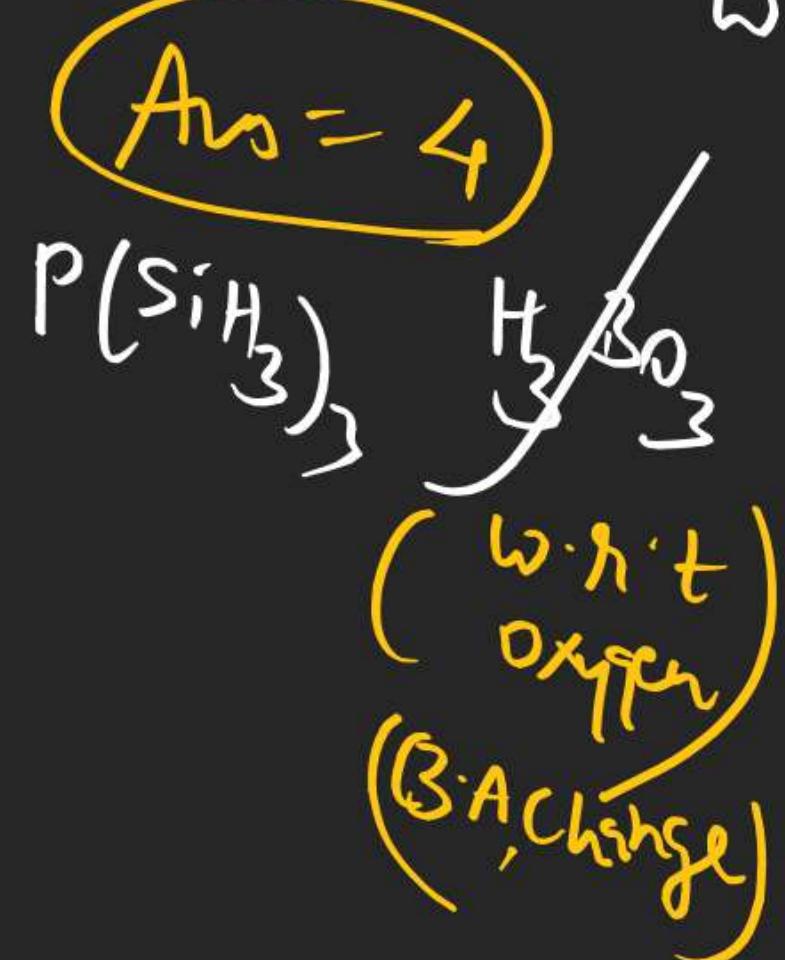


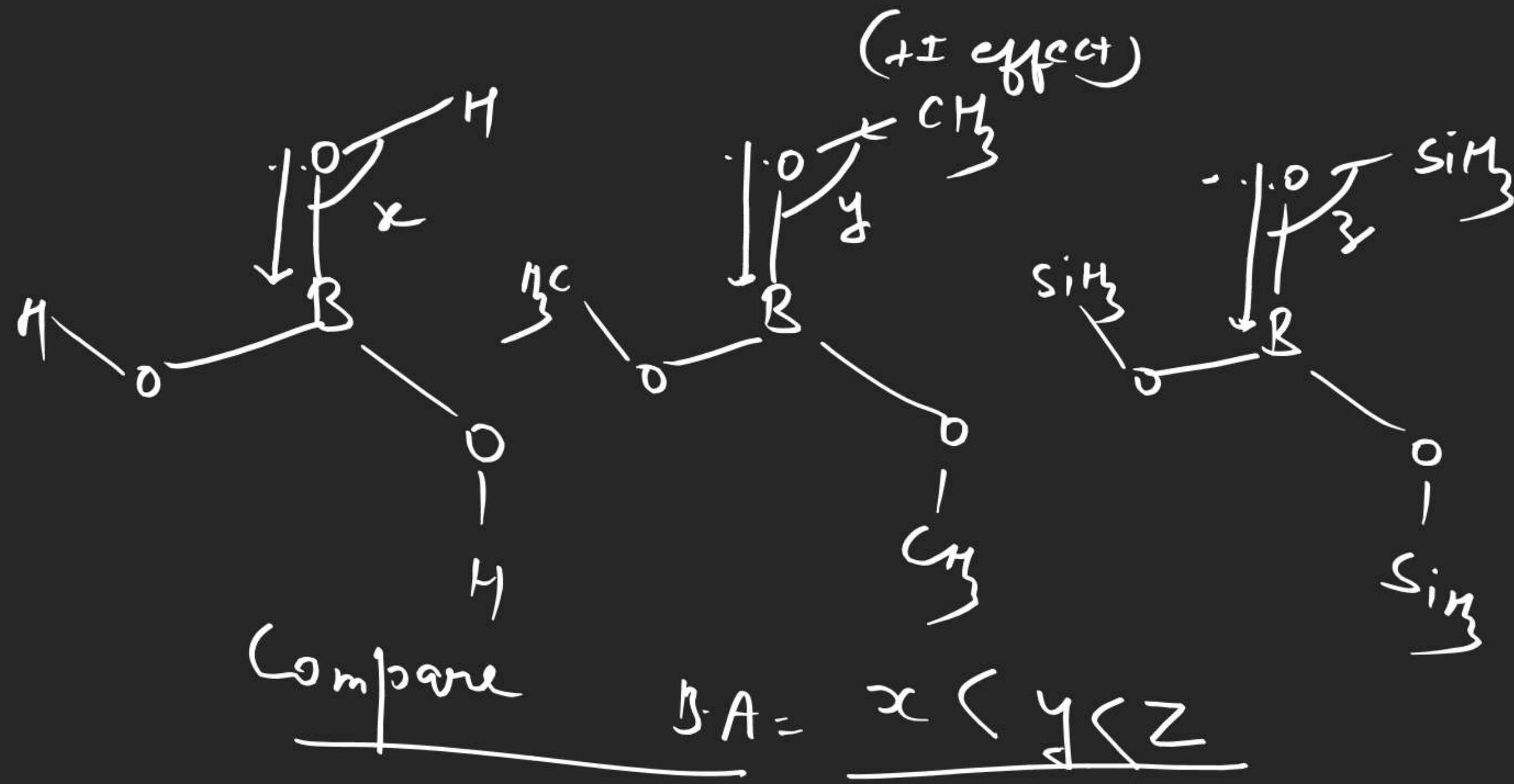


