PAGE NO.: 1
DATE: / /

Organic Chemistry is the study of compds. containing carbon.

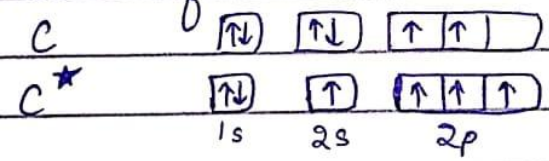
C

3 isotope

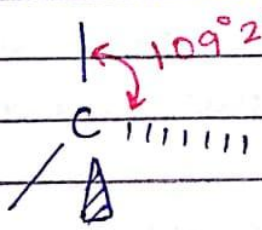
$^{12}_{6}C$	$^{13}_{6}C$	$^{14}_{6}C$
98.8 %	1.1 %	0.001 %
A = 12		half life = 5730 yrs
Z = 6		

Characterstic of C :-

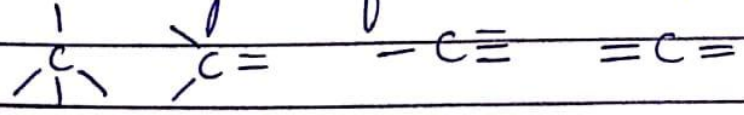
1. Tetravalent : C form 4 bond in excited state.



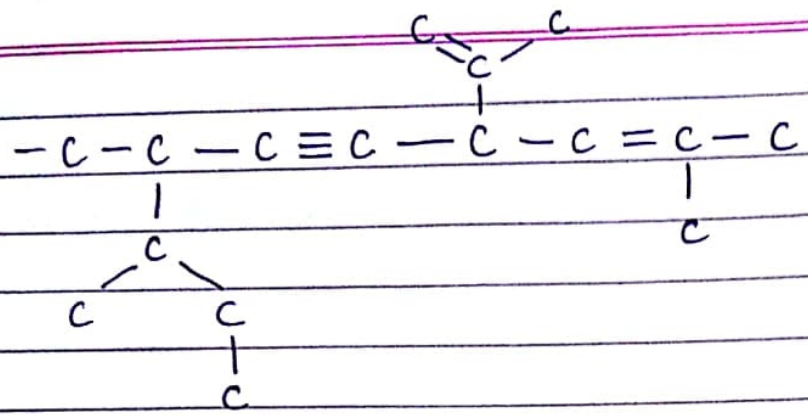
2. Tetrahedral : 4 bonds are directed towards corner of ~~of~~ tetrahr~~ed~~-on.



3. Tendency to form multiple bond :

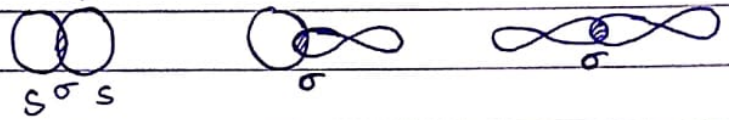


4. Catenation : Self linking property of C to form long chain.

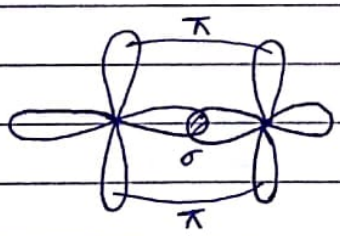


5. Hybridisation :-

σ bond : It is formed by coaxial overlapping of atomic orbitals.

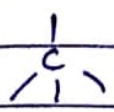


π-bond : formed by colateral / sidewise overlapping of atomic orbitals.

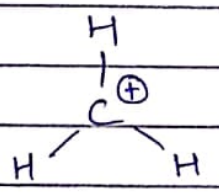


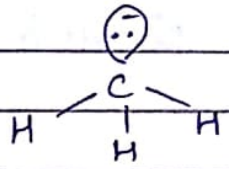
Note :- ① σ bond is stronger than π-bond becoz extent of overlapping is more in σ bond.

② π-bond is more reactive than σ-bond.

Str.	σ, π bond	hyb.	B.A.	Shape
	4, 0	sp^3	109° 28'	tetrahedral
	3, 1	sp^2	120°	trigonal planar
$-C\equiv$	2, 2	sp	180°	linear
$=C=$	2, 2	sp	180°	linear

S.No. = no. of σ bond + no. of lp.

CH_3^+  sp^2

CH_3^-  sp^3

E.N. ∝ % S-Char

	sp	sp^2	sp^3
% s char	$\frac{1}{2} \times 100$	$\frac{1}{3} \times 100$	$\frac{1}{4} \times 100$
	= 50 %	= 33.3 %	= 25 %

E.N. → sp (3.25) > sp^2 (2.75) > sp^3 (2.5)

B.L. $\propto \frac{1}{E.N.}$ $sp < sp^2 < sp^3$

B.S. ∝ E.N $sp > sp^2 > sp^3$

or

B.S. $\propto \frac{1}{B.L}$

Ques Find hyb. of underlined C ?

