

Organic Chemistry

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ORGANIC CHEMISTRY

EARNING OBJECTIVES

- Classification of organic compounds
- Homologous series
- Characteristics of a homologous series

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- Nomenclature, IUPAC System
- Nomenclature of Branched hydrocarbons
- Nomenclature of (A) Cyclo-alkanes, Cyclo-alkenes, Cyclo-alkynes
- Isomerism, structural isomerism types.

Real life applications:

Organic chemistry is a subject with which many people struggle. It is important because it impacts our every day lifes.

Classification of organic compounds

A) Classification based on structure : The simplest organic compounds containing only carbons and hydrogens are called hydrocarbons. Other organic compounds are thought to have been derivative of hydrocarbons. All the known organic compounds have been broadly divided into two categories depending upon the nature of their carbon skeleton.

- (A) Open-chain or Acyclic compounds.
- (B) Closed-chain or Cyclic compounds.

(A) Open-chain or Acyclic compounds : These are the compounds in which the carbon atoms are linked to each other in such a manner that the molecule is having an open-chain structure. The carbon chain may be either straight chain or branched-chain. Such compounds are known as aliphatic compounds.

Examples are :

$$CH_3 - CH_2 - CH_2 - CH_2 - CH_3$$

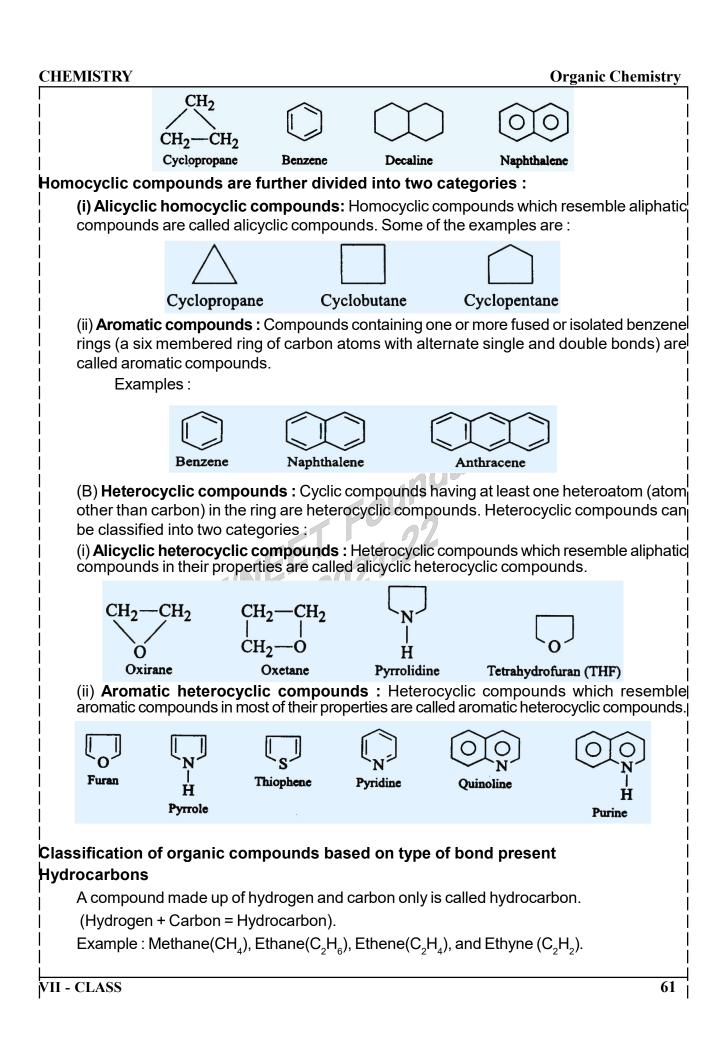
n-pentane
 $CH_3 - CH_2 - CH_2$
1-Butene
 CH_3

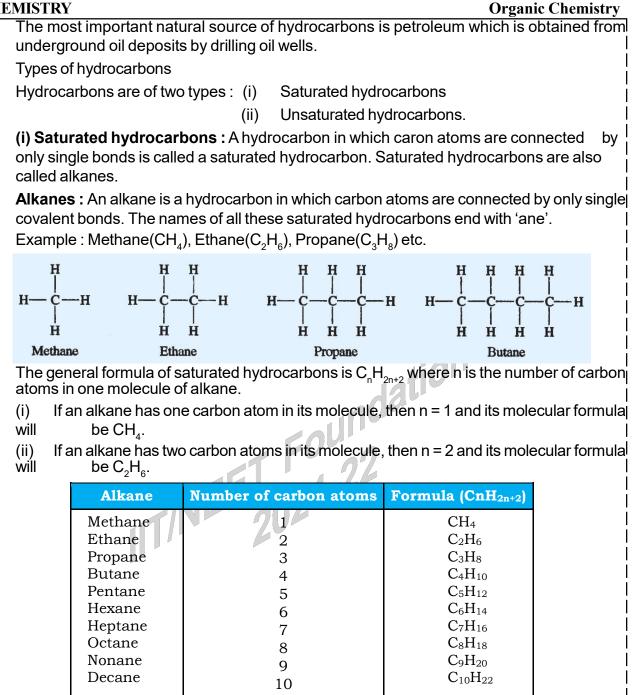
$$\begin{array}{c|c} CH_{\overline{3}} - CH - CH_3 & | \\ | & CH_{\overline{3}} - C - CH_3 \\ CH_3 & | \\ CH_3 & CH_3 \\ Isobutane & Neopentane \end{array}$$

Neopentane

B) Closed-chain or Cyclic compounds : Compounds containing atleast one ring in their structure are called cyclic compounds. If the ring contains one type of atoms, the compound is homocyclic, and if the ring contains atleast one hetero atom, the compound is called heterocyclic. Depending upon the nature of the ring, these compounds are further divided into the following two categories

(i) Homocyclic or Carbocyclic compounds : When all the atoms of the ring are carbons, the cyclic compounds are known as homocyclic compounds. For example,





The saturated hydrocarbons are chemically not very reactive.

(ii) Unsaturated hydrocarbons : A hydrocarbon in which two carbon atoms are connected by a double bond or a triple bond is called an unsaturated hydrocarbon. Alkenes and alkynes are called unsaturated hydrocarbons.

Example : Ethene(C_2H_4), Propene(C_3H_6), Butene(C_4H_8)

Alkenes : An alkene is a hydrocarbon in which carbon atoms are connected by double bonds. The names of alkenes end with 'ene'.

The general formula of alkenes is $C_n H_{2n}$ where n is the number of carbon atoms in one

molecule of alkane.

- (i) If an alkene has two carbon atoms in its molecule, then n = 2 and its molecular formulal will be C_2H_4 .
- (ii) If an alkene has three carbon atoms in its molecule, then n = 3 and its molecular formula will be C_3H_6 .

Alene	Number of carbon atoms	Formula (C _n H _{2n})
Ethene	2	C_2H_4
Propene	3	C_3H_6
Butene	4	C_4H_8
Pentene	5	$C_{5}H_{10}$
Hexene	6	$C_{6}H_{12}$
Heptene	7	$C_{7}H_{14}$
Octene	8	$C_{8}H_{16}$
Nonene	9	$C_{9}H_{18}$
Decene	10	$C_{10}H_{20}$

Alkynes : An unsaturated hydrocarbons in which the two carbon atoms are connected by a triple bond is called an alkyne. The names of alkynes and with 'yne'.

Example : Ethyne(C_2H_2), Propyne(C_3H_4), Butyne(C_4H_6) etc.

The general formula of alkynes is C_nH_{2n-2} where n is the number of carbon atoms in one molecule of alkane.

(i) If an alkyne has two carbon atoms in its molecule, then n = 2 and its molecular formula will be C_2H_2 .

(ii) If an alkane has three carbon atoms in its molecule, then n = 3 and its molecular formula will be C_3H_4 .

Alkyne	Number of carbon atoms	Formula (C _n H _{2n-2})
Ethyne	2	C_2H_2
Propyne	3	C_3H_4
Butyne	4	C_4H_6
Pentyne	5	C_5H_8
Hexyne	6	$C_{6}H_{10}$
Heptyne	7	$C_7 H_{12}$
Octyne	8	$C_{8}H_{14}$
Nonyne	9	$C_{9}H_{16}$
Decyne	10	$C_{10}H_{18}$

Alkyl group : By removing a hydrogen atom from an alkane, alkyl group is obtained. Alkyl group is represented by R –.

While deriving the name of an alkyl group "ane" of alkane is replaced with "yl". Thus

Alkane – H atom = Alkyl group

A unit called methylene group - CH₂

The difference between carbon and hydrogen atoms in the first four members of alkanes is,

- 1. Difference between ethane and methane = $[C_2H_6 CH_4) = CH_2$
- 2. Difference between propane and ethane = $[C_3H_8 C_2H_6] = CH_2$
- 3. Difference between butane and propane = $[C_4H_{10} C_3H_8] = CH_2$

The difference of carbon and hydrogen atoms between two consecutive members of alkanes is CH_2 . This ---CH₂ unit is called methylene group.