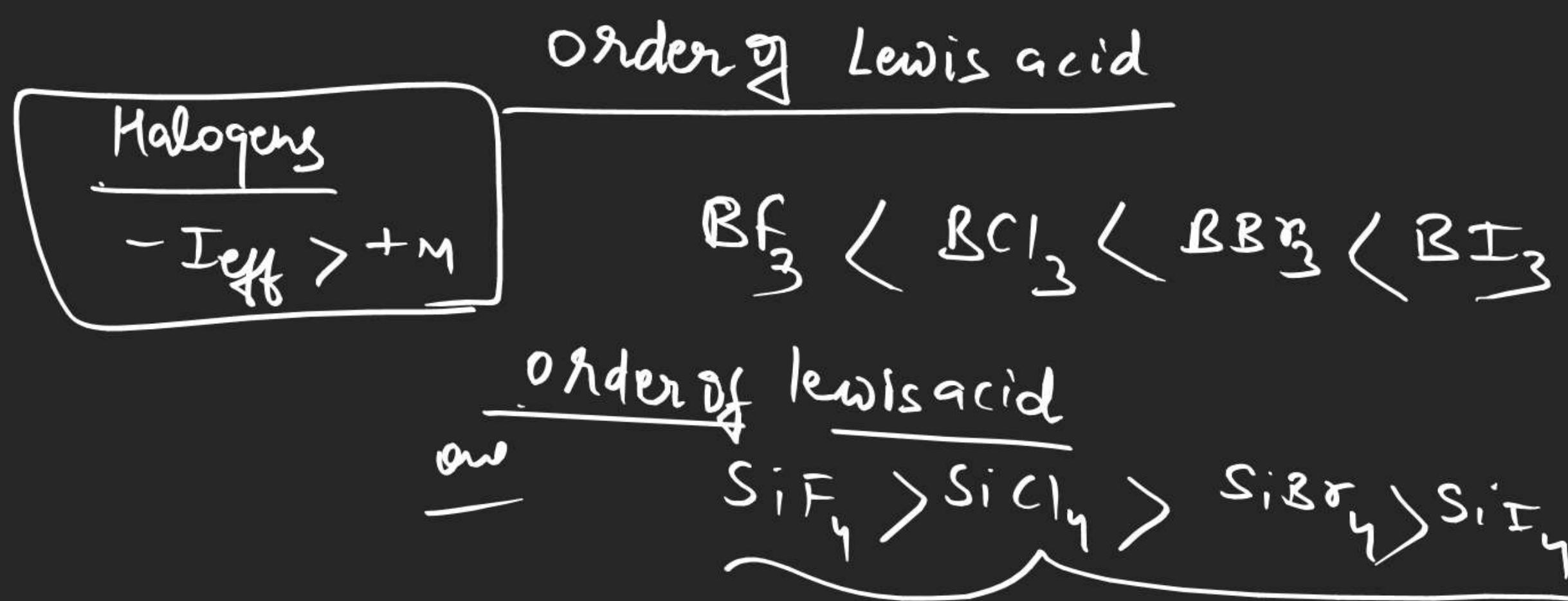


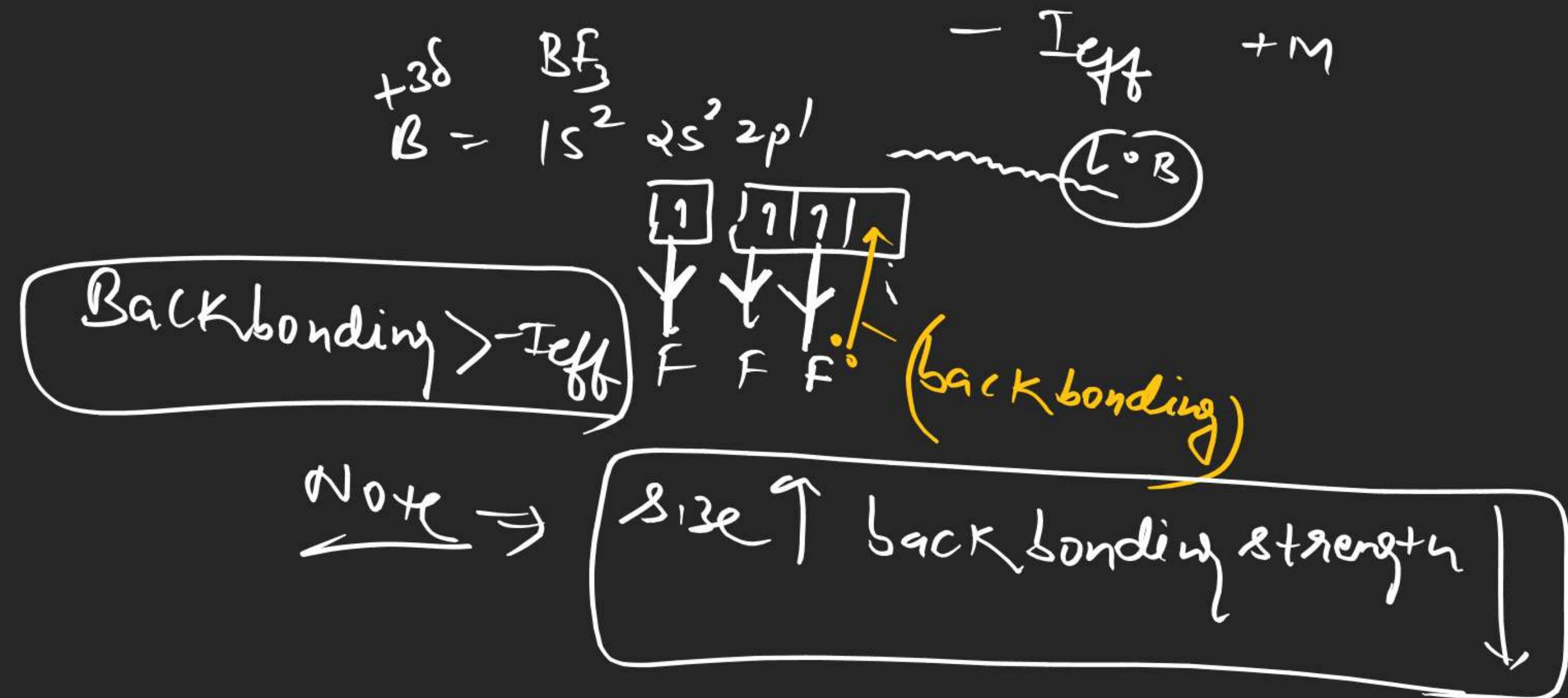
Find the numbering molecule in

