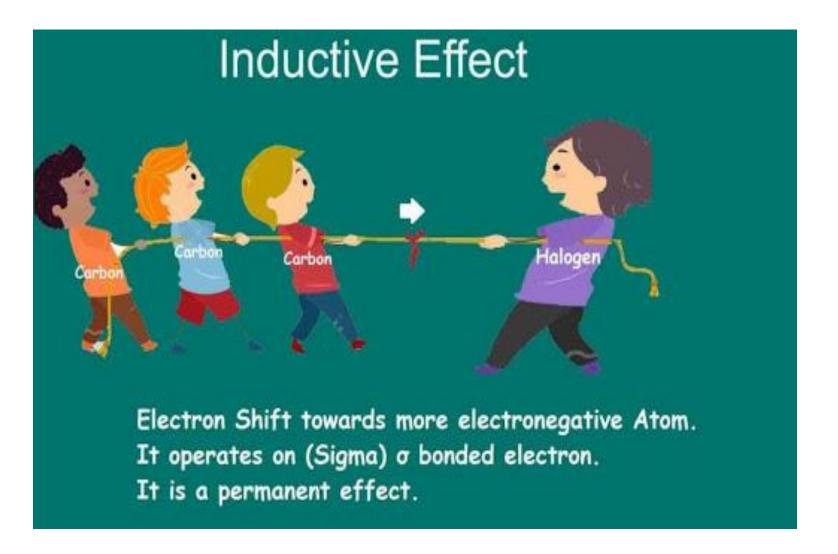
## **Inductive effect**

The effect of the sigma electrons displacement toward the highly electronegative atom is known as the inductive effect (I).







## **Types of Inductive Effect** System Group System Group More ∠Less Electronegative Electronegative Than system Than system -I effect +I effect "Y" is called + I effecting group "X" is called (- I effecting group



#### -I Effect

When an <u>electronegative atom</u>, such as a halogen, is introduced to a chain of carbon atoms the resulting unequal sharing of electrons generates a positive charge which is transmitted through the chain.

$$\delta \delta^{+} \qquad \delta^{-} \qquad \delta^{-}$$
 $CH_{3} \longrightarrow CH_{2} \longrightarrow Cl$ 



As the number of electronegative atoms or groups increase; the -I effect increase

(II) 
$$C1 \leftarrow CH_2 \leftarrow C \leftarrow \overset{\oplus \delta}{O} - H$$

(II)  $C1 \leftarrow CH \leftarrow C \leftarrow \overset{\oplus 2\delta}{O} - H$ 

(III)  $C1 \leftarrow C \leftarrow \overset{\oplus 3\delta}{C} - H$ 

the acidity order for the above compounds would be, III > II > I.



## +I Effect

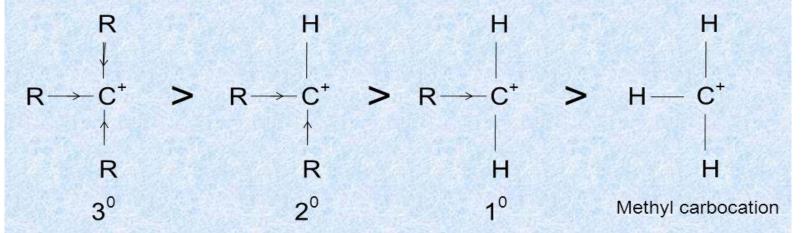
When a group tends to donate electrons, such as an <u>alkyl group</u>, is introduced to a carbon chain, the charge is relayed through the chain and this effect is called +I Effect

$$_{H-C-O-H}^{O} > _{CH_3}^{O} \rightarrow _{C}^{O} \rightarrow _{O-H}^{\Theta \delta}$$
(II)



#### (a) Inductive effect

More the number of alkyl group on the carbon atom carrying the +ve charge, greater would be the dispersal of the charge and hence more stable would be the carbocation. Thus, the stability of the carbocations decereases in the order: 3°> 2°> 1°>;



Stability decreases as +I-effect of the alkyl group decreases

$$(CH_3)^{\stackrel{+}{C}} > CH_3^{\stackrel{+}{C}}HCH_3 > CH_3^{\stackrel{+}{C}}H_2^{\stackrel{+}{C}} > CH_3^{\stackrel{+}{C}}H_2^{\stackrel{+}{C}} > CH_3^{\stackrel{+}{C}}H_3^{\stackrel{+}{C}}$$



## **Mesomeric effect**

It is defined as the polarity produced in the molecule by the interaction of two pi bonds or between a pi bond and lone pair of electrons present on an adjacent atom.