By removing one or more CH_2 groups from a given saturated hydrocarbon, we can obtain another saturated hydrocarbon with lesser number of carbon atoms.

Similarly, by adding one or more CH_2 groups in a given saturated hydrocarbon, we can obtain another hydrocarbon with more number of carbon atoms.

Furthermore, there is a difference of 14 amu (molecular mass of CH_2 is 14) between two consecutive hydrocarbons.

Homologous series

Alkanes have a general formula $C_n H_{2n+2}$ and each member differs from the next by a fixed group — CH_2 — (the methylene group), even they have similar chemical properties.

In other classes of compounds like alkenes, alkynes, alcohols or acids, the same thing happens. Such a series of compounds is known as a homologous series and the individual members are homologues.

Compounds having the same functional group and similar properties that differ from the adjacent members by a $- CH_2$ – group constitute a homologous series and the phenomenon is called homology.

The members of the same class of organic compounds, when arranged in the order of ascending molecular weights such that they differ from each other by $- CH_2$ group, are collectively called homologues.

Example of a homologous series for each class (or family) of compounds are :

Compound	Formula (C _n H _{2n+2})	Difference in formulae
Methane	CH ₄	CH_2
Ethane	C_3H_8	CH_2
Propane	C_2H_6	CH_2
Butane	$C_{4}H_{10}$	CH_2
Pentane	$C_{5}H_{12}$	CH_2

Compound	Formula (C _n H _{2n})	Difference in formulae
Ethene Propene Butene Pentene	$C_{2}H_{4}\ C_{3}H_{6}\ C_{4}H_{8}\ C_{5}H_{10}$	$egin{array}{c} \mathrm{CH}_2 \ \mathrm{CH}_2 \ \mathrm{CH}_2 \ \mathrm{CH}_2 \end{array}$

Compound	Formula (C _n H _{2n-2})	Difference in formulae
Ethyne Propyne Butyne Pentyne	$egin{array}{c} C_2H_2\ C_3H_4\ C_4H_6\ C_5H_8 \end{array}$	$\begin{array}{c} CH_2\\ CH_2\\ CH_2\\ CH_2\end{array}$

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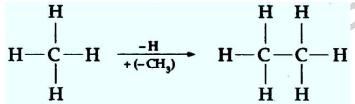
Compound	Formula (C _n H _{2n+2})	Difference in formulae
Methane	CH ₄	$ m CH_2$
Ethane	C_3H_8	CH_2
Propane	C_2H_6	CH_2
Butane	$C_{4}H_{10}$	CH_2
Pentane	$C_{5}H_{12}$	$ m CH_2$

Characteristics of a homologous series

The members of a homologous series can be represented by a general formula, e.g., the alkanes by C_nH_{2n+2} , alkenes by C_nH_{2n} , alkynes by C_nH_{2n-2} and alcohols by $C_nH_{2n+1}OH$, where 'n' stands for number of carbon atoms in one molecule of alkane.

Every member of a homologous series differs from the adjacent ones in composition by CH_2 .

This is because every succeeding member of a homologous series arises by the replacement of an H atom in the preceding member by a CH_3 group. For example, C_2H_6 (ethane) may be thought to arise from CH_4 (methane) as follows.



The molecular masses of two consecutive members of a homologous series differ by 14 amu (equivalent to one CH, group).

Generally, similar methods may be employed to prepare all the compounds belonging to a homologous series.

The physical properties of the compounds of a homologous series change gradually as the molecular mass changes. For example, the melting point, boiling point and density increase with molecular mass.

The chemical properties of the compounds belonging to the same homologous series are similar.

Nomenclature Of Organic Compounds

Naming of organic compounds is an important aspect in the study of organic chemistry. since some millions of compounds are present, remembering their names is an herculean task. There are mainly three types of naming.

They are i) Trivial naming, ii) Derived naming and iii) Systematic naming.

Trivial names are also called common names. In this system the compound is named after the source from which it is obtained. For instance citric acid (obtained from citrus plant), uric acid (obtained from urine), formic acid (obtained from red ants, formica means ants in Latin) et.

Derived names are based on the recognition of certain familiar common names. Thus CH_3 .OH is called carbinol and CH_3 .CH₂.OH is called methylcarbinol.

II. IUPAC System

With the rapid growth of organic chemistry, it was realised that the nomenclature of fast increasing number of organic compounds should be related to their structures in a systematic manner as to avoid undue strain on memory. The first rational system of naming organic compounds was evolved in 1892 by the International Chemical Congress at Geneva under the name of Geneva system of nomenclature. However, the system could not be successfully applied to complicated and multifunctional organic compounds. This system was partially revised in 1931 and the revised system was referred to IUPAC system.

Modifications in this system have been made from time to time by the International Union of Pure and Applied Chemistry and in its present form, the system is known as IUPAC system (pronounced as eye-you-pack). It is the most modern and largely used system. This system has been evolved in 1957 and some additions and subtractions were done in 1967. The IUPAC system has set rules for naming organic compounds on the basis of their structures.

These rules underwent further modifications in 1979 and later revised in 1993 (A Guide to IUPAC Nomenclature of Organic Chemistry by R. Panico, W.H. Powell and J.C. Richer). The name assigned on the basis of latest IUPAC rules to an organic compound is known as its systematic name.

Salient Features of IUPAC System

- 1. A given compound can be assigned only one name.
- 2. A given name can clearly direct in writing of one and only one molecular structure.
- 3. The system can be applied in naming complex organic compounds.
- 4. The system can be applied in naming multifunctional organic compounds.
- 5. This is a simple, systematic and scientific method for nomenclature of organic compounds.

Basic rules of nomenclature : For naming simple aliphatic compounds, the normal saturated hydrocarbons have been considered as the parent compounds and the other compounds as their derivatives obtained by the replacement of one or more hydrogen atoms with various functional groups. Each systematic name has two or three of the following parts:

(i) Root word, (ii) Primary suffix (iii) Secondary suffix.

Which indicate linear or continuus chains or carbon atoms are known by special root words while chains from C_5 onwards are known by Greek number roots.

Chain length	Root word	Chain length	Root word
$\begin{array}{c} C_{1} \\ C_{2} \\ C_{3} \\ C_{4} \\ C_{5} \\ C_{6} \\ C_{7} \\ C_{8} \\ C_{9} \\ C_{10} \end{array}$	Meth- Eth- Prop- But- Pent- Hex- Hept- Oct- Non- Dec-	$\begin{array}{c} C_{11} \\ C_{12} \\ C_{13} \\ C_{14} \\ C_{15} \\ C_{16} \\ C_{20} \\ C_{30} \\ C_{40} \\ C_{50} \end{array}$	Undec- Dodec Tride- Tetradec- Pentadec- Hexadec- Eicos- Tnacont- Tetracont- Pentacont-

In general, the root word for any carbon chain is alk-. (ii) Primary suffixes: Primary suffixes are added to the root words to show saturation or unsaturation in a carbon chain.

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Classification of Nomenclature of organic compound

Nature of carbon chain	Primary suffix	IUPAC name
Saturated (C-C)	-ane	Alkane
\int unsaturated (C = C)	-ene	Alkene
\with one double bond		
\int unsaturated (C = C)		Alkyne
with one triple bond	-yne	hikyite
∫unsaturated with	-adiene	Alkadiene
$\int two (C = C) bonds$		
(unsaturated with	a alianna a	Alkadiyne
$\int two (C = C) bonds$	-adiyne	
unsaturated with		
three (C = C) bonds	-atriene	Alkatriene

Naming of hydrocarbons

i) The number of carbon atoms in a hydrocarbon is indicated as

Number of carbon atoms	Name	Number of carbon atoms	Name
1	Meth	6 0 0	Hex
2	Eth	7	Hept
3	Prop	8	Oct
4	But	9	Non
5	Pent	10	Dec

ii) A saturated hydrocarbon containing single bonds is indicated by writing the word 'ane' at the end.

iii) An unsaturated hydrocarbon containing double bonds is indicated by writing the word 'ene' at the end.

iv) An unsaturated hydrocarbon containing triple bonds is indicated by writing the word 'yne' at the end.

