

# NEET - UG

# NATIONAL TESTING AGENCY

# Chemistry

**Organic Chemistry** 



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# Organic Chemistry

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Unleash the topper in you



# "General Organic Chemistry"

\* Urea was the first organic compound synthesized in laboratory by Wohler.

$$NH_4CNO \stackrel{\Delta}{\longleftarrow} H_2N \stackrel{O}{\longleftarrow} NH_2$$

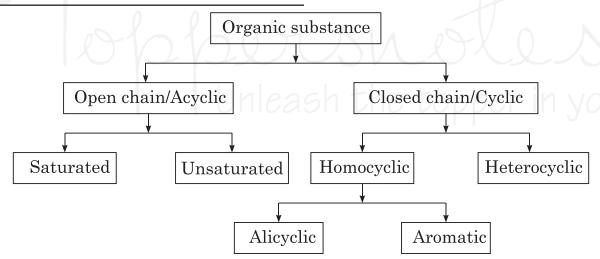
\* Hydrocarbons and their derivatives are called organic substances.

$$R$$
—H  $\xrightarrow{\Delta G}$   $R$ —G

Hydrocarbon Hydrocarbon derivative

# Classification of Organic Substances

# A. On the basis of carbon chain



#### 1. Open chain substances

(i) Saturated substances: Substances in which all the carbon atoms are joined by single bonds only.

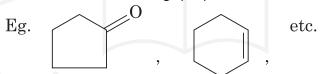
(ii) Unsaturated compounds: Substances having at least 1 C—C multiple bond.



- \* Multiple bond containing organic substances may be saturated or unsaturated.
- \* Olefinic bond: Carbon-carbon double bond is also called olefinic bond.
- \* Acetylenic bond : C—C triple bond is also called acetylenic bond.
- \* Open chain substances are also called aliphatic substances.

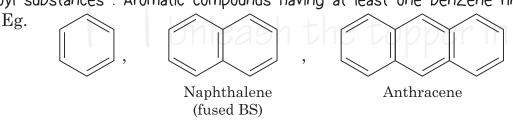
#### 2. Closed chain substances:

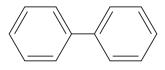
- (i) **Homocyclic substances**: [Carbocyclic substances] substances in which ring system is made up of same type of atoms.
  - (a) Alicyclic substances: Substances having properties similar to aliphatic substances.



#### (b) Aromatic substances:

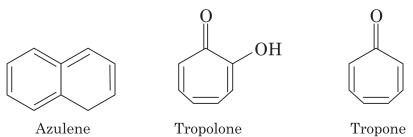
I. Benzenoyl substances: Aromatic compounds having at least one benzene ring.





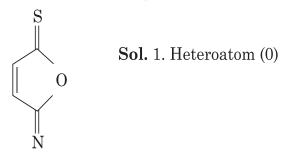
Diphenyl

II. Non-Benzenoyl substances: Aromatic substances without benzene ring.

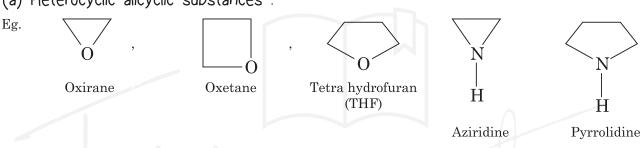




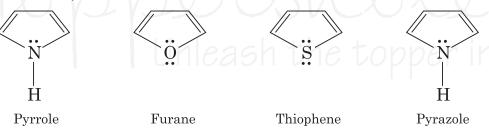
- (ii) Heterocyclic substances: Substances in which ring system having at least 1 different atom another than C atom, like N, S, etc.
  - Q. How many heteroatoms are present in given compounds?



(a) Heterocyclic alicyclic substances :



(b) Aromatic heterocyclic substances :



# B. Classification based functional group

#### 1. Homologous series:

 $_{\circ}$  Series of organic substances having same functional group and difference of at least 1 —CH  $_2$  unit 14 mol weight.

#### 2. Characteristics:

- o An individual member of any homologous series is called homologue.
- o Homologous have same chemical properties but different physical properties.
- o All the members of homologous series can be represented by same general formula.



