

# General Chemistry.1

The title of Lecture: Alkenes and Alkynes

Year One; Semester. 1

Presented by Dr. Sameer



#### Alkenes (Olefins):

Carbon-hydrogen combination of CnH2n. Alkene family consists of chain molecules that contain double carbon-to-carbon bond and are therefore unsaturated.

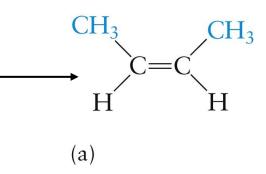
- ☐ Also known as olefins, aliphatic, unsaturated
- ☐ C=C double bonds
- $\square$  Formula for one double bond =  $C_nH_{2n}$ .
- ☐ Cis-trans isomerism ←

**ethene** = **ethylene** 

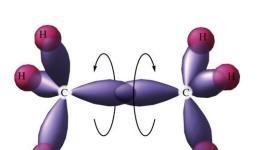
**propene** = **propylene** 

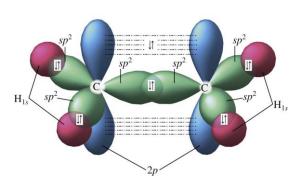
$$H$$
 $C = C$ 
 $H$ 

$$\begin{array}{c} H \\ C = C \\ H \end{array}$$



$$CH_3$$
  $C=C$   $CH_3$ 





(b)

### **Alkynes**

Alkynes: hydrocarbons containing at least one carbon—carbon triple bond.

General formula is: CnH2n-2

Begins with ethyne (acetylene) 5. For Example: C2H2 H—C=C—H

- 1. Identify the parent chain, which should include the C≡C triple bond
- 2. Identify and Name the substituents
- 3. Assign a locant (and prefix if necessary) to each substituent giving the C≡C triple bond the lowest number possible
- 4. List the numbered substituents before the parent name in alphabetical order. Ignore prefixes (except iso) when ordering alphabetically
- 5. The C≡C triple bond locant is placed either just before the parent's name or just before the -yne suffix

TABLE 20.7 Alkenes							
n	Name	Molecular Formula C <sub>n</sub> H <sub>2n</sub>	Structural Formula	Condensed Structural Formula			
2	Ethene	$C_2H_4$	H $C = C$ $H$	$CH_2 = CH_2$			
3	Propene	$C_3H_6$	$\begin{array}{c} H \\ \downarrow C = \begin{array}{c} H \\ \downarrow \\ - \\ \downarrow \\ - \\ H \end{array} $	CH <sub>2</sub> == CHCH <sub>3</sub>			
4	1-Butene*	C <sub>4</sub> H <sub>8</sub>	$\begin{array}{c c} H & C = C - C - C - H \\ \downarrow & \downarrow & \downarrow \\ H & H \end{array}$	$CH_2 = CHCH_2CH_3$			
5	1-Pentene*	C <sub>5</sub> H <sub>10</sub>	$\begin{array}{c ccccccccccccccccccccccccccccccccccc$	$CH_2 = CHCH_2CH_2CH_3$			
6	1-Hexene*	C <sub>6</sub> H <sub>12</sub>	$\begin{array}{c ccccccccccccccccccccccccccccccccccc$	$CH_2 = CHCH_2CH_2CH_2CH_3$			

<sup>\*</sup> These alkenes have one or more isomers depending on the position of the double bond. The isomers shown here have the double bond in the 1 position, meaning the first carbon–carbon bond of the chain.

# of Carbons	Prefix	Suffix	Structural Formula	Molecular Formula
2	Eth	yne	H-C=C-H	CzHz
3	Prop	yne	H-C=C-C-H	C3H4
4	But	yne	H-C=C-C-H	C4H6
5	Pent	yne	H-C=C-C-H H-T-C-C-H	C5H8
۵	Hex	yne	H-C=C-C-C-C-H	Cenio
7	Hept	yne	M H H H H H H H H H H H H H H H H H H H	C <sub>4</sub> H <sub>12</sub>
8	Oct	yne	H-CEC-C-C-C-H	C8 H14

#### **Physical Properties of Alkenes**

#### o **Physical State**

Alkenes and alkynes occur at room temperature are gases, liquids, and solids.

- C2 to C4 are gases,
- C5 to C17 are liquids,
- C18 and larger alkenes are wax —like solids.

#### o Solubility

- Alkenes are nonpolar compounds.
- Alkenes are soluble in the nonpolar solvents; CCl<sub>4</sub> and benzene,
- Alkenes are insoluble in polar solvents like water.

#### Boiling Points & Melting Points

■ The boiling points and melting points of normal hydrocarbons increase with increasing molecular weight.

#### Nomenclature of Alkenes and Alkynes

☐ To name alkenes, select the longest carbon chain which **includes** the carbons of the double bond. Remove the **-ane** suffix from the name of the alkane which corresponds to this chain. Add the suffix **-ene**.

#### The IUPAC Rules

The IUPAC rules for naming alkenes and alkynes are similar to those for alkanes, but a few rules must be added for naming and locating the multiple bonds.