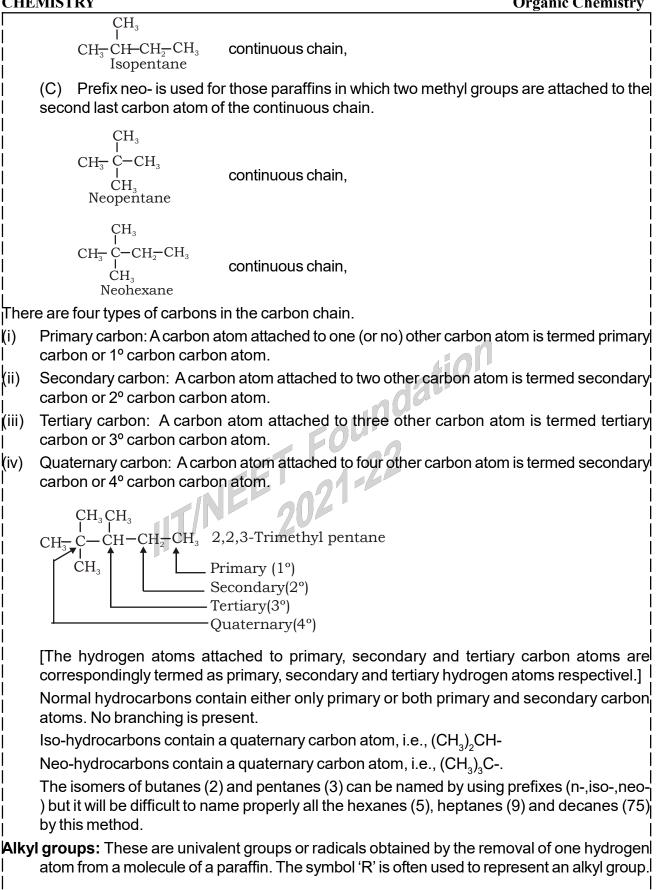
All isomeric paraffins have same parent name. The names of various isomers are distinguished by prefixes.

(a) Prefix n- is used for those paraffins in which all the carbon atoms are in one continuous chain. The prefix n-stands for normal or stright chain

Ex:	$CH_{3}CH_{2}CH_{2}CH_{3}$,	n-Butane
	$CH_{3}CH_{2}CH_{2}CH_{2}CH_{3}$,	n-Pentane
	$CH_3CH_2CH_2CH_2CH_2CH_3$	n-Hexane

(b) Prefix iso- is used fro those paraffins in which methyl group is attached to the second last carbon atom of the continous chain.

CH_3 CH_3 $CHCH_3$ Isobutane	continuous chain,
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$$(\text{Alkane})C_{n}H_{2n+2} \xrightarrow{-H} C_{n}H_{2n+1} (\text{Alkylgroup})$$

Nomenclature of Branched hydrocarbons IUPAC Rules

- 1. For saturated hydrocarbons and their substituent derivatives
- (i) Longest chain rule :
- The longest continuous chain of carbon atoms is selected as the parent hydrocarbon.
 The compound is then named as a derivative of the parent hydrocarbon.
- (b) If more than one set of longest possible chain are possible then the selected longest chain should have(i) maximum number of side chains, or (ii) minimum number of branched side chains.

$$\begin{array}{c}
1 & 2 & 3 & 4 & 5 & 6\\
CH_2 - CH - CH_2 - CH_2 - CH_2 - CH_3 & (Wrong)\\
CH_2 \\
CH_3 \\
CH_3 - {}^3CH - {}^4CH_2 - {}^5CH_2 - {}^6CH_2 - {}^7CH_3(Right)\\
2 {}^{CH_2} \\
1 {}^{CH_2} \\
1 {}^{CH_3} \\
\end{array}$$

Carbon atom chain with one side chains or 7 carbon atoms chain with one side chains having no branching (right)

6 Carbon atom chain with one side chains or 6 carbon atoms chain with one branched side chain (wrong)

i) Lowest sum rule:

(a) The selected chain is numbered in terms of arabic numbers one end to

other 1 2 3 4 (Right)

$$CH_3-CH_2-CH_2-CH_3$$
 (Right)
4 3 2 1

- (b) Lowest number is assigned to the first side chain (alkyl group) or a substitutent (Cl,-Br,-I,-NO₂)
- (c) If two different alkyl groups are at the same position from opposite ends, lowest number is given in order of their alphabets (IUPAC 1993)

eg: 1 2 3 4 5 6 7

$$CH_3-CH_2-CH-CH_2-CH-CH_2-CH_3$$
 (Wrong)
7 6 5 4 3 2 1 (Right)
 $CH_3 C_2H_5$

(d) If two different substitutents are at the same position from opposite ends, lowest number is assigned in order of their alphabets,

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2 3 4 1 eg: CH₃-CH-CH-CH₃ (Wrong) 4 3 2 1 (Right) CI Br If a substitutents and side chains are at the same position ends, opposite eneds, (e) lowest number is assigned to the substitutents. If more than two substitutents and side chain are present, the sum of their (f) numbers should be lowest at the first preference, irrespective of the naturte of 4 3 2 1 CH₃-CH-CH-CH₃(Right) 2 3 4 (Wrong) CH₃ CI side chain or substitutents. Note: See lowest sum at the first preference, as over lined and never see total sum of the numbers. (iii) Arrangements of prefixes: The names of a saturated hydrocarbon in general, may (a) be divided into two parts: word root and suffix. The word root designates table1) the number of carbon atoms in the chain. The suffix (ane) is added to the word to indicate the saturation in hydrocarbon Alkyl nature of side chain or substituent gp. is identified and reported as prefix (b) (table 4 and 5) with its number (locant) in hydrocarbon name in alphabatic order an d while doing so prefixes, di, tri, tetra are not to be considered. If more than one similar alkyl chains or substituents are present, prefix names are (c) suitably modified by putting di, tri....terms. CH₃—CH—CH—CH₃ ĊH, ĊH, 2,3-Dimethyl butane If more than one similar alkyl gps. or substituents are present, at same position, their (d) no. is also repeated. In case side chain is also branched, it is also numbered from the carbon atom (e) attached to main chain and reported in paran thesis. Rules for Naming Complex Aliphatic Compounds Containing One Functional Group: (A) Longest chain: The parent carbon chain in so chosen as to include the functional group even if it is not the actual longest continuous chain. ${\stackrel{5}{C}}H_3 - {\stackrel{4}{C}}H_2 - {\stackrel{3}{C}}H_2 - {\stackrel{2}{C}}H - CH_2 - CH_3$ ¹CHOH VII - CLASS 70

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The longest chain (horizontal) consists of six carbon atoms but this does not include the functional group. This chain is rejected and the chain having five carbon atoms is selected as parent chain as it includes the functional group.

(B) Numbering of parent chain:

The numbering of the parent carbon chain is done in such as way that the carbon linking to functional group gets the lowest number even if there is violation of saturated hydrocarbon rules.

$${\stackrel{5}{C}}H_{3} - {\stackrel{4}{C}}H_{2} - {\stackrel{3}{C}}H_{2} - {\stackrel{2}{C}}H - CH_{2} - CH_{3}$$

The longest chain (horizontal) consists of six carbon atoms but this does not include the functional group. This chain is rejected and the chain having five carbon atoms is selected as parent chain as it includes the functional group. Hence, named as, Numbering of parent chain: The numbering of the parent carbon chain is done in such as way that the carbon linking to functional group gets the lowest number even if there is violation of saturated hydrocarbon rules.

$$\begin{array}{c} CH_{3} - CH_{2} - CH_{4} - CH_{2} - CH_{3} \\ CH_{2} - CH_{2} - CH_{2} - CH_{3} \\ CH_{2} - CH_{2} - CH_{2} - CH_{2} - CH_{3} \\ CH_{2} - CH_{2} - CH_{2} - CH_{2} - CH_{2} - CH_{3} \\ CH_{3} - CH_{2} - CH_{2} - CH_{2} - CH_{3} \\ CH_{3} - CH_{2} - CH_{2} - CH_{3} \\ CH_{3} - CH_{3} \\ CH_{3} - CH_{2} - CH_{2} - CH_{3} \\ CH_{3} - CH_{3} \\ CH_{3} - CH_{2} - CH_{2} - CH_{3} \\ CH_{3} \\ CH_{3} - CH_{3} \\ CH_{3$$

Note: When a chain terminating group such as – CHO, –COOH, COOR, – CONH₂ – C = N, etc., is present as the functional group, it must be assigned number 1. This does not apply to non-terminal groups such as >CO, – NH₂ and – OH which may or may not be assigned 1.

$$\overset{4}{C}H_{3} - \overset{3}{C}H_{2} - \overset{2}{C}H - CH_{2} - CH_{3} \qquad \overset{6}{C}H_{3} - \overset{5}{C}H_{2} - \overset{4}{C}H_{2} - \overset{3}{C}H - \overset{2}{C}H_{2} - \overset{1}{C}H_{3}$$

(C) The last 'e' of the primary suffix is dropped and the secondary suffix representing the functional group is added. The number giving the position of the functional group is inserted in the name.