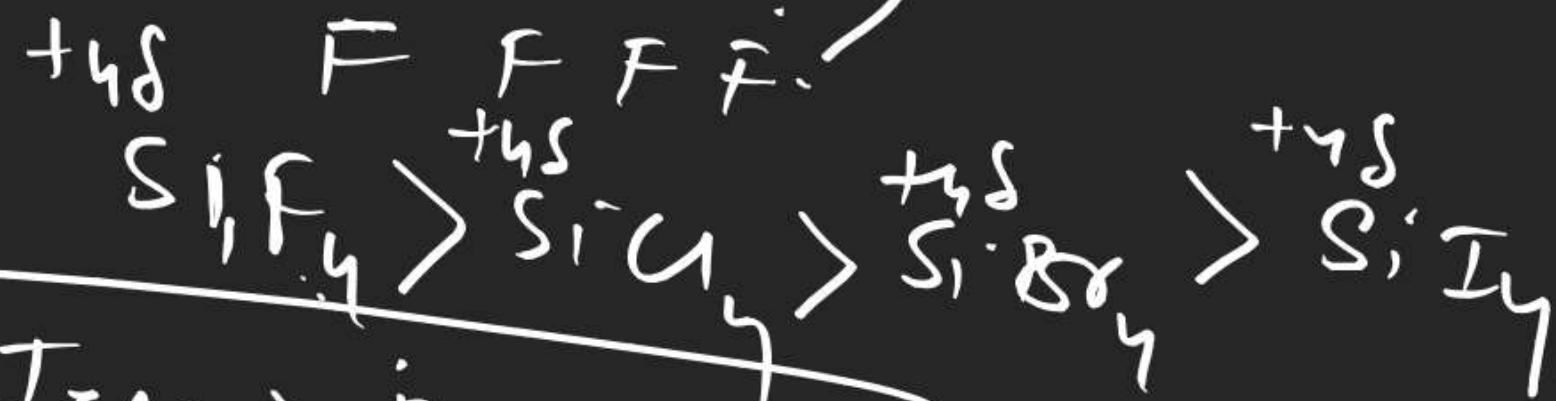
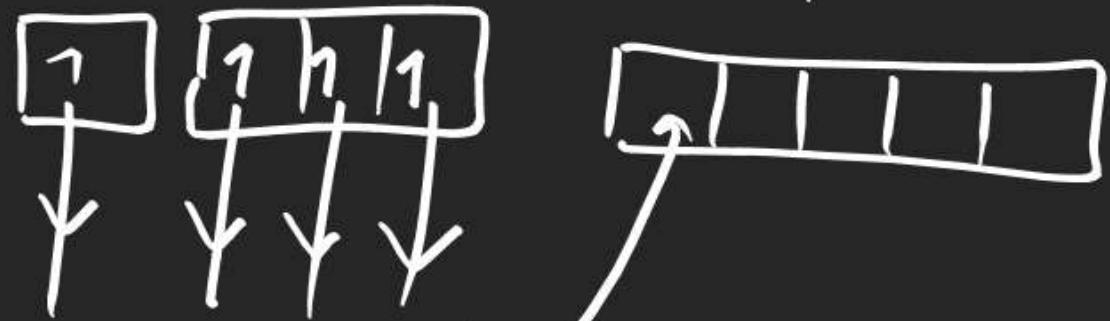