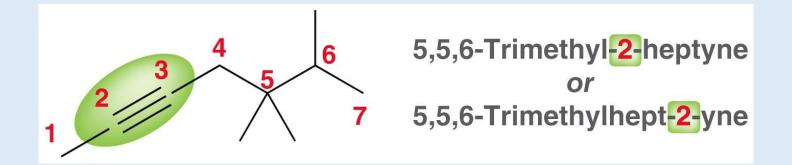
- 1. The ending *-ene* is used to a carbon–carbon double bond. The ending *-yne* is used for a carbon–carbon triple bond
- 2. Select the longest chain that includes both carbons of the double or triple bond.
- 3. Number the chain from the end nearest the double or triple bond so that the carbon atoms in that bond have the lowest possible numbers.

TABLE 12.7 Comparison of Names for Alkanes, Alkenes, and Alkynes

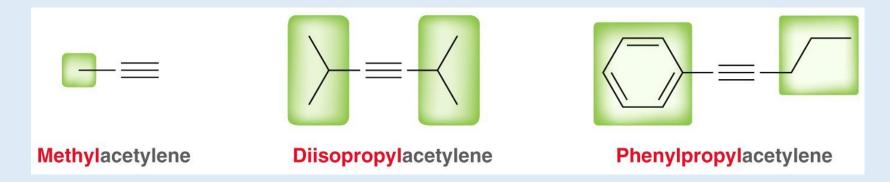
Alkane	Alkene	Alkyne
CH <sub>3</sub> —CH <sub>3</sub>	$H_2C = CH_2$	НС≡СН
Ethane	Ethene (ethylene)	Ethyne (acetylene)
$CH_3-CH_2-CH_3$	$CH_3-CH=CH_2$	$CH_3-C\equiv CH$
		_=
Propane	Propene	Propyne

#### **Nomenclature of Alkenes and Alkynes**





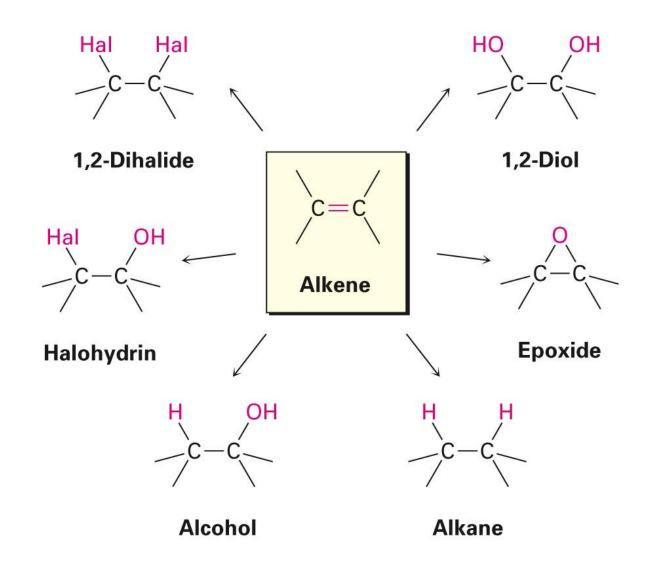
• common names derived from **acetylene** are often used as well



• Alkynes are also classified as **terminal** or **internal** 



# Reactions of Alkenes



#### **Reactions of alkenes**

☐ Heat of hydrogenation

✓ the heat liberated during this reaction.  $\Delta H$  is ~125 kJ/mol for each double bond in the compound.

☐ Addition of hydrogen halides

$$HX = HCI, HBr, HI$$

 $H_2C=CH_2 + HCI \longrightarrow CH_3CH_2CI$ only one product is possible, chloroethane but....

$$CH_3$$
- $CH$ = $CH_2$   $CH_3$ - $CH$ - $CH_3$   $CI$ - $CI$ 

Only 2-chloropropane is formed

#### Markovnikov's rule

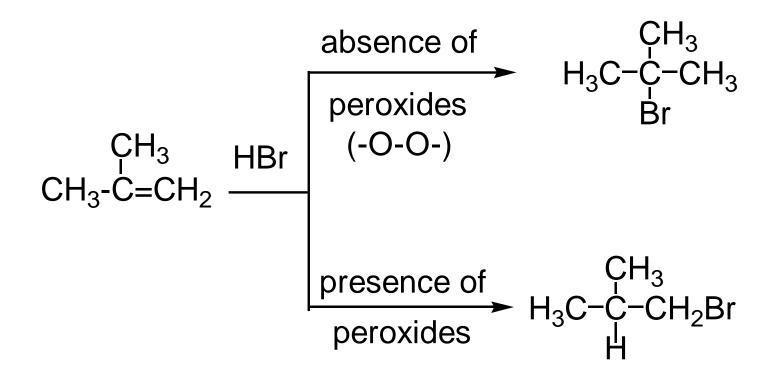
$$H \rightarrow CH_3 \rightarrow H_3C-C-CH_3$$
 $H \rightarrow CH_3 \rightarrow CI$ 

☐ In 1869, Markovnikov proposed that in the addition of an acid to an alkene, the hydrogen of the acid bonds to the carbon which is already bonded to the greater number of hydrogens.