The compound ${}^{6}_{CH_3}{}^{5}_{CH_2}{}^{4}_{CH_2}{}^{3}_{CH_2}{}^{2}_{CO}{}^{1}_{CH_3}$ is named as:Hexan-2-one

Similarly, the compound ${}^{1}_{C}H_{3} {}^{2}_{C}H_{2} {}^{3}_{C}HOH {}^{4}_{C}H_{2} {}^{5}_{C}H_{3}$ may be named as: Pental-3-ol

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The names of the substituents are prefixed to the parent hydrocarbon according to (D) IUPAC rules with alphabetical order without considering the presence of functional group. Halo and nitro groups are considered as substituents. (E) Numerical prefixes di-, tr-, tetra-, etc., are attached before the designations of functional group if two or more identical groups are present, e.g., ${}^{1}_{CH_{2}OH} - {}^{2}_{CH_{2}OH}$ is named Ethane-1, 2- diol. ²CHCOOH is named Butane - 1, 4 - dioic acid. CHÇOOH Examples: (A) $\overset{5}{C}H_{3} - \overset{4}{C}H_{2} - \overset{3}{C}H_{2} - \overset{2}{C}H - CH_{2} - CH_{3}$ lation The name of the compound is given in the following way: Secondary suffix 2 - ethylpentan - 1 0l(functional group) Locant for Locant Prefix Root functional group word Primary suffix i.e., 2-Ethylpentan-1-ol. $\overset{2}{C}H_{3} - \overset{1}{\overset{1}{C}}H - \overset{2}{\overset{1}{C}}H - \overset{3}{\overset{1}{C}}H_{2} - \overset{4}{\overset{1}{C}}H - \overset{5}{\overset{1}{C}}H$ ĊH₃ ĊHO ĊH₂ 2(1'-Methyl ethyl)-4-methylpentanal or 2-Isopropyl-4-methylpentanal Rules for Naming Aliphatic Compounds Having Polyfunctional Groups. A compound is said to be polyfunctional compound if it contains more than one functional group. The multiple bond (> C = C < or - C = C -) is also considered as a functional group. In IUPAC system, one of the functional groups is chosen as the principal functional group (secondary suffix) and the remaining functional groups (secondary functional groups) are treated as substituents and indicated by prefixes. For example, in the following structure. OH ← Substituent $CH_3 - CH - CH_2 - CH_2 - COOH - Principal functional group$

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The – COOH group is the principal functional group, while the –OH group is a substituent. The principal functional group is mentioned with its suffix name while secondary functional groups are mentioned only with their prefix name.

The choice of the principal functional group is made on the basis of the following order of preference.

[Carboxylic acids > sulphonic acids > acid anhydrides > esters > acid chlorides > acid amides > cyanides > aldehydes > ketones > alcohols, phenols, thiols > amines > ethers > alkenes, alkynes > halo, nitro, alkyl.]

The order of seniority among the principal groups is given according to the following table along with their prefix and suffix names. The functional group which occurs higher up in the table is the principal functional group.

Seniority Table for Principal Groups

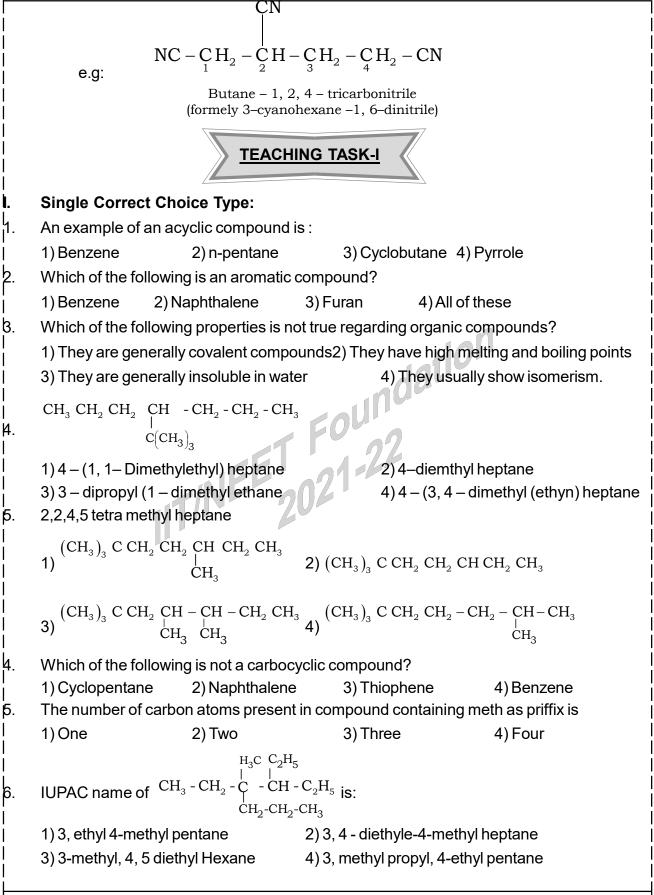
(Higher Priority Group at the Top)

Group	Prefix name	Suffix name
– COOH	Carboxy	-oic acid
– SO ₃ H	Sulpho	sulphonic acid
– COOR	Alkoxy carbonyl or Carbalkoxy	alkyl oate
– COX	Haloformyl or halocarbonyl	–oyl halide
$- CONH_2$	Carbamoyl	amide
– CN	Cyano	nitrile
– NC	Isocyano	carbylamine
– CHO	Formyl or aldo	-al
> C = O<	Keto or oxo	-one
– OH	Hydroxy	–ol
– SH	Mercapto	thiol
$- \mathrm{NH}_2$	Amino	amine
– OR	Alkoxy	—
-C - C -	Epoxy	—
0		
> C = C<	—	– ene
$-C \equiv C -$		– yne
-N = N -	Azo	—
$-NO_2$	Nitro	—
– NO	Nitroso	—
- X	Halo (Chloro, bromo, iodo)	—

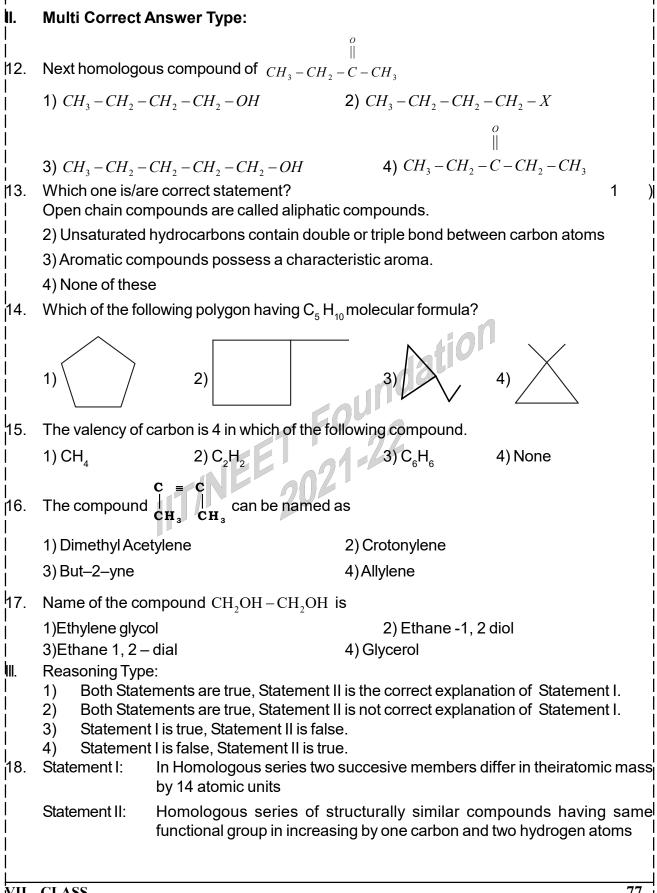
- The first step in the naming of polyfunctional compounds is the selection of principal functional group. The principal functional group gives the class name of the structure.
- 2. The second step is the selection of parent chain. The parent chain is so elected that it includes the maximum number of functional groups including the principal group.
- B. The third step is the numbering of parent chain. The parent chain is numbered from the side of principal functional group, i.e., it gets lowest number. The following decreasing order of preference for giving the lowest numbers is followed.

Principal functional group > Double bond or Triple bond > Substituents

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1.	Substituents, side chains and secondary functional groups are named in alphabetical order.
	On the basis of the above rules, the following general procedure is followed in naming an aliphatic compound.
a)	Locate the longest chain containing the principal functional group and as many of the
	secondary functional groups and multiple bonds are possible.
b)	Select the corresponding root word for the chain selected.
c)	Number the carbon atoms of the chain selected from the end so that the various
	groups get as small numbers as possible and the principal functional group gets the
	lowest number. The number of chain terminating group
	(i.e., $-COOH$, $-COOR$, $-COCI$, $-CONH_2$, $-CHO$, or $-CN$) must always be 1. When
	no functional group is present, the substituents should get the lowest numbers.
d)	Attach to the root word, the primary suffix (-ane, -ene or -yne) representing
. ,	the nature of carbon-carbon bonds in the selected chain.
e)	Add suitable prefixes (alphabetical order) and secondary suffixes with their
. /	respective numbers of attachment to parent chain as to denote the number and position for each substituent or functional group.
f)	
1)	In case the substituent has a branched chain, number the carbon atoms in the chain also from the carbon atom attached to the main chain. The name of such substituent
	is written in bracket and a hyphen (-) is put between the locant and the bracket. In such cases carbon of substituent attached to main chain is given number 1', next one number 2 and so on.
g)	Write the name of the compound as one word, separating the numbers from t h e
97	designations or prefixes or suffixes by hyphens and separating the numbers from each other by commas.
Som	ne Specific Rules for IUPAC Nomenclature
1.	If the substituent group is named as an oxygen derivative of $CH_3 - or CH_2$ group, the carbon atom of the substituent group is also counted towards parent chain. example: H
	O = C + C + C + C + C + OOH 3-Formylpropanoic acid
2.	Nomenclature of polyfunctional compounds containing more than two like functional groups.
	According to IUPAC convention 1993, if an unbranched carbon chain is directly linked to more than two like functional groups, the organic compound is named as a derivative of parent alkane which does not include the carbon atom of the functional groups. These are named by use of suffix such as tricarboxylic acid (for three – COOH groups), tricarbaldehyde (for three – CHO – groups) or tricarbonitrile (for three – CN groups).



