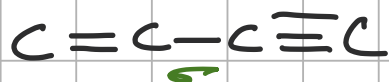


\* Polyenes →  $\beta$  carotene

→ lycopene

multiple bond

\* Conjugated



\* non conj



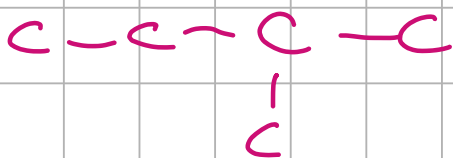
cumulated



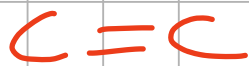
\* more than double bond

→ diene

triene



iso



tetrahedral

trigonal

Linear

relatively free

restricted

$180^\circ$

$109.5^\circ$

$\pi$  bond

planar

"Acetylenes"

$(sp^3)$

$120^\circ$

$(sp)$

$(sp^2)$

nucleophilic



electrophilic



\* Hydrogenation = syn addition  
+ H

# Alkynes

→ Hydrogenation (+ H)

Catalyst Ni/Pt → —

special palladium (Pd)  
= (Lindlar's catalyst) → 1 mol H<sub>2</sub>  
Cis alkene

Na/Li → trans alkene

→ Halogenation (+ X<sub>2</sub>)

First trans →  $\begin{array}{c} \text{Cl} & & \text{Cl} \\ | & & | \\ \text{C} & - & \text{C} \\ | & & | \\ \text{Cl} & & \text{Cl} \end{array}$

→ Hydrogen Halide

Markovnikov's Rule

H  $\rightarrow$   $\text{C}(\text{H})_3$  is H  $\neq$

# Alkenes \* Electrophilic addition.

→ Hydrogenation (+H)

syn addition (same side)

cis

→ Halogenation (+X<sub>2</sub>)

Anti addition (opposite side)

trans

Cl<sub>2</sub> ✓

Br<sub>2</sub> ✓

I<sub>2</sub> unreactive

F<sub>2</sub> too reactive

→ Addition of acids

HF / HCl / HBr / H-O-SO<sub>3</sub>H

→ Oxidation → Ozonolysis O<sub>3</sub>

→ \*two carbonyl groups:  $\begin{matrix} \text{O} \\ \parallel \\ \text{C} - \text{H} \end{matrix}$

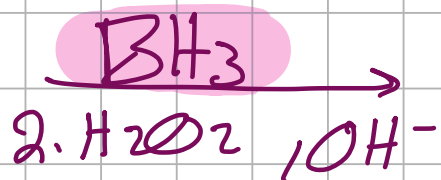
purple → brown-black (MnO<sub>2</sub>)

permanganate

KMnO<sub>4</sub>

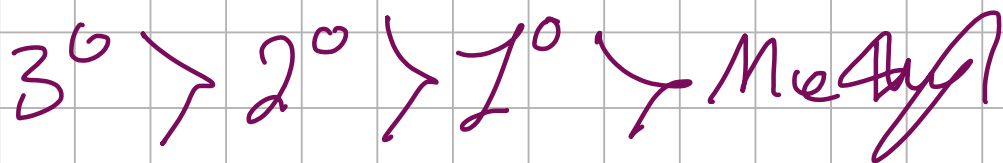
cis

→ aglycol  
⇒ cis-cyclopentane-1,2-diol



anti M.V

!!  $\text{C}^\ominus$  &  $\text{H}^\oplus$



\* most stable

\* least stable

# Alkanes

physical properties of —  
||  
|||

\* physical states

$C_1 \rightarrow C_4$  gases  
 $C_5 \rightarrow C_{17}$  liquids  
 $C_{18} \rightarrow$  wax like solids

\* Solubility  $\Rightarrow$  "Like dissolve like"

\* Alkane  $/= / \equiv$  )  $\rightarrow$  Soluble in nonpolar solvents

$\rightarrow$  Carbon tetrachloride  
 $CCl_4$   
 $\rightarrow$  benzene.

Water X

\* Boiling points B.P

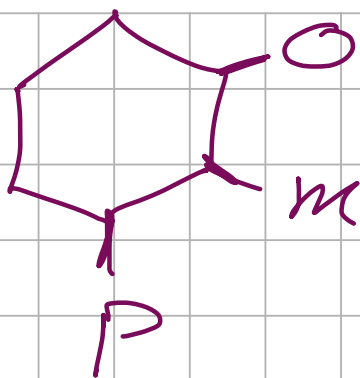
→ more branch less B.P

\* Intermolecular interaction

→ H-bond  
→ dipole-dipole  
→ London.

B.P → ↑ increases — ↑ M. weight  
→ ↓ decrease — ↑ branches





\*  $M \Rightarrow$  withdraw  $e^-$  from the ring  
 $\hookrightarrow$  ring deactivating groups

\* substituents withdraw  $e^-$

Decrease  $e^-$   
density

slow down the  
reaction.

\*  $O, P \Rightarrow$  supply  $e^-$  to the ring.

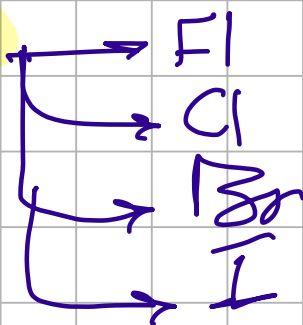
$\hookrightarrow$  ring activating

\* substituents donate  $e^-$

increase  $e^-$   
density and

speed up the  
reaction

\* Halogens



withdrawing  $e^-$  (and they are strong w.d)

But (X) unshared electrons pairs so they are

Ortho, para-directing

\* X<sub>2</sub> ⇒ deactivating.

exception