1. $CH_3-\underset{\underset{OH}{|}}{\underline{C}H}-CH_2-\underset{\underset{sp^2}{|}}{\underline{C}H}=CH_2$ ($\underline{C}H$: sp^3)

2. $CH_3-CH_2-\underline{C}\equiv N$ ($\underline{C}$: sp)

3. $CH_3-CH=CH-\underset{\underset{sp^3}{|}}{\underline{C}H_2}-\underset{\underset{O}{||}}{\underline{C}}-CH_3$ ($\underline{C}H_2$: sp^3; $\underline{C}$: sp^2)

4. $CH_3-\underset{\underset{O}{||}}{\underline{C}}-CH_2-\underline{C}OOH$ ($\underline{C}$: sp^2; $\underline{C}OOH$: sp^2)

5. $CH_3-\underset{\underset{OH}{|}}{\underline{C}}=CH-\underline{C}H_2-Cl$ ($\underline{C}$: sp^2; $\underline{C}H_2$: sp^3)

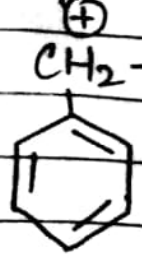
Ques Find hybridisation of C-C single bond

$HC\equiv\underset{\underset{sp}{|}}{\underline{C}}-\underset{\underset{sp^2}{|}}{\underline{C}H}=CH_2$

Ques Find hyb. of C_2 & C_3 in following

$HC\overset{6}{\equiv}\overset{5}{C}-\overset{4}{C}H_2-\overset{3}{C}H_2-\overset{2}{H}C=\overset{1}{C}H_2$ (C_3: sp^3; C_2: sp^2)

Ques Find hyb. of carbonium ion ?



$CH_2 \rightarrow sp^2$

Ques find no. of σ & π-bond in following?

1. $CH_3 - CH_2 - CH_2 - OH$ 11σ

2. $CH_3 - CH_2 - CH = CH_2 - OH$ 13σ, 1π

3. $CH_3 - CH_2 - CN$ 8σ, 2π

4. $CH_3 - CHO$ 6σ, 1π

5. $CH_3 - COOH$ 7σ, 1π

6. $CH_3 - \underset{\underset{O}{\|}}{C} - CH_2 - COOH$ 12σ, 2π

7. $CH_3 - \underset{\underset{OH}{|}}{C} = CH - COOH$ 12σ, 2π

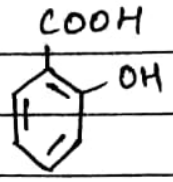
NT 8. $CH_2 = CH - C \equiv CH$ 7σ, 3π

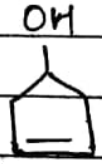
NT 9. $CH \equiv C - CH_2 - C \equiv C - CH_3$ 11σ, 4π

ANIMS 10. $CH_2 = CH - CH_2 - CN$ 9σ, 3π
allyl cyanide

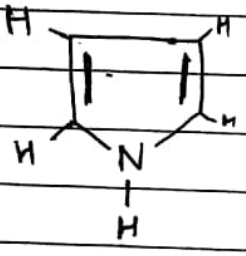
IIT 11. Toluene  15σ, 3π

DEC 12. Styrene  16σ, 4π

13.  $16\sigma, 4\pi$

14.  $14\sigma, 1\pi$

15.  $23\sigma, 2\pi$

16.  $10\pi, 2\pi$

Ques Arrange following in ↑ing order of B.L.?

$\underset{x}{CH \equiv CH}$ $\underset{y}{CH_3 - CH_3}$ $\underset{z}{CH_2 = CH_2}$

① $\underset{sp}{CH \equiv CH}$ $\underset{sp^3}{CH_3 - CH_3}$ $\underset{sp^2}{CH_2 = CH_2}$

B.L. $\propto \frac{1}{E.N}$

$y > z > x$

② $\underset{sp}{CH \equiv C - \overset{a}{H}}$ $\underset{sp^3}{CH_3 - CH_2 - \overset{b}{H}}$ $\underset{sp^2}{CH_2 = CH - \overset{c}{H}}$

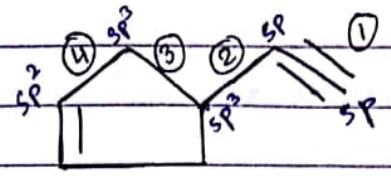
$b > c > a$

③ $CH_3 - \overset{\alpha}{CH} = CH_2$ $CH_3 - \overset{\beta}{C} \equiv CH$ $CH_3 - \overset{\gamma}{CH_2} - CH_2$

$\gamma > \alpha > \beta$

IIT

(4)



~~2 = 3 > 4 > 1~~

3 > 4 > 2 > 1

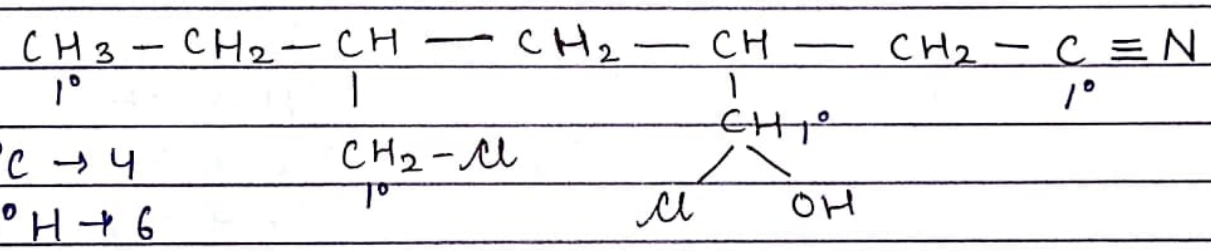
B.S. ① > ② > ④ > ③

Type of Carbon and Hydrogen :-

1. Primary Carbon & Primary Hydrogen :-

1°C → C which is attached to only 1° carbon.