Which of the following structure is 2, 2, 5-trimethyle heptane? 7. 1) $\begin{pmatrix} (CH_3)_3 C CH_2 CH_2 CH_2 CH_2 CH_3 \\ \downarrow \\ CH_2 \end{pmatrix}$ 2) $(CH_3)_3 C CH_2 CH_2 CH_2 CH_2 CH_3$ CH₃ The IUPAC name of the compound is $CH_3 - CH_2 - CH - CH_2 - CH_$ ß. ĊH-CH₃ CH₂ , Сн₂-Сн3 1) 2, 2, 5 – trimethyl – 4 – (1 – methyl propyl) hexane 2) 2, 2, 7 - trimethyl - 4 - (1 - methyl propyl) nonane 3) 2, 2, 7 – dimethyl – 4 – (1 - methyl propyn) octane 4) 2, 2, 5 – trimethyl– 4 – (1 – methyl propyn) oxtane CH₃ The IUPAC name of the compound $ch_3 - ch_2 - ch - ch_2 - ch - ch_3$ is 9 **Н₃С**—С--CH₃ CH. 1) 4 - (1, 2 - Dimethyl ethyl) - 5 - ethyl (-2 - methyl octane)2) 4 - (1, 1 - Dimethyl ethyl) - 3,6 - dimethyl octane 3) 5 - (1, 1 - Dimethyl ethyl) - 6 - ethyl (-5 - methyl octane)4) 4 - (1, 2 - Dimethyl ethyl) - 6 - ethyl (-4 - methyl octane)10. .The substituent group for prefixes Amino, Bromo, Methyl, Nitro are respectively, 2) $-\overset{\oplus}{N} \equiv N_{2} - CH_{2} - Br_{2} - NH_{2}$ 1) $-CH_3$, $-NH_2$, -Br, $-NO_2$ 3) $-NH_2$, -Br, $-CH_3$, $-NO_2$ 4) $-NH_2 - C\ell$, $-NO_2$, $-CH_3$ CH₂ OH IUPAC name of is CH-CH-CH-CH-CHO 1) 4-hydroxy-2-methyl pentanal 2) 2-hydroxy-4-methyl pentanal 4) 2-hydroxy-4-methyl pentanol 3) 4–hydroxy–2–methyl pentanol h1.. IUPAC name of $HO - CH_2 - CH - COOH$ is : NH₂ 1) 3–hydroxy–2–amino butanoic acid 2) 2-amino-3-carboxy propanal 3) 2–amino–3–hydroxy propanoic acid 4) 2-amino-3-hydroxy butanoic acid

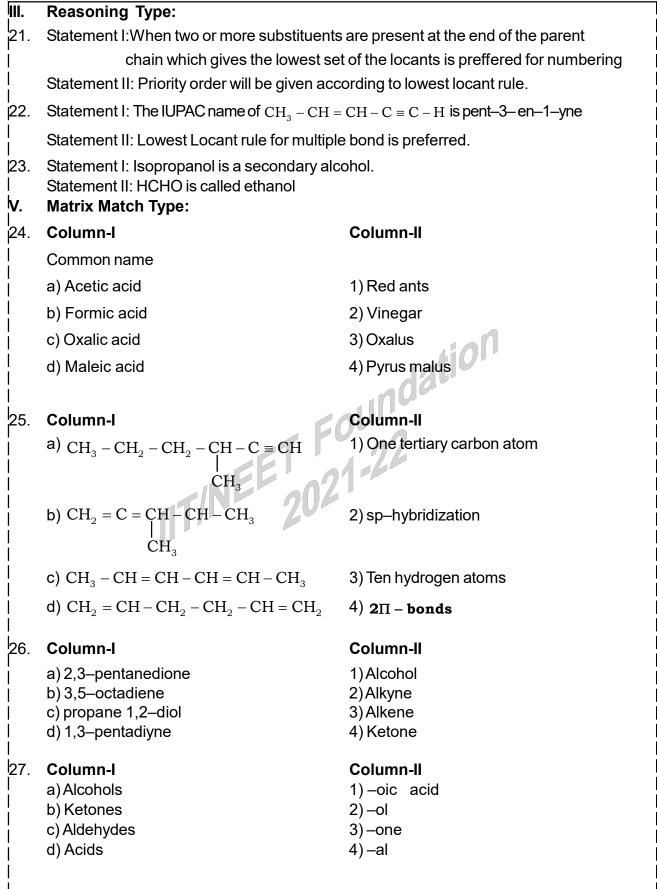


19.	Statement I:	The IUPAC name of $CH_3 - C(CH_3)_2 - CH_2 - CH = CH_2$ is 4,4 – dimethyl- 1–pentene.						
	Statement II:	The IUPAC name of $C_2H_5 - C_2 - CH_2 - OH$ is 2-ethyl prop-2-en-1-ol.						
V. 20.	Matrix Match T Column-I	Гуре:			Column-II			
	a) Next homologous compound of CH_4				1) <i>C</i> ₂ <i>H</i> ₄ <i>O</i>			
	b) Next homologous compound of $C_2 H_5 CI$			2) $C_5 H_{11} O$				
	c) Next homolog	gous compound (of HCHO		3) $C_2 H_6$			
	d) Next homolo	gous compound	of	∕_он	4) <i>C</i> ₃ <i>H</i> ₇ <i>Cl</i>			
21.	Column-I			Column II	5) C ₂ H ₄			
	. —		Column-II					
	a) 1) Pentane							
	b) c) d)	ITNEE	2021	3) Nonan	thyl butane e nethyl butane			
/. \.	Comprehension Type: Homologous series is a series of structurally similar compounds having same functional group in increasing number of carbon atoms.							
23.	The molecular weight of one compound is 16. Next homologous compound weight weight of one compound is 16.							
	1) 28	2) 30	3) 32	4) 29				
24.	Next homologous compound of $CH_3 - CH = CH_2$ is							
	$1) CH_3 - CH = 0$	$CH - CH_3$	2) <i>CH</i> ₃ -	$-CH = CH_2$				
	3) $CH_2 = C =$	CH ₂	4) <i>CH</i> ₂ =	$=CH_2$				
25.	Next homologous compound of $CH_2 = CH - OH$ is							
	1) $CH_3 - CH =$	CH – OH	2) CH ₃ -	$-CH_2 - CH_2$	-ОН			
	0) <i>CH CH</i>	CH = CH - OH	A) CU	$-CH_2 - OH$				

CHEMISTRY **Organic Chemistry** I.U.P.A.C System is a conference of the well known chemists of nearly all the Scientifically Β. advanced countries was convened in Geneva in 1892 to derive a uniform scheme of naming aliphatic organic compounds. This international body was eventually formalised as the International Union of Pure Applied Chemistry (IUPAC) at a meeting in Paris in 1957. This official system of naming organic compounds is referred to as IUPAC System of Nomenclature or simply IUPAC Nomenclature. CH₃-CH-CH₂-CH₂-CH₃ 26. The IUPAC name of the following compound is $H_3C - CH - CH_3$ 1) 2- isopropyl pentane 2) 2, 3 – dimethyl hexane 4) 2, 4 - dimethyl hexane 3) isononane Which of the following compound is 2, 2 - di methyl pentane 27. CH₃ СН3 2) CH₃ - CH₂ - C - CH₂ - CH₃ 1) сн₃ - с - сн₂ - сн₂ - сн₃ СН₃ CH₃ CH₃ CH, CH, CH **3)** C H ₃ - C H ₂ - C - C H ₃ -CH-CH-CH₂-CH₃ CH₃ The IUPAC name of the following compound is 28. 1) 1, 1 – diethyl 1–2, 2– dimethyl pentane 2) 4, 4 dimethyl 5, 5 – diethyl pentane 3) 5, 5 – diethyl 4, 4 diemthyl pentane 4) 3 - ethyl 4, 4 - dimethyl heptane C. The numbering of the parent carbon chain is done in such a way that the carbon linking to funcitonal group gets the lowest number even of there is violation ofl saturated hydrocarbon rules. The IUPAC name of $\begin{array}{c} CH_3-CH_2-CH_2-CH_2-CH_2-CH_3 \end{array}$ is 29. $\dot{C}H_2 - CH_2OH$ 1) 4–Ethyl–6–hexan–1–ol 2) 3-Ethyl-3-propyl propanol 3) Octanol 4) 3-Ethyl hexan-1-ol $\beta 0.$ The IUPAC name of $CH_3 - CH_2 - CH_2 - CH_2 - CH_3$ is Br VII - CLASS 79