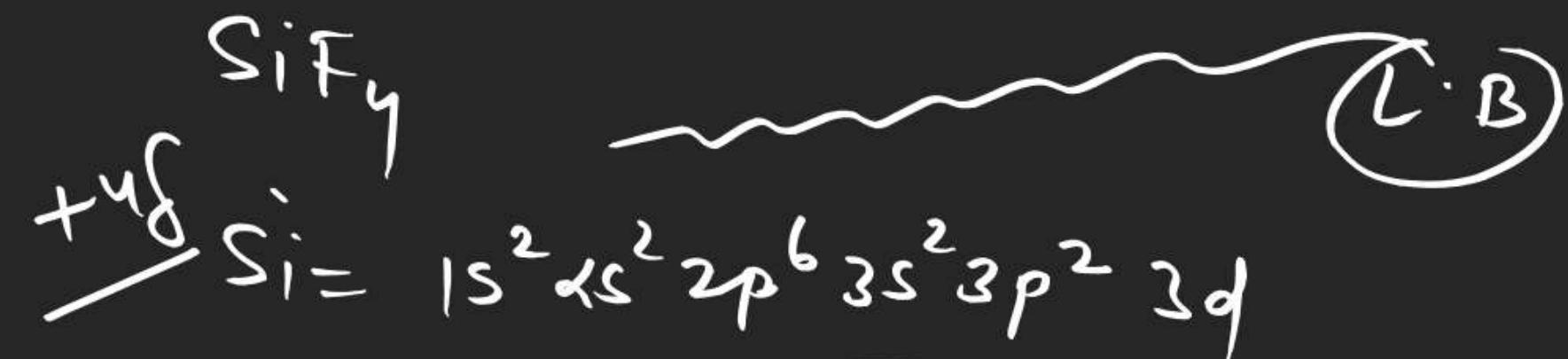
which $B \cdot A$ charge due to
Back bonding w.r.t \textcircled{B} , O, N