#### Homologous series

General formula

(i) Alkane

 $\mathbf{C}_n\mathbf{H}_{2n+2}$ 

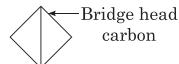
(ii) Alkene, Cycloalkane

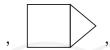
 $C_nH_{2n}$ 

(iii) Alkyne, Alkandiene, Bicuclo, Spiro

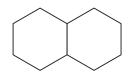
- $C_n C_{2n-2}$
- (a) Bicyclo Compounds: The substance in which 2 cyclic rings are fused at 2 common carbon atoms.

Eg.





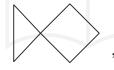


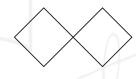


(b) Spiro Compounds: The compounds in which 2 cyclic rings are fused at single carbon.

Eg.







Spiro carbon

#### Homologous series

General formula

(iv) Alcohol, Ether

 $C_nH_{2n+2}O$ 

(v) Carbonyl compounds

(Aldehyde, Ketones)

Unsaturated alcohols & Ether

 $\mathbf{C}_n\mathbf{H}_{2n}\mathbf{O}$ 

(double bond containing)

Cyclic alcohol & Cyclic ether

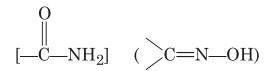
(vi) Carbocylic acid & Ester

 $C_nH_{2n}O_2$ 

Hydroxy carbonyl compounds

(vii) Amide oxime

 $C_nH_{2n+1}NO$ 





(viii) Aromatic alcohol & phenolic compounds

$$C_nH_{2n-6}O$$

(ix)  $1^{\circ}/2^{\circ}/3^{\circ}$  amine

$$C_nH_{2n+3}N$$

$$(-NH_2)(-NH-)(-N-)$$

Q. An alkane can be represented by following general formula.

1. 
$$C_n H_{2n+2}$$

$$n(n-1)$$

2. 
$$C_{n-1}H_{2n}$$

$$C_{n-1}H_{2(n-1)+2}$$

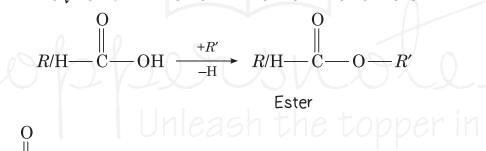
3. 
$$C_{n+1}H_{2n+4}(n=0)$$
  $C_{n-1}H_{2n}-2+2$ 

$$C_{n-1}H_{2n} - 2 + 2$$

Sol. All

$$I \rightarrow \boxed{n=2}$$

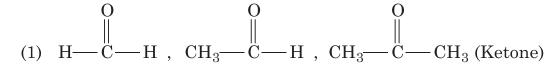
Find minimum required C for the second member of ester series. Q.



Sol.

IInd 
$$CH_3$$
— $C$ — $CH_3$  or  $H$ — $C$ — $CH_2$ — $CH_3$ 

Q. Select triad of Homologous.



(2) 
$$CH_3$$
— $NH_2$ ,  $CH_3$ — $NH$ — $CH_3$ ,  $CH_3$ — $N$ — $CH_3$  (1°/2°/3° amine)



(4) None

Sol. (4) None

# C. Classification of Carbon & Hydrogen Atoms

1. Types of Carbon: Carbon atom is classified on the basis of total next joining carbon atoms.

(i) Primary/1°-carbon : Eg. 
$$\overset{1^{\circ}}{C} = \overset{1^{\circ}}{C} - O - \overset{1^{\circ}}{C}$$
(Super 1°)

(ii) Secondary/2°-carbon : Eg. 
$$\overset{1^{\circ}}{C} \overset{2^{\circ}}{=} \overset{1^{\circ}}{C}, \quad \overset{1^{\circ}}{C} \overset{2^{\circ}}{=} \overset{1^{\circ}}{C}$$

$$NH_{2}$$

(iii) **Tertiary/3°-carbon :** Eg. 
$$\overset{1^{\circ}}{C} = \overset{3^{\circ}}{C} = \overset{1^{\circ}}{C}, \quad \overset{1^{\circ}}{C} = \overset{1^{\circ}}$$