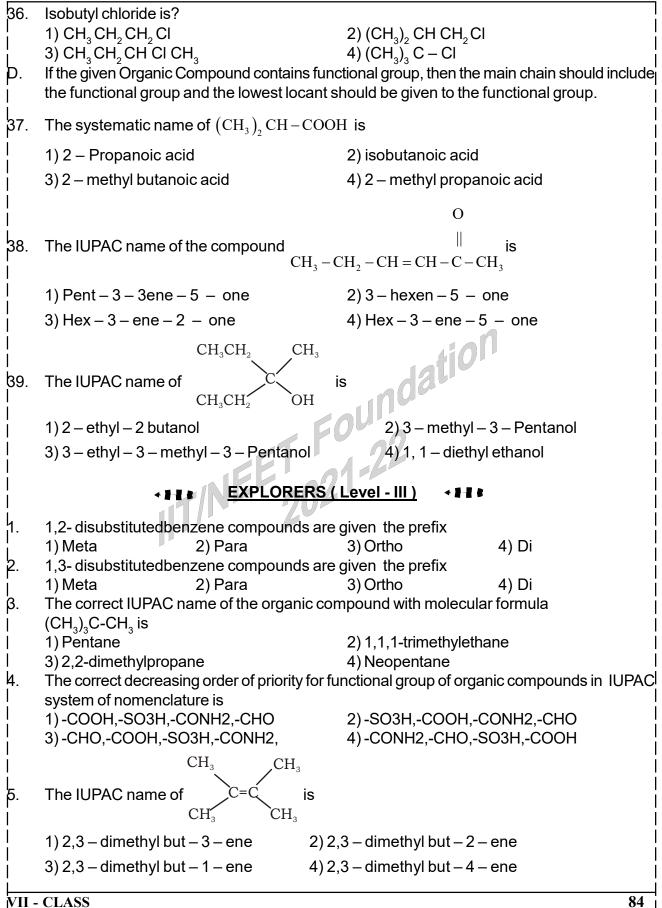
CHI	LIVIISIKY				Organic Chem	ustry
	1) 1–Bromo pe	entane	2) 3-	-Bromo pentane		
	3) 2–Bromo pe	entane	4)4-	-Bromo pentane		
Ι β1.	The IUPAC na	me of $CH_3 - CH_2$	$-CH_2 - CH_1$	$I - CH_2 - CH_2 - CH_3$	is	
			CH	$-CH_3$		
			 СН	3		
İ	1) 4–(1 ¹ –meth	yl ethyl) heptane	2)4-	–(21–methyl ethyl) he	eptane	
	3) 4–(3 ¹ –meth	yl ethyl) heptane	4)4-	–(4 ¹ –methyl ethyl) he	eptane	
		LEAR	NER'S TAS	SK- I		
İ		◆∎∎ ◆ <u>BEGINI</u>	NERS (Le	<u>vel-l)</u> ◆∎∎≯		
μ.	An example of	f alicyclic compour	nd is		4	
l L	1) Benze	,	exane	3) Cyclohexane	4) Furan	
2.		ollowing is an aron			4 \ 4 11	
۱ ß.	1) Benze	,	nthracene	3) Pyridine	4)All	
6. 	1) Benze	ollowing is an aron		3) Pyridine	4)All	
 4.	,	for unsaturated hy				
	1)-ane	2)-e		3)-yne	4) Both 2&3	
Ι þ.	,	ave a general form		-, ,	,	
	1) $C_{n}H_{2n}$		H_{2n+4}	3) $C_{n}H_{2n-2}$	4) $C_{n}H_{2n+2}$	
6.		ave a general form		, u 5u-5	, u 50+5	
		2) C		3) C _n H _{2n-2}	4) $C_{2n}H_{2n+1}$	
۲.		ollowing statemen				
İ	•	name of alkenes e				
	•	name of alkynes e name of alkanes e		-		
	4) All of these					
8.	Formula of alk	ane		AC name of alkyl rac	dical formed	
	i) CH ₄		p)	Butyl		
Ì	ii) C_2H_6		q)	Methyl		
	iii) C ₃ H ₈		r)	Ethyl		
į	iv) $C_4 H_{10}$	(::)	s)	Propyl		
	(i) 1) p	(ii)	(iii) r	(iv)		
	1) p 2) a	q	r	S		
	2) q - CLASS	r	S	р		80
1						

CHI	CHEMISTRY Organic Chemistry								
	3) s	r	р	q					
	4) p	r	S	q					
9. 	IUPAC names of $CH_3 - CH_3 - CH_3 - CH_3$ is: $\begin{matrix} \downarrow \\ C_2H_5 \end{matrix} = \begin{matrix} \downarrow \\ C_2H_5 \end{matrix}$								
	1) 3, 4 dimethyl hexane2) 2, 3 diethyl butane3) 2, ethyl 3-isopropyl propane4) 2, 3 dimethyl heptane								
10.	The hydrocarbon residue derived by removing a hydrogen atom from an alkene is called : 1) Alkenyl group 2) Alkyle group 3) Alkynyl group 4) Aryl group								
1 11.	,	ated hydrocarbons v	with $C = C$ are ca		4) None				
12. 	The correct of 1) Primary pro 2) Secondary	order of arrangemer efix+Rootword+Prin / prefix+Primarypref / prefix+Rootword+F	nt of rootword, s narysuffix+Secc ïx+Rootword Pr	uffixes and prefixe ondarysuffix+Seco imarysuffix+Seco	es is ondaryprefix . ndarysuffix.				
h3.	CH_3COCH_2C	CN has the IUPAC n	ame	4:01					
1	1) 3–oxo–but	ane nitrile		2) 1–cyando pro	pane				
i	3) 2–oxo proj	pane	4) 1	-cyano butanone	İ				
14.	-	prop–2–ynal is :	COU						
1	$1) CH_3 - C \equiv$	C – OH		$CH \equiv C - CH_2OH$					
i	3) $CH \equiv C - C$	СНО	4) ($CH_2 = CH - CHO$	İ				
			VERS (Level -	´ <u>)</u>	►				
U.	Multi Correc	ct Choice Type:	60		l				
<u>1</u> 5.		following is/are aror	-						
l ko	1) Furan	2) Pyrrole	,	hiophene	4) None				
16. 	1) ane	(for unsaturated hyd 2) ene	arocarbons 3) y	ne	4) none				
1 17.	,	following is/are com	, .		+)none				
1	1) Acetic acio	-		ormic acid	4) 1-butene				
18.		nd $(CH_3)_2 - C(OH)$ -			ļ				
	•	butan – 2 – ol	,	soamyl alcohol	. !				
 ИО	3) Ethyl dime	-	4) te	ert – pentyl alcoho					
19. 	1) Alkynes	general formula of 2) Alkenes	3) (Cyclo Alkanes	4) Cyclo Alkenes				
20.	, .	correct for a homol	,						
	1)All membe	rs have a general fo	ormula.						
İ	2) All membe	ers have similar chei	mical properties	3.	ļ				
	3) All membe	rs have same physi	ical properties.						
	4) All membe	rs have same functi	onal group.						
			-						
	CLASS								



CHEMISTRY **Organic Chemistry** V. **Comprehension Type:** Depending on the nature of the ring carbon compounds are divided into carbo cyclic and Ά. heterocyclic compounds. Which of the following is an example of alicyclic homocyclic compound? 28. 1) Benzene 2) Pentane 3) Cyclo pentane 4) Furan 29. Which of the following are aromatic heterocyclic compound? 1) Furan 2) Pyrrole 3) Thiophene 4) All 80. Which of the following is/are alicylic heterocyclic compound? 2) Pyrrolidine 1) Pyrrole 3) Benzene 4) Butane Β. The parent carbon chain is numbered in a manner so as to give lowest number to that carbon atom linked by double (or) triple bond even if it Violates the rules of saturated hydrocarbons. CH₃ β^{I} . The IUPAC name of $CH_3 - C \equiv C - CH_2 - CH - CH_3$ is 2) 2 – Methyl – 4 – Hexyne 1) 5 - Methyl - 2 - Hexyne3) 2 - yne - 5 - Methyl Hexane4) 1,1 – Dimethyl – 3 – Pentyne The IUPAC name of $CH_2 = C - CH = CH_2$ is β2. CH₃ 2) 2 - Methyl buta - 1,3 - diene 1) 3 - Methyl buta - 1, 3 - diene3) Penta diene 4) 2 – Methyl pentene CH₂ CH, B3. The IUPAC name of is CH. CH 1) 2,3 - dimethyl but - 3 - ene2) 2,3 - dimethyl but - 2 - ene4) 2,3 - dimethyl but - 4 - ene3) 2,3 – dimethyl but – 1 – ene C. According to IOPAC system of nomenclature; the longest possible continous chain of carbon atioms containing the funthonal group and carbon - carbon muthple bonds is first selected and the root word corresponding to it is noted. 34. IUPAC name of $H - \stackrel{|}{C} - \stackrel{|}{C} - \stackrel{|}{C} - Cl$ is 1) 1,2 - dichloro ethane 2) 2,2 - dichloro ethane 3) 1,1 - dichloro ethane 4) Dichloro ethane IUPAC name of $CH_2 = CH - CH(CH_3)_2$ is β5. 1) 1,2 - dimethyl - 2 - propene 2) 3 - methyl - 1 - butene 4) 1-isoprophyl ethylene 3) 2. vinyl propane VII - CLASS 83