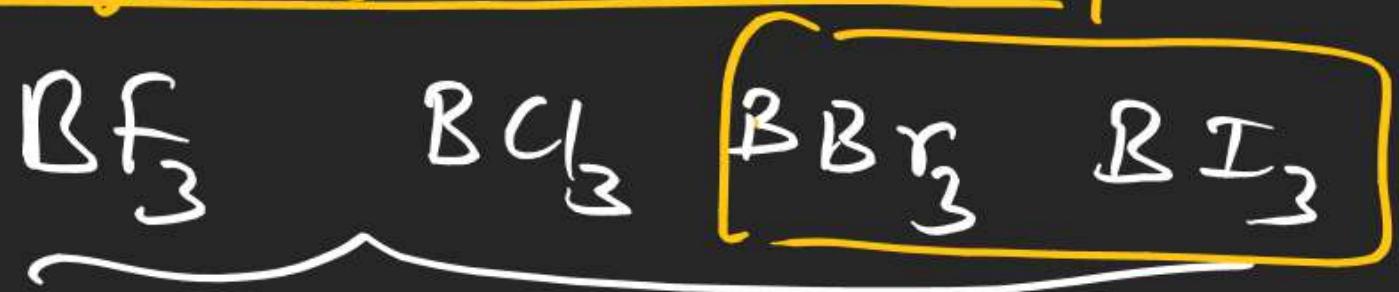






$-I_{eff} > \text{Back bond}$

Strength of back bonding



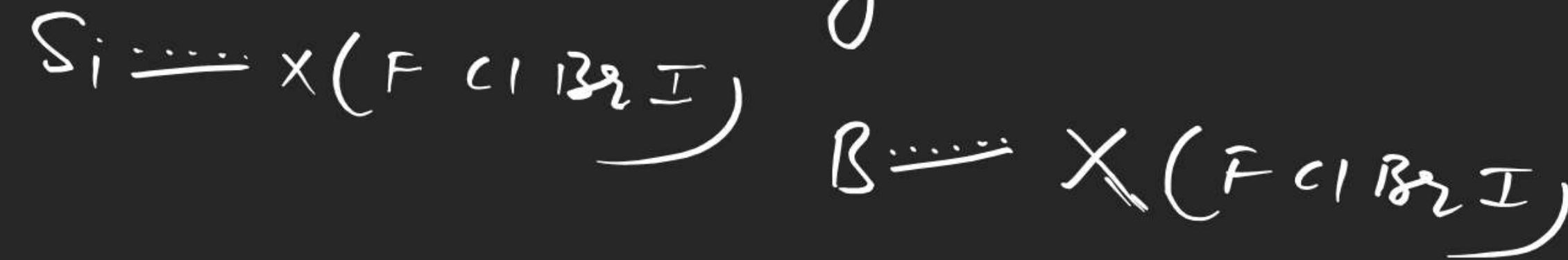
tri Halides of boron

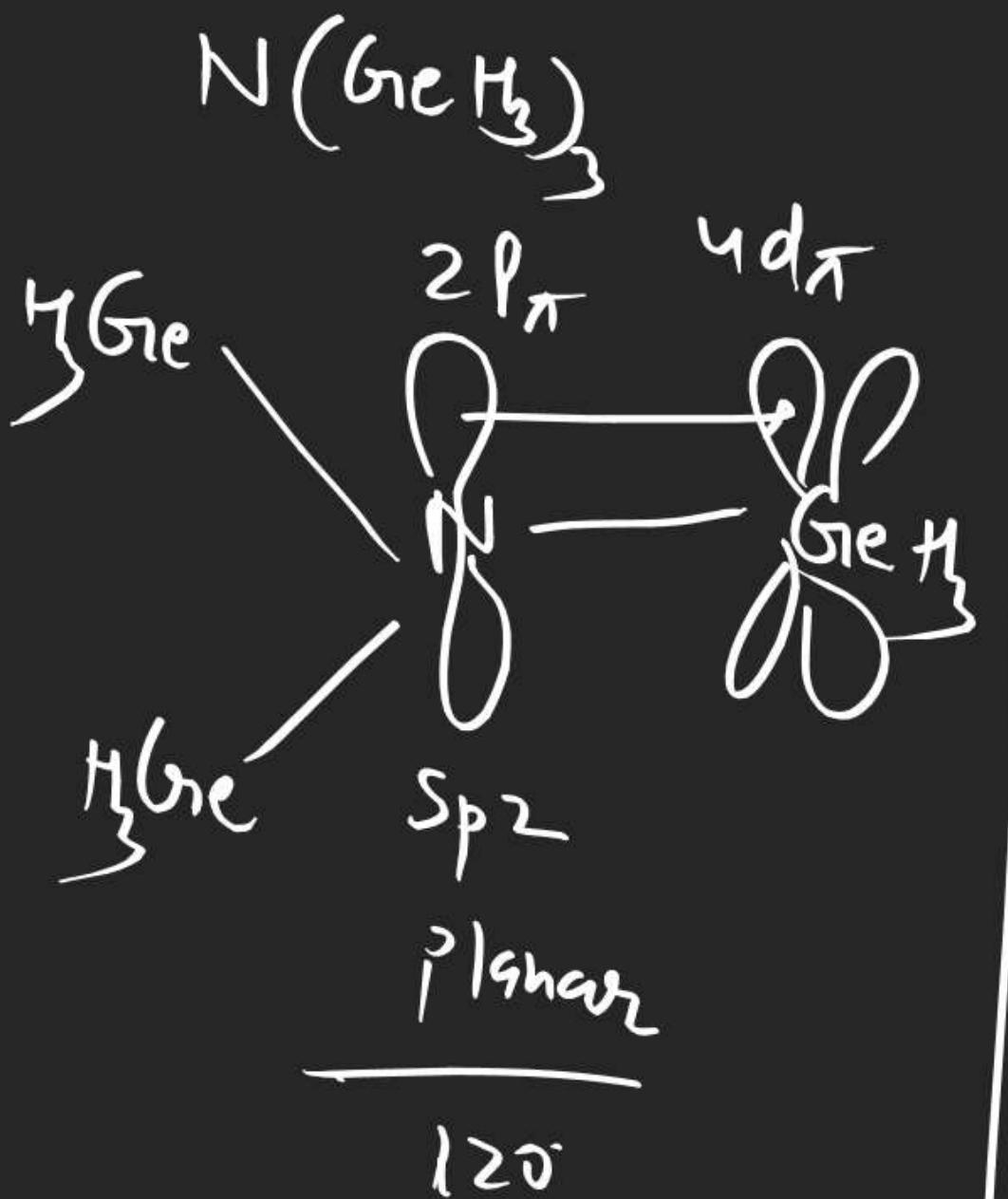
Note \Rightarrow Boron tri Halides have partial double bond character due to