Eg. 
$$CH_3$$
  $CH_3$   $CH_$ 

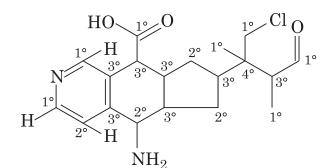
#### 2. Types of Hydrogen atom:

- (i) H atom is classified on the basis of C atom.
- (ii) H which is directly attached to  $1^{\circ}$ C/ $2^{\circ}$ C/ $3^{\circ}$ C atom is respectively called as primary H/ sec H/ tertiary H.



(iii) 1º/2º/3º Carbon atom may be saturated or unsaturated but quaternary Carbon atom is always saturated.

Q.



No of 1°C, 2°C,  $3^{\circ}$ C &  $4^{\circ}$ C and H = ?

 $3^{\circ} \rightarrow 7$ 

 $4^{\circ} \rightarrow 1$ 

(H)

 $\begin{array}{c} 1^{\circ} \rightarrow 11 \\ 2^{\circ} \rightarrow 6 \end{array}$ 

D. State of Hybridisation and Geometry

Structure

Hybridisation

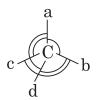
Geometry

Bond angle

٦.

 $sp^3(4\sigma)$ 

Tetrahedral

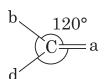


109°28' 6

[ab, bd, dc, ac, ad, bc]

 $sp^2(3\sigma)$ 

Trigonal planar



120°

3.  $-C \equiv$   $sp(2\sigma)$ 

Linear

180°

180°



4. 
$$= C =$$

$$sp(2\sigma)$$

Linear

180°



Eg.



$$sp^{3} \xrightarrow{Sp^{3}} CH_{3} \xrightarrow{CH_{3}} CH_{3} \xrightarrow{CH_{3}} CH_{3} \xrightarrow{Sp^{2}} CH_{3} \xrightarrow{CH_{3}} CH_{3} \xrightarrow{Sp^{3}} CH_{3}$$

 $\sigma$  bonds = 23 and  $\pi$  bonds = 4

$$sp^3 - 5sp^3 = 2$$

$$sp^3 - s = 13$$

$$< 109^{\circ}28' = 5 \times 6 = 30$$

$$sp^3 - sp^2 = 3$$

$$sp^2 - s = 0$$

$$< 120^{\circ} = 2 \times 3 = 6$$

$$sp^2 - sp = 3$$

$$sp - s = 1$$

$$< 180^{\circ} = 3 \times 1 = 3$$

$$sp - sp = 1$$

Nomenclature of Organic Substances:

[IUPAC/Geneva System]

A. Format for IUPAC Name

 $2^{\circ}$ -prefix +  $1^{\circ}$ -Prefix + Root word +  $1^{\circ}$ -suffix +  $2^{\circ}$ -suffix

(1)  $2^{\circ}$ -Prefix  $\rightarrow$  It is used according to substituent

#### Substituent

#### 2°-Prefix

$$-X(-F, -Cl, -Br, -I)$$

Halo

$$-NO_2$$

Nitro

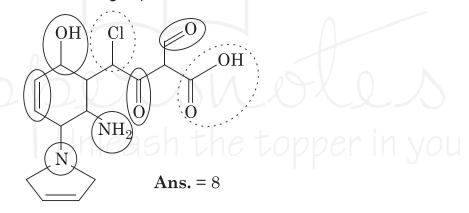
Nitroso



| —OR  | Alkoxy        |
|--|---------------|
| 0  | Epoxy         |
| -R   | Alkyl         |
| $CH_2X$                                    | Halomethyl    |
| O  |               |
| $-0$ — $\overset{\circ}{\mathrm{C}}$ — $R$ | Alkanoyloxy   |
| O<br>                                      |               |
| -NH-C-R                                    | Alkanoylamino |

# Note According to IUPAC system for naming —X, —OR — $NO_2$ , —NO,  $O \rightarrow$  All are considered as a substituent.