Organic Chemistry





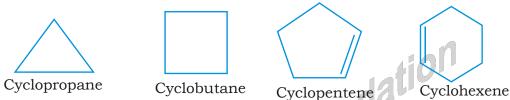
Nomenclature of (A) Cyclo-alkanes, Cyclo-alkenes, Cyclo-alkynes Nomenclature of alicyclic compounds containing functional groups

Nomenclature of Alicyclic compounds : These compounds contain one or more rings of three or more carbon atoms and resemble aliphatic compounds in their characteristics. These are, therefore, called aliphatic cyclic or alicyclic compounds

Monocyclic compounds :

- The name of alicyclic compounds are based on the following rules.
- 1. The names of the alicyclic compound are obtained by adding there primary prefix 'cyclo' to the word root that corresponds to the number of carbon atoms in the ring.

For the cyclic compounds containing all single bonds in the ring, primary suffix 'ane' is added to the word root. For those containing one double or triple bond, the primary suffix ene or yne is added.



2. If only one substituent is attached to the ring, its position is note mentioned. If two or more substituents are present, their positions are indicated by arabic numerals i.e., 1,2,3,4 etc. which are used for numbering the carbon atoms in the ring. The numbering is done in such a way (clockwise or anticlockwise) that the substituents get the lowest set of locants. All other rules relating to aliphatic or acyclic compounds are then followed. For example.

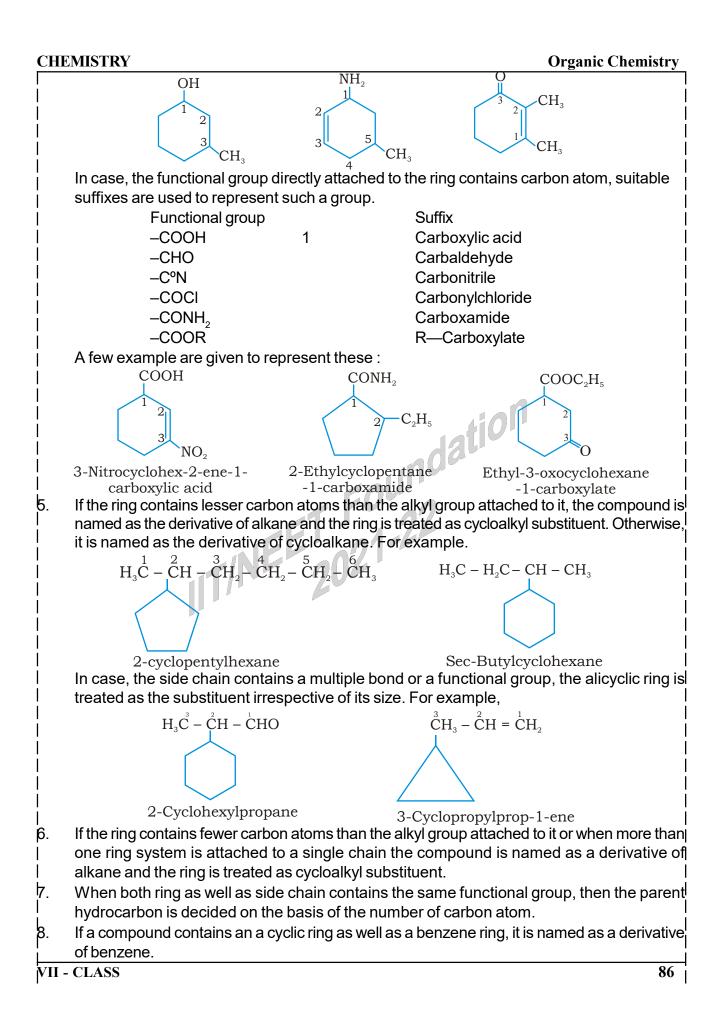


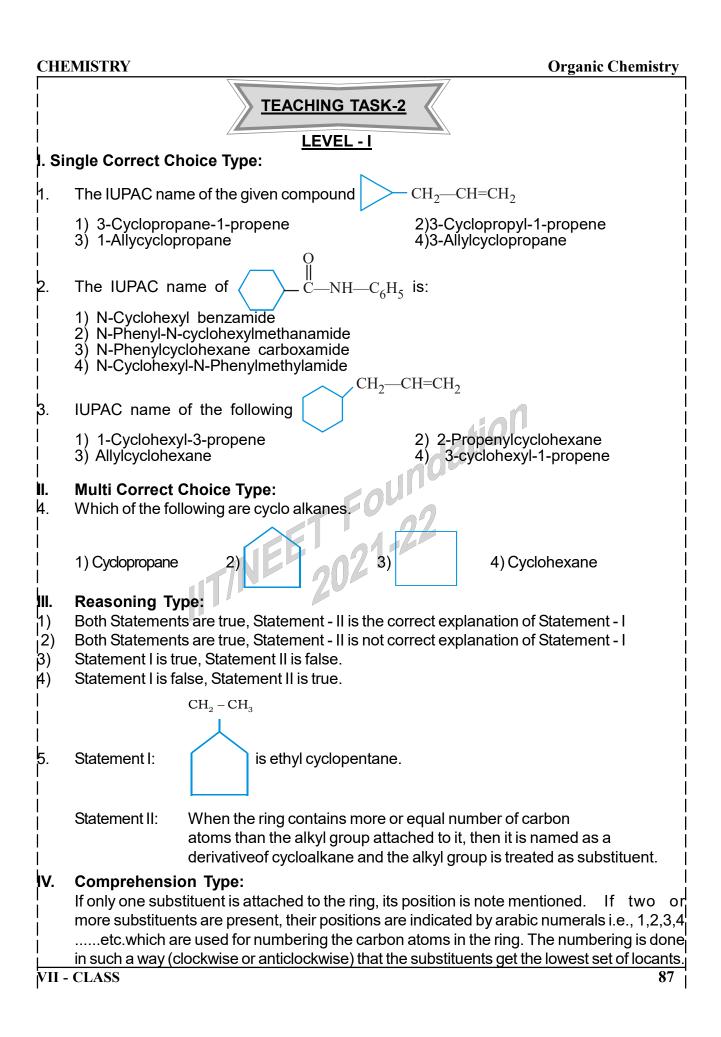
- Methylcyclopentane 1-Ethyl-2-Methylcyclopentane
- B. If a multiple bonds and some other substituents are present in the ring, the numbering is
 done in such a way so as to assign lowest number to the multiple bond. For example