Strength Back bonding

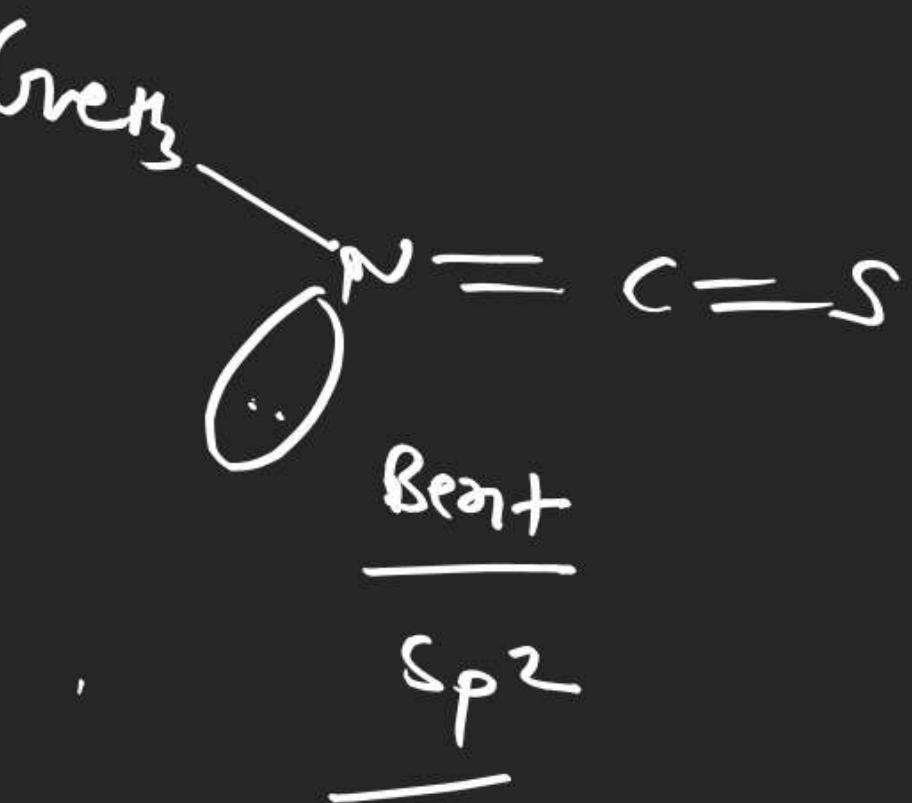


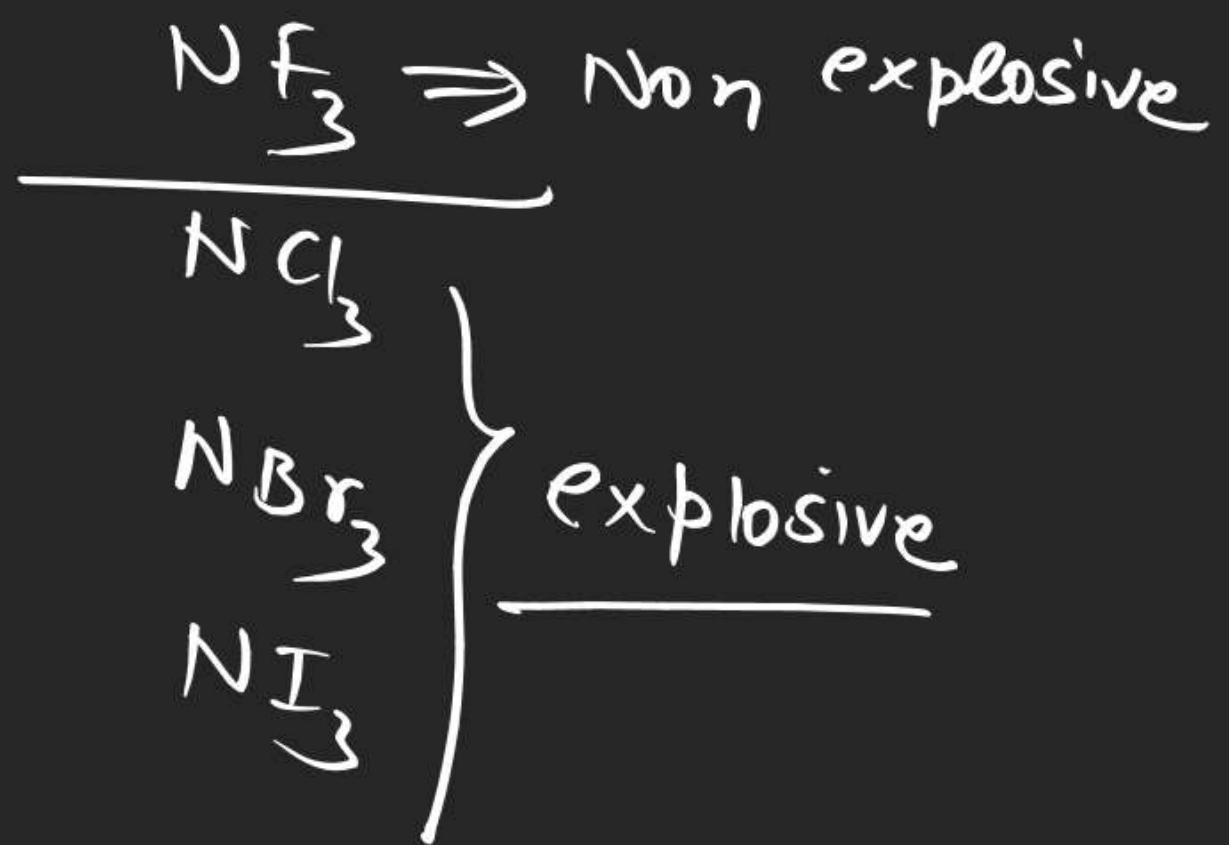