Q. Identify the total types of functional group.



- 2.  $1^{\circ}$ -prefix: It is a cyclo & used only for cyclic substance.
- 3. Word Root: It is used according to total no. of carbon atoms present in parent carbon chains.
- **4.**  $1^{\circ}$ -Suffix: It is used according to saturation or unsaturation.

C—C → ane
$$C = C \rightarrow \text{ene}$$

$$C \equiv C \rightarrow \text{yne}$$

$$2C = C \rightarrow \text{adiene}$$

$$2C \equiv C \rightarrow adiyne$$

Eg. 
$$C = C - C \equiv C$$

But-1-ene-3-yne ×

But-1-en-3yne or 1, 3-Butenyne ✓

**5.**  $2^{\circ}$ -suffix: It is used according to principal functional group.

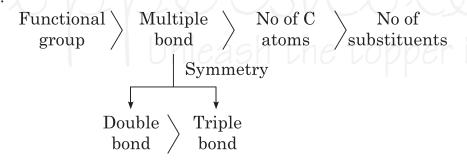
# <u>Note</u> In any compound, when more than one different functional groups are possible, then according to priority series, only one of them (having comparatively top rank in priority series) is considered as main functional group, while remaining are considered as substituent.

# B. Rules for IUPAC System

### 1. Selection of parent carbon chain :

\* We select that carbon chain as a parent including functional group, multiple bond & substituent.

#### Priority order:

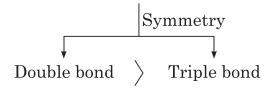


## 2. Numbering in parent carbon chain :

(i) In parent carbon chain, numbering can be done from that side from which multiple bond, substituent and functional group gets lowest number.

#### Priority order:

Functional group Multiple bond Substituents





| S.No. | Functional Group  | Prefix                     | Suffix                        |
|-------|---|----------------------------|-------------------------------|
| 1.    | (c) OOH [Carboxylic acid]<br>—COOH  | ×<br>Carboxy               | Oic acid carboxylic acid      |
| 2.    | —SO <sub>3</sub> H [Sulphonic acid]   | Sulpho                     | Sulphonic acid                |
| 3.    | O   | ×                          | Oic Anhydride                 |
| 4.    | -(C)OOR [Ester] $-COOR$   | ×<br>Alkoxy                | Alkyloate<br>Alkyl            |
|       | Alkanoyloxy $\left( \begin{array}{c} \mathbf{O} \\ \mathbf{II} \\ -\mathbf{O} - \mathbf{C} - R \end{array} \right)$ | Carbolkoxy  leash the to   | Carboxylate                   |
| 5.    | —(C)OX ∑Acid halide)<br>—COX  | ×<br>Halo formyl           | Oyl halide<br>Carbonyl halide |
| 6.    | $-(C)ONH_2$ [Amide] $-CONH_2$   | ×<br>Carbomoyl             | Amide<br>Carbocamide          |
| 7.    | —(C)N [Cyanide]<br>—CN  | ×<br>Cyano                 | Nitrile<br>Carbonitrile       |
| 8.    | $-N \stackrel{ ightharpoonup}{=} (C)$ [Isocyanide] $-NC$  | ×<br>lsocyano/Carbyl amino | Isonitrile<br>×               |



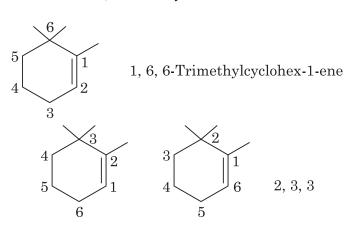
| S.No. | Functional Group         | Prefix   | Suffix                |
|-------|--------------------------|----------|-----------------------|
| q.    | —(C)HO [Aldehyde]        | formyl   | al, Carbaldehyde, one |
| 10.   | —(C)— [Ketone]<br>  <br> | keto/oxo | one                   |
| 11.   | —OH [Alcohol]            | Hydroxy  | OI                    |
| 12.   | —SH [Thio alcohol]       | mercapto | Thiol                 |
| 13.   | -NH <sub>2</sub> [Amine] | Amino    | Amine                 |
| 14.   | —OR [Ether]              | Alkoxy   | ×                     |

- \* When 2 different substituents are present at symmetric positions, then numbering can be done acc. to alphabet rule.
- \* If more than 2 substituents are present in parent Carbon chain, then we follow lowest locant rule.