1°H → H which is attached to only 1° C.



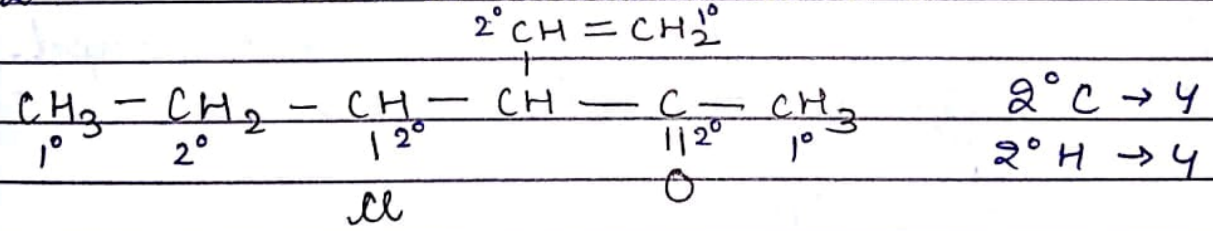
1°C → 4

1°H → 6

2. Secondary Carbon & Secondary Hydrogen :-

2°C → C which is attached to 2° carbon

2°H → H which is attached to 2° C.



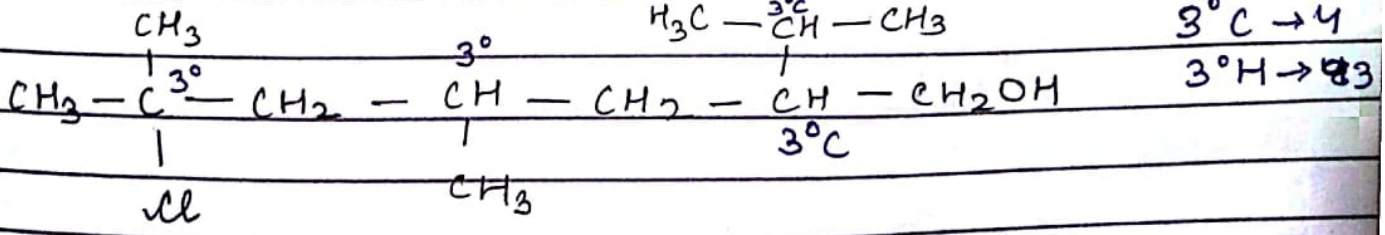
2°C → 4

2°H → 4

3. Tertiary Carbon & Tertiary Hydrogen :-

3°C → C which is attached to 3 carbon.

3°H → H which is attached to 3° C.