Each carbon of the double bond is bonded to one H therefore both isomers are formed.

# HBr - the peroxide effect

## 1933, Kharasch and Mayo



#### **Addition of sulfuric acid**

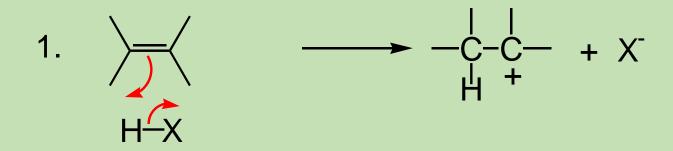
$$CH_{3}CH=CH_{2} \xrightarrow{\text{Cold} \atop \text{H}_{2}SO_{4}} \xrightarrow{\text{CH}_{3}CHCH_{3}} \xrightarrow{\text{CH}_{3}CHCH_{3}} \xrightarrow{\text{DSO}_{3}H} \xrightarrow{\text{DSO}_{3}H} \xrightarrow{\text{CH}_{3}CHCH_{3}} \xrightarrow{\text{CH}_{3}CHCH_{3}}$$

### **Hydration**

$$H \rightarrow CH_3 + H_2O \rightarrow H^+ \rightarrow H^+C^-C^-CH_3$$
  
 $H \rightarrow CH_3 + H^+OH$ 

a Markovnikov addition

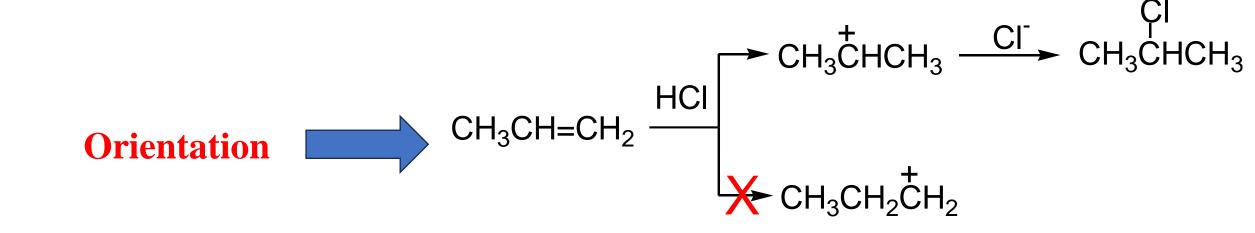
#### The mechanism of the addition



2. 
$$-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-\overset{|}{C}-$$

$$HX = HCI, HBr, HI, H2SO4, H3O+$$

# An example



# **Addition of halogens**

$$X_2 = CI_2$$
,  $Br_2$  usually iodine does not react

$$CH_3CH=CH_2 \xrightarrow{Br_2 \text{ in}} CH_3CHBrCH_2Br$$

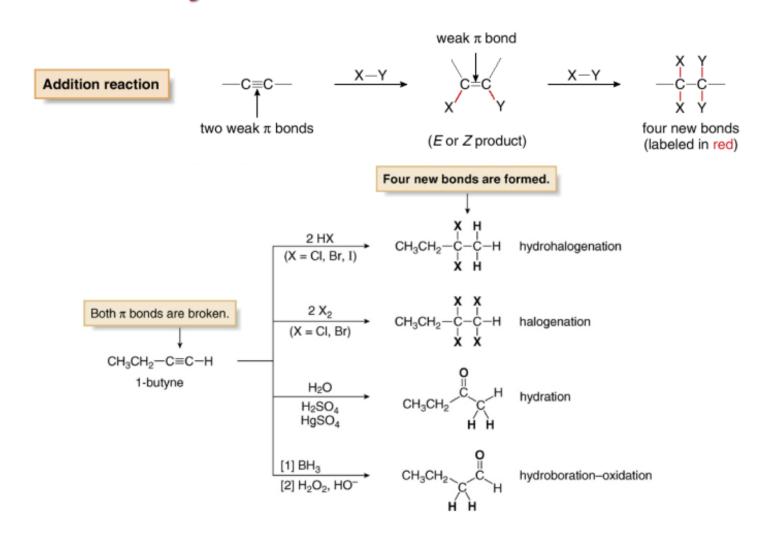
$$CCl_4$$

1,2-dibromopropane

# Mechanism of X<sub>2</sub> addition

# **Alkynes Reactions**

# **Alkyne Reactions—Additions**



# **Preparation of Alkynes**

- Like alkenes, alkynes can also be prepared by elimination
- Need a **dihalide** to make an alkyne

An alkyl halide

An alkyl dihalide

- Such eliminations usually occur via an E2 mechanism
- ☐ Geminal or vicinal dihalides can be used

Vicinal dihalide 
$$R-C-C-R$$
 Strong base  $R-C\equiv C-R$   $H$   $H$ 

# The End of Lecture