The mesomeric effect is negative (**-M**) when the substituent is an electron-withdrawing group and the effect is positive (**+M**) when the substituent is an electron releasing group.

#### +M EFFECT ORDER:

$$-0^{-} > -NH_{2} > -NHR > -OR > -NHCOR > -$$
  
OCOR > -Ph > -F > -Cl > -Br > -I

#### -M EFFECT ORDER:

$$-NO_2 > -CN > -S(=0)2-OH > -CHO > -C=O >$$
 $-COOCOR > -COOR > -COOH > -CONH_2 > -COO^-$ 



## +M effect

$$CH_{2} = CH - CH = CH - CH = CH - CH = MH_{2}$$

$$CH_{2} = CH - CH = CH - CH = MH_{2}$$

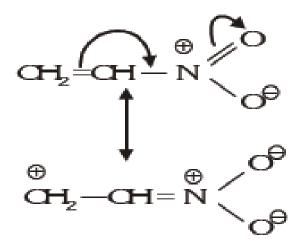
$$CH_{2} = CH - CH = CH - CH = MH_{2}$$

$$CH_{2} = CH - CH = CH - CH = MH_{2}$$

$$CH_{2} - CH = CH - CH = CH - CH = MH_{2}$$



## -M effect

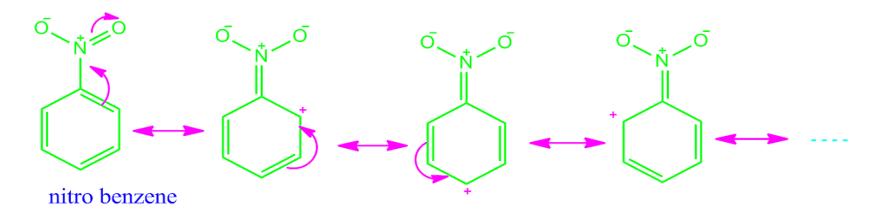




#### + M effect:

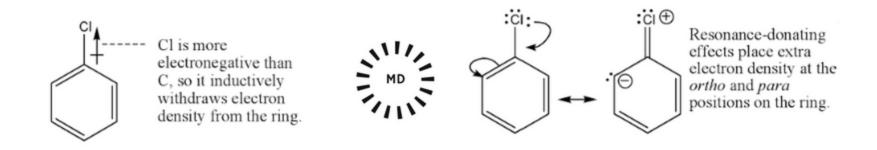


#### - M effect:





#### Inductive Effect vs Resonance Effect

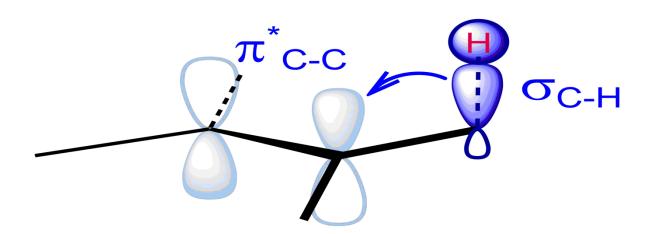


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# Hyperconjugation

It the interaction of the electrons in a  $\sigma$ -bond (usually C-H) with an adjacent empty or partially filled porbital or a  $\pi$ -orbital





# Hyperconjugation empty 2p orbital of carbon H H



$$H \xrightarrow{H} CH \xrightarrow{C} CH_2 \rightleftharpoons H \xrightarrow{C} CH \xrightarrow{C} CH_2 \rightleftharpoons H \xrightarrow{C} CH \xrightarrow{C} CH_2 \rightleftharpoons H \xrightarrow{C} CH \xrightarrow{C} CH_2 \rightleftharpoons H \xrightarrow{H} H \xrightarrow{H} H \xrightarrow{H} C \xrightarrow{C} CH = \overline{C}H_2$$

$$(I) \qquad \qquad H \xrightarrow{H} C \xrightarrow{C} CH = \overline{C}H_2$$

$$(IV)$$



# Hyperconjugation: No bond resonance

- ▶ The electrons of the sigma bond between C and H are involved in delocalization.
- ► In structure to the right: No bond between C and H due to migration of the sigma bond.

  Hence Hyperconjugation is also called as 'NO BOND RESONANCE'.
- ➤ This does not indicate that hydrogen is completely detached from the structure, but some degree of ionic character in the C – H bond and some single bond character between carbon – carbon double bond.