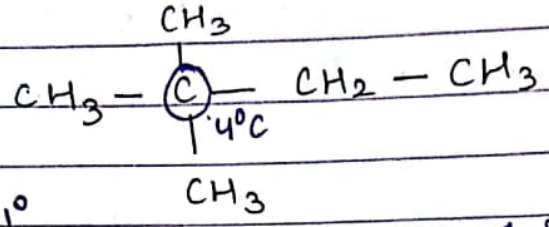


3°C → 4

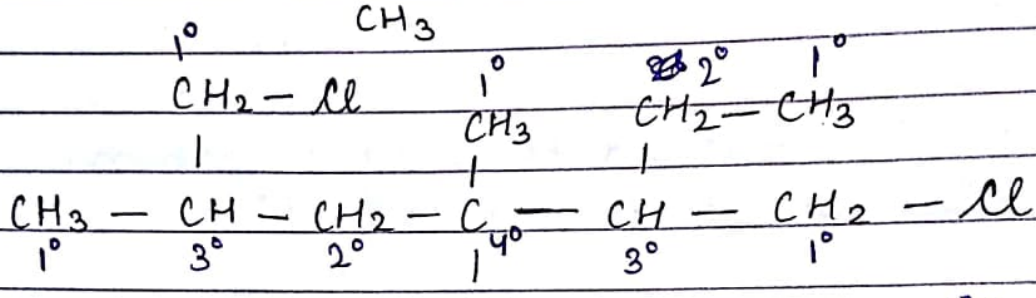
3°H → 13

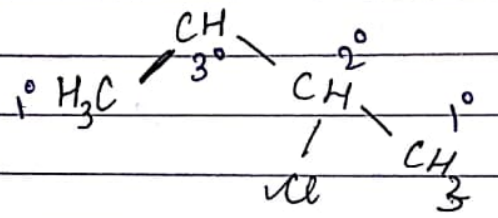
4. Quaternary Carbon :-

4°C → Carbon which is attached to 4 carbon atom



ex





1°C → 7C
1°H → 19H
2°C → 3C
2°H → 5H
3°C → 3
3°H → 3
4°C → 1

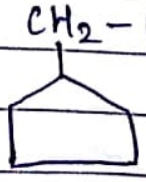
Type of Alcohol :-

1. Primary Alcohol :- (1° OH) The alcoholic compd. in which OH gp. is attached to 1°C.

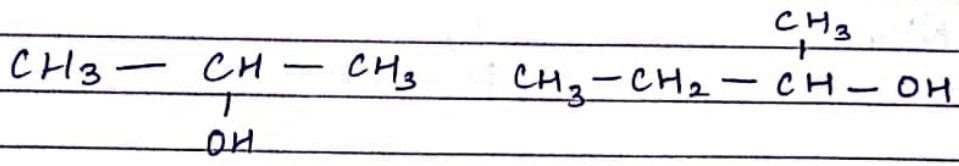
CH_3-CH_2-OH

$CH_3-CH_2-CH_2-OH$

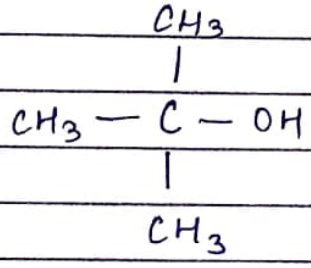
CH_3-OH



2. Secondary Alcohol : The compd in which OH gp. is attached to 2°C.

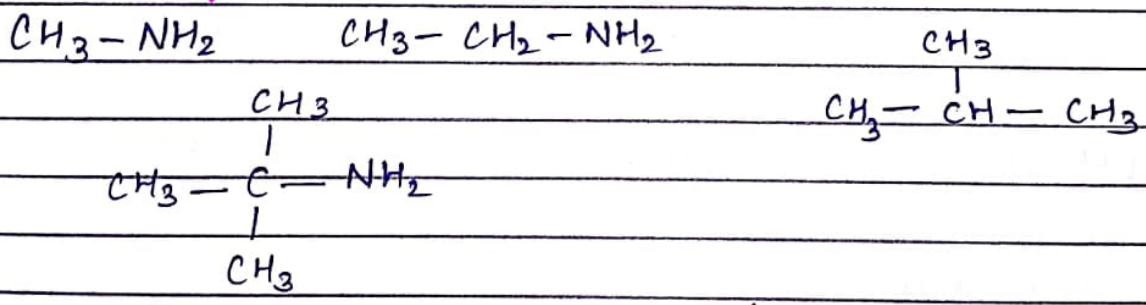


3. Tertiary Alcohol : The alcoholic compd. in which OH gp is attached to 3°C

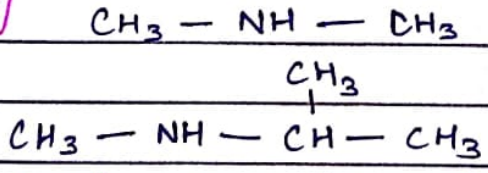


Type of Amine :-

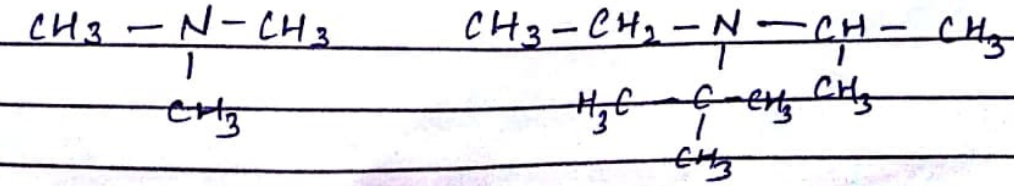
1. Primary Amine : N is attached to only 1 C atom.



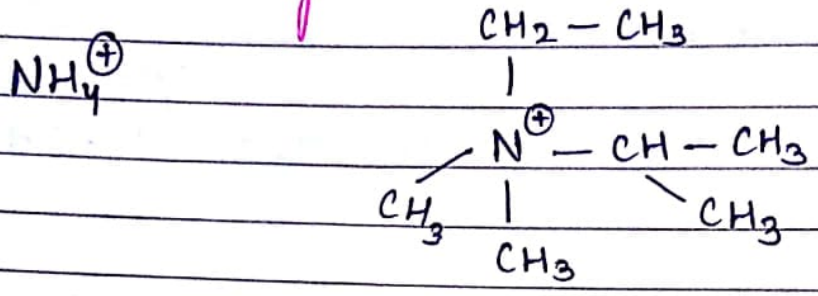
2. Secondary Amine :- N is attached to 2 C atom.