Q. 1

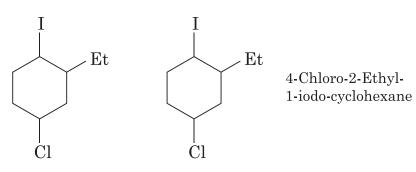
#### 5-Nitrohex-4-en-1-yne or 5-Nitro-4, 1-Hexenyne

Q. 2



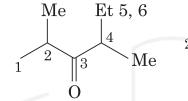




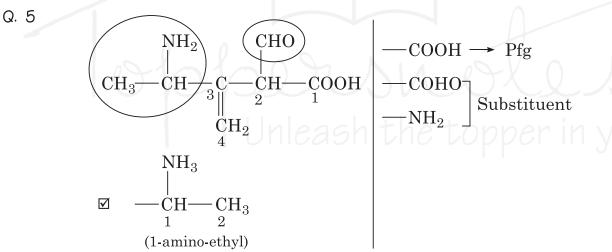


(Lowest locant rule)

(Lowest locant Rule)



2, 4-Dimethyl-hexan-3-one



#### Sol. 3-(1-Aminoethyl)-2-formyl but-3-enoic acid

$$\begin{array}{c|c} O & NH_2 \\ \hline OH & 3 & 2 & 1 \\ \hline CHO & O \end{array}$$

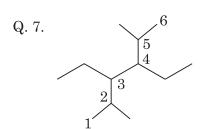
-COOH 
$$\rightarrow$$
 Pfg

-CHO 
$$\rightarrow$$
 Substitute

$$\text{-NH}_2 \, \to \, \text{Substitute}$$

Sol. 2-Amino-3-formyl-but-2-en-1, 4-dioic acid

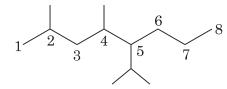




Sol. 3, 4-Diethyl-2, 5-Dimethyl hexane

\* If no. of C-atoms are same go for more number of substitutes.





$${\operatorname{CH}_3} \ | \ {\operatorname{CH-CH}_3} \ {\operatorname{CH-CH}_3} \ {\operatorname{Methyl}}$$

#### Sol. 2, 4-Dimethyl-5-(methyl ethyl) octane.

Q. 9.

$$8$$
 $6$ 
 $4$ 
 $3$ 
 $2$ 
 $1$ 

Sol. 5-(Dimethylethyl)-6-ethyl-2-methyl octane

- \* When Di Tri, Tetra are part of a name (like here then considered alphabetically.
- # <u>Note</u> Primary letter of extra prefix di, tri, tetra etc are not considered alphabetically, but whenever they are part of name, then we consider alphabetically.
  - \* In Hydrocarbon group or substituted substituents minimum number goes to that carbon which have free valency.
  - \* To indicate the total number of substituted substituents, we use some extra prefix like bis, tris, tetrakis etc.

Q. 10.

Sol. 6-Ethyl-2-Methyl-4, 5-bis-(Methyl ethyl) octane



Sol. N-ethyl methanamide

Sol. N-ethyl-N, 3-dimethyl butan-2-amine

# <u>Note</u> Whenever ester group is present as a main functional group, then we always write the name of Alkyl group.

Format  $\rightarrow$  **Alkyl.....oate.** 

or

Alkyl .....carboxylate

**V** 

$$(2^{\circ} p + 1^{\circ} - p + wR + 1^{\circ} - 5$$

Q. 13. O  $\operatorname{NH}_2$ H 4 3 2 1 0

$$\begin{array}{c} --\text{COOCH}_3 \rightarrow \text{Pfg} \\ --\text{CHO} \\ --\text{NH}_2 \end{array}$$
 Substituent

Sol. Methyl-2-amino-3-methyl-4-oro but-2-enoate.

Sol. 2-mercapto-3-methoxy carbonyl-butanoic acid