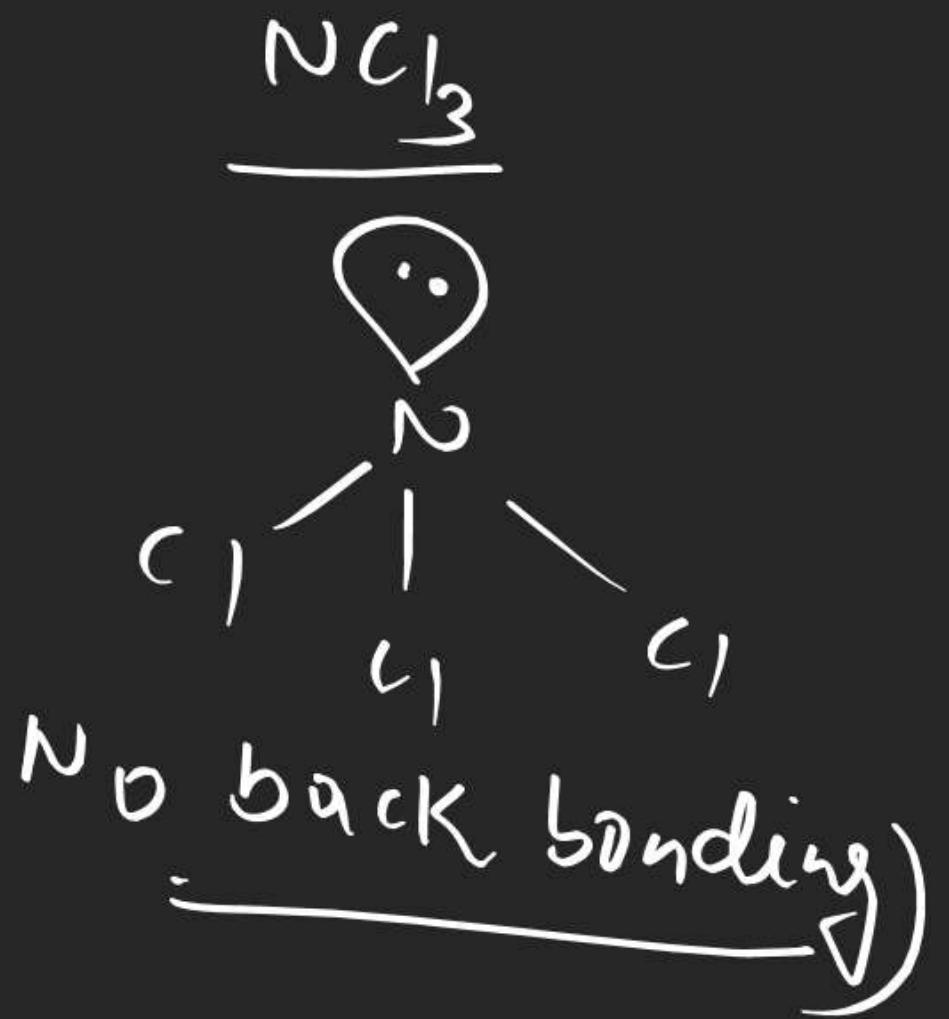
Silicon tetrahalide have partial double bond character due to back bonding