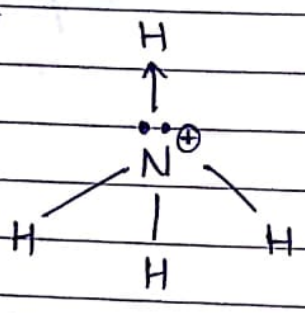


3. Tertiary Amine :- ~~in which~~ N is attached to 3 C atom.

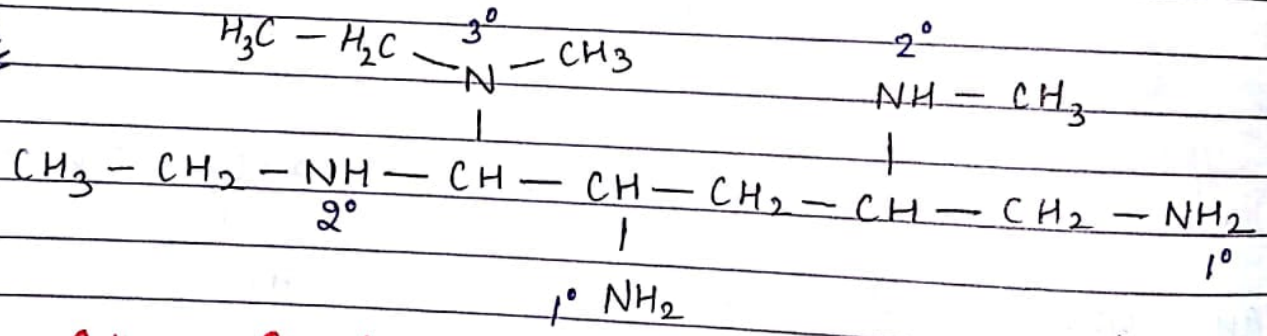


4. Quaternary amine : N is attached to 4C atom.





ex



Classification and Nomenclature

Classification :-

1. Based on structure
2. Based on Group
3. Based on Homologous

① Based on structure :-

- Open chain / aliphatic / acyclic
  - Saturated
    - C–C
    - + alkanes
    - C_nH_{2n+2}
    - Paraffins
    - $CH_3-CH_2-CH_3$
    - $CH_3-\overset{\overset{O}{\|}}{C}-OH$
    - $CH_3-CH=O$
  - Unsaturated
    - C=C
    - Olifinic bond
    - alkene
    - C_nH_{2n}
    - C≡C
    - acetylinic bond
    - alkyne
    - C_nH_{2n-2}
- Closed chain / Cyclic
  - Homocyclic
    - alicyclic
    - aromatic
  - Heterocyclic
    - Alicyclic
    - Aromatic
    - Pyridine
    - Pyrole

Ques which of the following is unsaturated compo

1. C_3H_8 ×
2. CH_3-CHO ×
3. CH_3-COOH ×
4. C_3H_4
5. $CH_3-CH_2-CH_2-OH$

6. $CH_3 - CH = CH - CH_3$

7. $CH_3 - \underset{\underset{O}{\|}}{C} - NH_2$

8. $CH_3 - \underset{\underset{O}{\|}}{C} - O - CH_3$

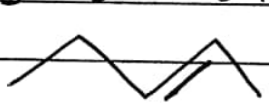
9. $CH_3 - C \equiv CH$

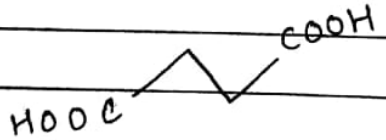
10. $CH_3 - CH_2 - \underset{\underset{\underset{O}{\|}}{C - O - CH_3}}{\overset{|}{CH}} - CH_3$

11. $CH_3 - CH_2 - CH = \underset{\underset{Cl}{|}}{C} - CH_3$

12. $CH_3 - CH_2 - CH_2 - NH_2$

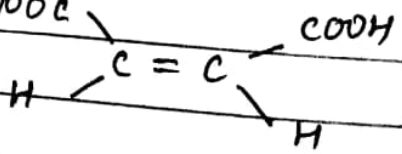
13. $CH_3 - CH = CH - CH_2 - NH - CH_3$

14. 

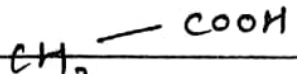
15. 

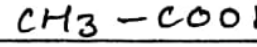
16. $CH_2 \equiv CH$

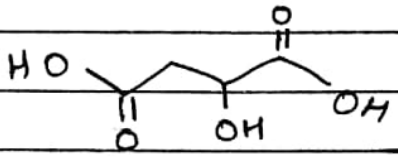
Ques Which of the following is unsaturated?

1. Maleic acid 

2. Succinic acid $\begin{matrix} CH_2 - COOH \\ | \\ CH_2 - COOH \end{matrix}$

3. Malonic acid 

4. acetic acid 

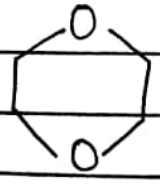
5. Malic acid 

A :- Acetic acid is not an unsaturated compd. ✓
R :- acetic acid do not have olefinic ~~comp~~ bond. (2)

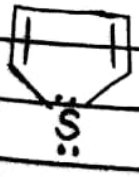
A :- acetic acid is saturated compd. ✓
R : acetic acid have  bond. ✓ (2)

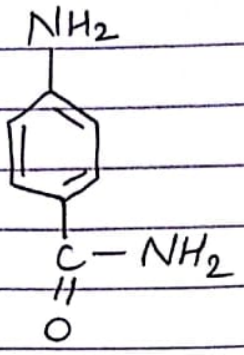
A : acetic acid have  bond.
R : acetic acid is saturated compd. (2)

Ques find no. of heteroatoms in following ?