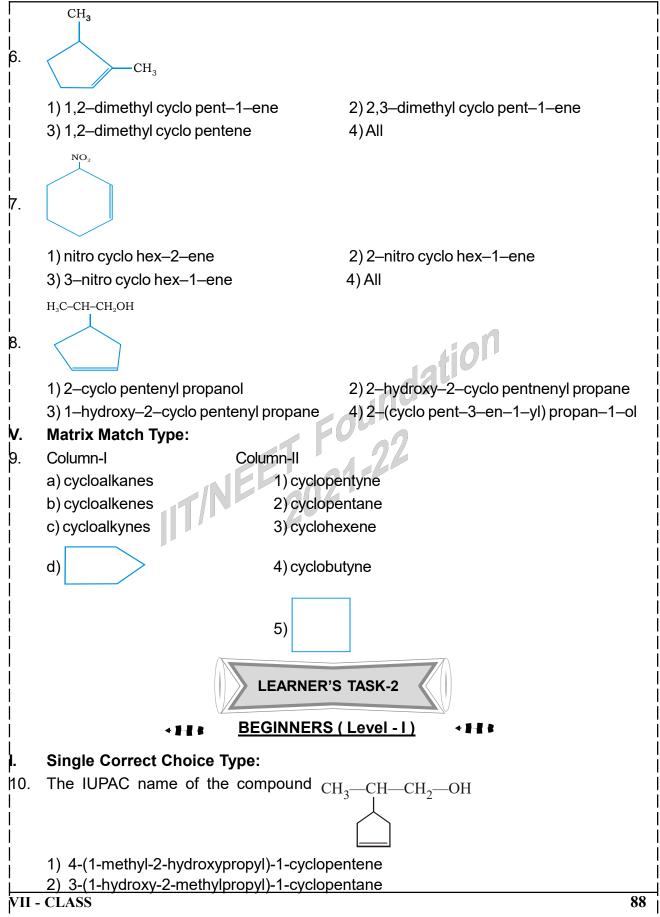
3-Nitrocyclohex-1-ene

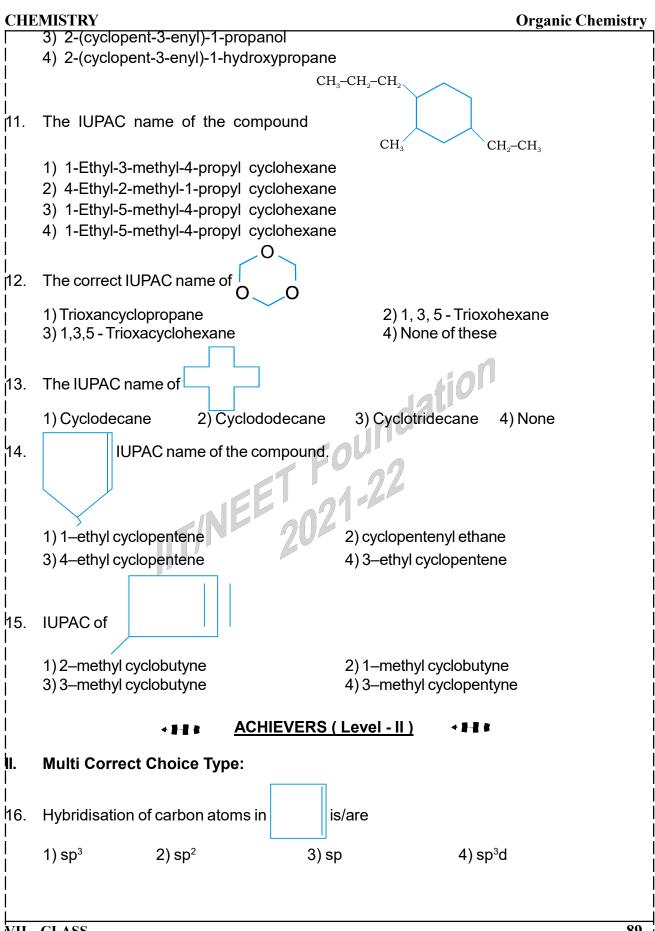
In case, some functional group along with some substituents are present in the ring, the
 numbering of the carbon atoms should be done in such a way so that the functional group
 gets lowest number.

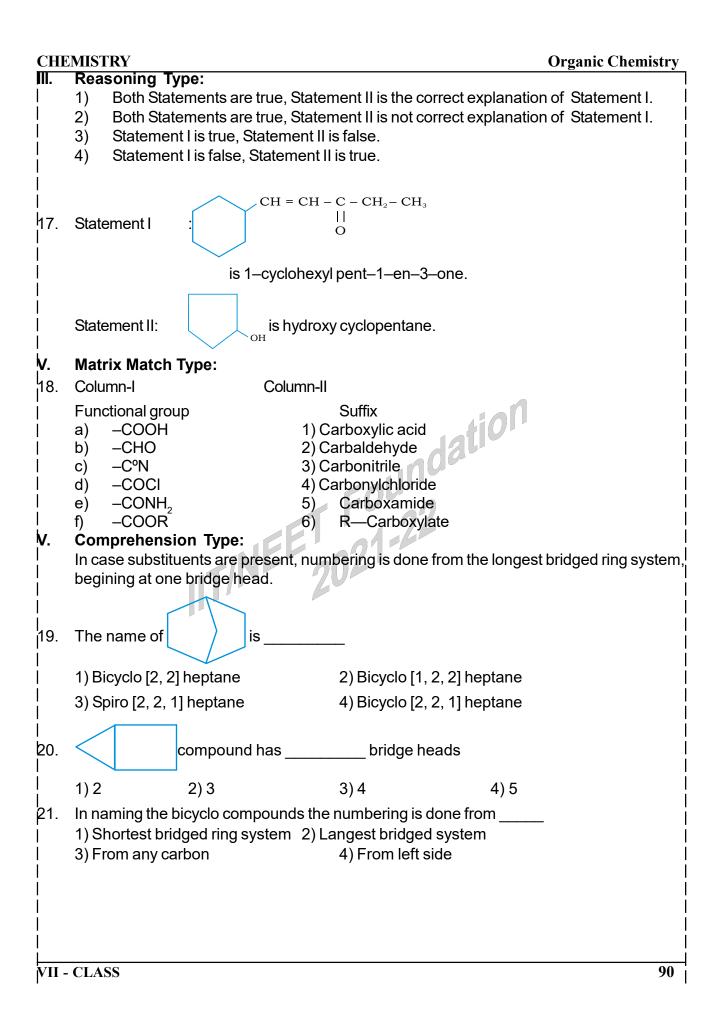




Organic Chemistry







İ	* 1	EXPL	<u>ORERS</u>	<u> 6 (Level - III)</u>	∢ ∦-	8	
bes	SCRIPTIVE TYPE						
կ. 	 Write the cyclic com 1) 2-methyl cyclobu 3) 3-methyl cyclobu 4) 3-methyl cyclope 2) 3-(1-hydroxy-2-n 3) 2-(cyclopent-3-e 	tyne tyne ntyne)4-(1-me nethylpropyl)-1	2) thyl-2-h I-cyclop	1-methyl cyclok ydroxypropyl)-1	outyne	pentene	
	4) 2-(cyclopent-3-e					_	
2.	Write the rules for	writing IUPA	.C nam	ies Of cyclic o	compo	unds.	
 	×1	∎∎≁ <u>Res</u> e	ARCH	ERS (Level -	<u>IV)</u>	┽┠╂╏ ≯	
Н.	Which of the followi	ng compounds	has wr	ong IUPAC nar	ne		(AIEEE2002
	1) CH3-CH2-CH2-C	COO-CH2CH3	Ethyl	butanoate			
į	2) CH3-CH(CH3)-C	H2-CHO	3-Me	thylbutanal	110	n	
	3) CH3CH(OH)-CH	(CH3)-CH3	2-Me	thyl-3-butanol			
į	4) CH3CH(CH3)-C	O-CH2-CH3	2-Me	thyl-3-pentanor	ne		
1 2.	The functional grou	p which is foun	d in Am	iino Acid is			(AIEEE2008
İ	1)-COOH	2)-NH2	ſſ	3)-CH3	4) bot	th 1&2	
β.	In which of the follow	ving species is	the un	der lined carboi	n havin	ig sp3 hy	bridisation
	ď		200				(AIEEE2008
	1) CH3COOH	2) CH3C H2	20H	3) CH3COC	CH3	4) CH3	=CH-CH3
4.	In allene (C3H4), the	e type of hybric	lisation	of the carbon a	atom is		(IIT-JEE -2011
	1) sp & sp3	2) sp & sp2		3) only sp3		4) sp2	& sp3
5.	An isomer of ethanc	l is(EAMCET-	93)(IIT-	·86)			
	1) Methanol	2) Diethyl et	her	3)Acetone		4) Dime	ethyl ether
6.	n-Propyl alcohol & i	so propyl alco	hol are	example ofiso	omerisi	m	(AIEEE2012
	1) Position	2) Chain		3) Functional		4) Taute	omers
ל. ו	The IUPAC name of 1) 2,2-Dimethylprop	•		ylpropane			(AIEEE2009
İ	3) 2,2-Dimethylbuta	ne 4)2-	Dimeth [,]	ylbutane			
į							
İ							

Organic Chemistry



<u>ΦΦ TEACHING TASK - I :</u>

3-2 4-1 5-3 2-4 6-3 7-1 8-2 9-1 10-2 11-2 12-3 1-2 13-1 14-3 15-3&4 16-1,2,3 17-1,2,3,4 18-1,2,3 19-1,3,4, 20-1,2,4 21-B 22-B 23-3,4,1,2 24-2,3,4,1 25-2 26-1 27-1 28-2 29-1 30-4 31-432-2 33-1. $\Phi\Phi$ LEARNER'STASK - I : 12-j 2-4 4-4 5-3 6-1 8-4 1-3 3-4 7-4 9-1 10-1 11-2 213-1 14-3 15-1,2,3 16-2,3 17-1,2,3 18-2 19-2,3 20-2,3,4 21-1 22-1 23-3 24-2,1,3,4 25-2,1,4,3 26-4,3,1,2 27-2,3,4,1 28-329-4 30-2 31-1 32-2 33-2 34-3 35-2 36-2 37-2 38-3 39-2. EXPLORERS : 1-3 3-3 4-1 5-2 2-1 $\Phi\Phi$ TEACHING TASK-2 : 4-1,2,3,4 5-1 6-2 9-5,3,4,2 1-2 2-2 3-4 7-2 $\Phi\Phi$ LEARNER'STASK - 2 : 11-1 12-1 13-2 14-4 15-3 16-1,217-2 18-1,2,3,4,5,6 19-4 20-2, 21-2 10-3 5-4 6-2 7-1 20 **RESEARCHERS**: 1-1 2-4 3-2 4-1