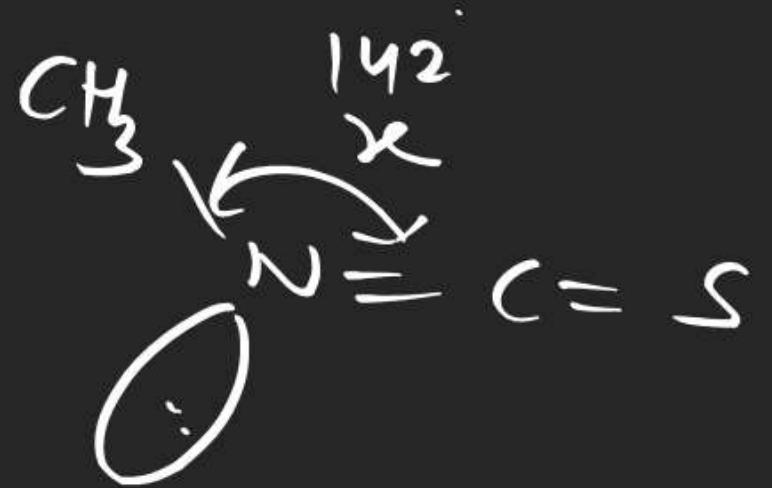




Get H_3 - NCS Draw the structure





Most important question

one $\beta\text{-A}$ x will be

~~A~~ $x > 120^\circ$

B) $x < 120^\circ$

C) $x = 120^\circ$

D) none