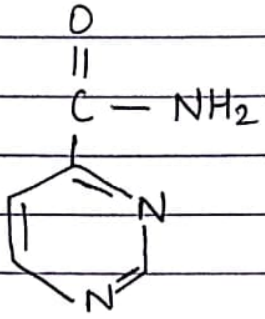
(1)  2

2.  1

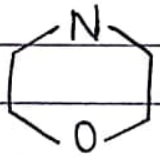
3.  1

4. 

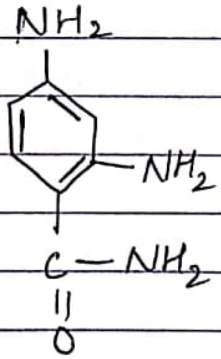
O

5. 

2

6. 

2

7. 

O

(2) Classification based on Group :-

$R-H \xrightarrow[-H]{+G_1} R-G_1$

Hydrocarbon → hyd.-radical → (G1) func^n gp.

phy. property → chemical property

B.P., M.P. & density etc

- $-OH$
- $-NH_2$
- $-X$
- $-COOH$
- $-CH=O$

Pentyl = amyl

PAGE NO.: 15
DATE: / /

Hydrocarbon Radical (R–) :-

Saturated :

Monovalent :-

alkane $\xrightarrow{-H}$ alkyl

CH_4 $\xrightarrow{-H}$ CH_3-

Methane — Methyl

CH_3-CH_3 $\xrightarrow{-H}$ CH_3-CH_2-

ethane — ethyl

$CH_3-CH_2-CH_3$ (Propane) $\xrightarrow{-H}$ $CH_3-CH_2-CH_2-$ n-propyl

$CH_3-CH_2-CH_3$ (Propane) $\rightarrow$ $CH_3-\underset{|}{CH}-CH_3$ iso-propyl

Normal :- • Straight chain hydrocarbon
• no branching

eg

$CH_3-\underset{|}{CH_2}-CH_3$

n-butane

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

n-hexane

$CH_3-\underset{|}{CH_2}$

propane

n-pentane

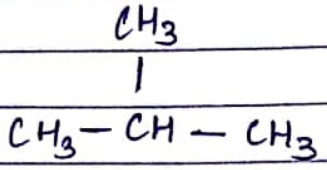
$CH_3(CH_2)_5CH_3$

n-heptane

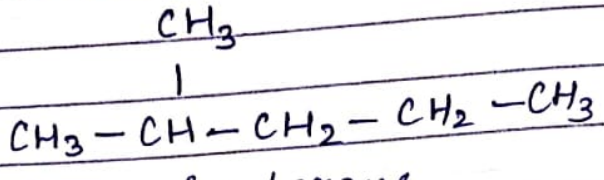
$CH_3-CH_2-CH_2-$ n-propyl

$CH_3-CH_2-CH_2-CH_2-$ n-butyl

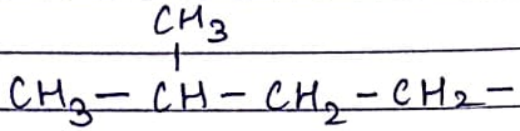
Iso : When methyl gp. is attached at 2nd last C or 2nd C (in case of alkanes) is c/a iso-comp.

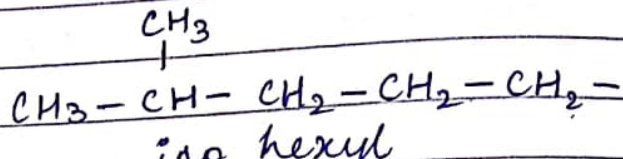


iso-butane

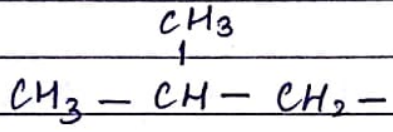


iso hexane

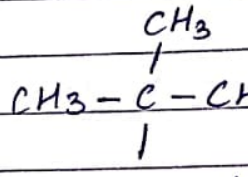




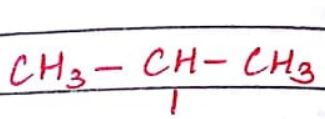
iso hexyl



iso-butyl

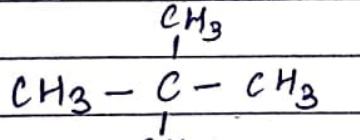


t-butyl

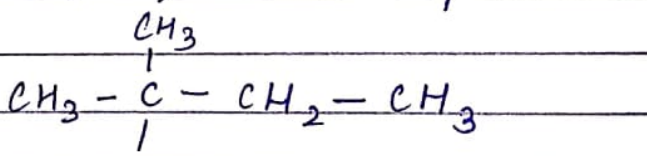


isopropyl

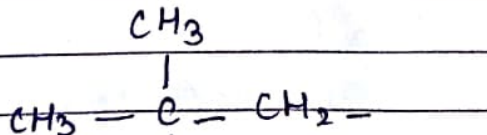
Neo :- When 2 methyl gp. is attached to 2nd last C or 2nd C (in case of alkane) c/a neo compd.



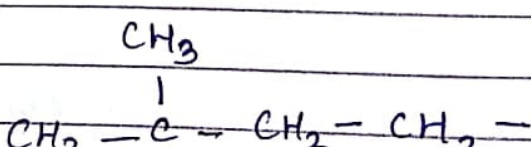
neo-Pentane



neo-hexane

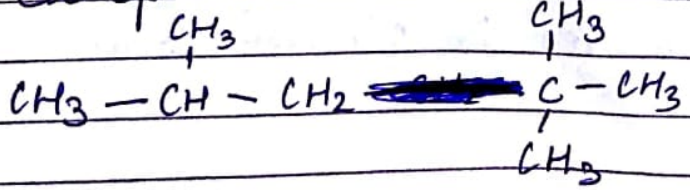


neopentyl

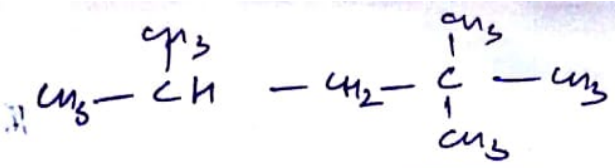


neo hexyl

Exception



Iso-octane



iso octane

PAGE NO.: 17
DATE: / /

$CH_3 - CH_2 - CH_2 - CH_2 - CH_2 -$ n-pentyl

$CH_3 - CH_2 - CH_2 - CH(CH_3) -$ Sec. pentyl

$CH_3 - CH(CH_3) - CH_2 - CH_2 -$ iso · pentyl

$CH_3 - C(CH_3)_2 - CH_2 -$ neo pentyl

Bivalent :- 2 valency free
from same C - atom → alkylidene
from vicinal C - atom → alkylene
terminal C - atom → polymethylene

$CH_4 \xrightarrow{-H} CH_3 - \xrightarrow{-H} CH_2 <$

Methylidene
or
Methylene

$CH_3 - CH_3 \longrightarrow CH_3 - CH_2 - \rightarrow CH_3 - CH <$ ethylidene

$CH_2 - CH_2$ ethylene

$CH_3 - CH_2 - CH_3 \longrightarrow CH_3 - CH_2 - CH <$ propylidene

$CH_3 - CH_2 - CH_2 -$ propylene

$CH_3 - C(<) - CH_3$ isopropylidene

$-CH_2-CH_2-CH_2-CH_2-$

$CH_2=CH_2 \longrightarrow CH_2=CH- \longrightarrow CH_2=C\overset{\cdot\cdot}{}$

Vinyl or ethenyl — ethenylidene

$CH_3-CH=CH_2 \longrightarrow CH_3-CH=CH-$ propenyl

$\longrightarrow CH_3-\underset{|}{C}=CH_2$ isopropenyl

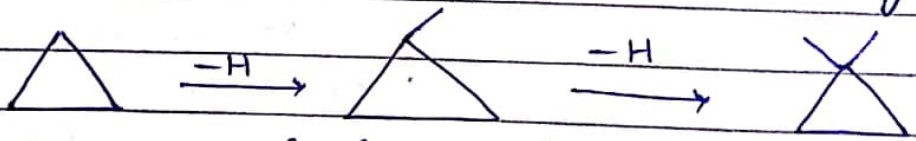
$\longrightarrow -CH_2-CH=CH_2$ Allyl

$CH\equiv CH \longrightarrow CH\equiv C-$ ethynyl

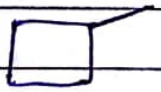
$CH_3-C\equiv CH \longrightarrow CH_3-C\equiv C-$ propynyl

$\longrightarrow -CH_2-C\equiv CH$ propargyl

$CH_3-CH=CH-CH_2-$ Crotyl



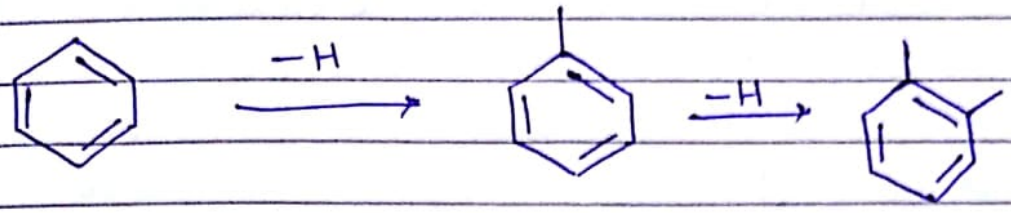
Cyclopropane — Cyclopropyl — cyclopropylidene



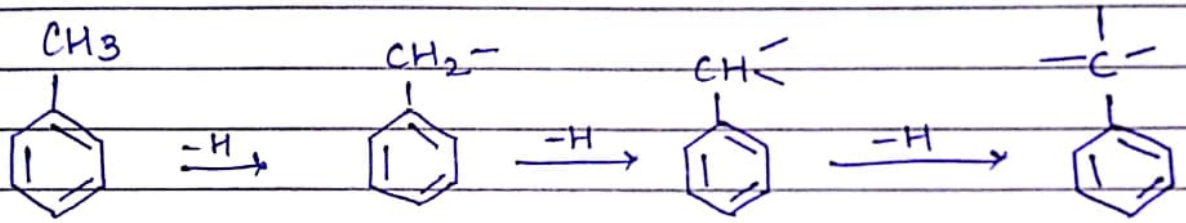
cyclobutyl

cyclopentyl

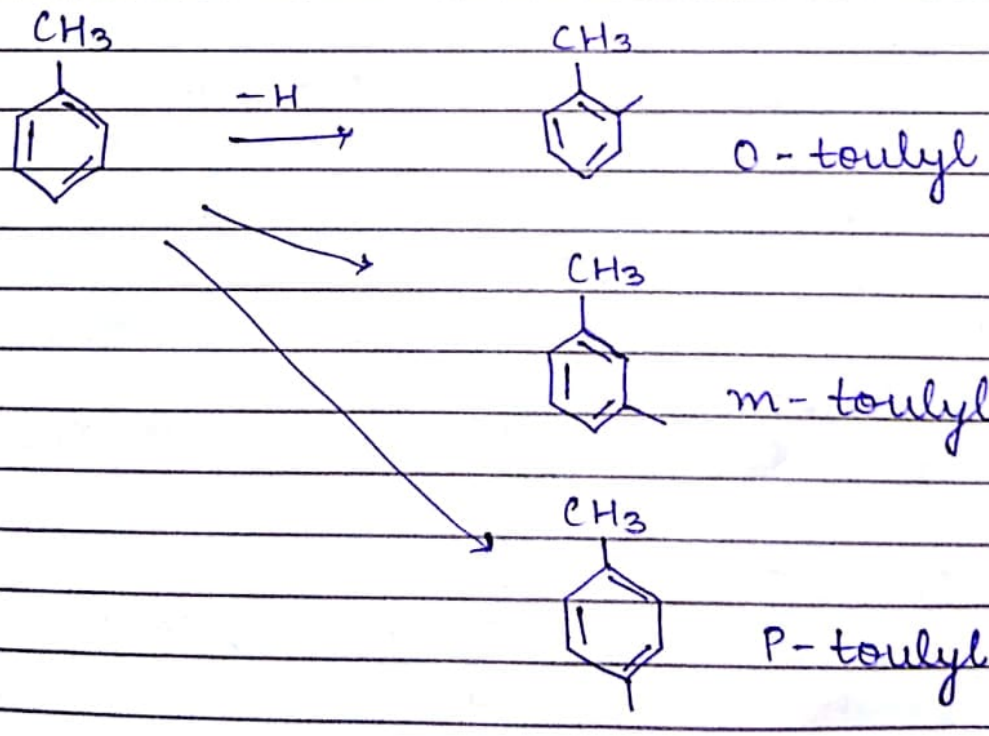
Aromatic :-



benzene or phene — phenyl — phenylene



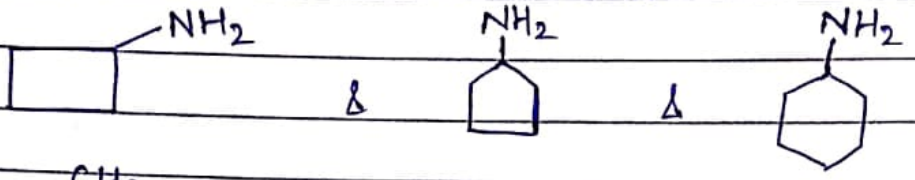
toluene — benzyl — benzal — benzo

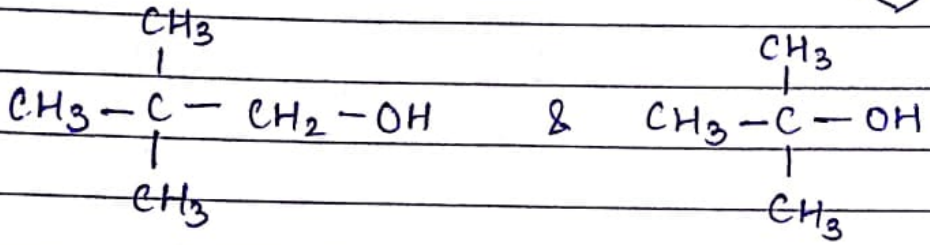


Concept of Homology :-

Those compd's which have same structural features as well as same general formula but differ by $-CH_2-$ group c/a homologs & series is c/a homologous series.

CH_3-CH_2-OH & $CH_3-CH_2-CH_2-OH$





	G.F.	
alkane	C_nH_{2n+2}	CH_4, CH_3-CH_3, $CH_3-CH_2-CH_3$
alkene	C_nH_{2n}	$CH_2=CH_2$, $CH_3-CH=CH_2$, $CH_3-CH_2-CH=CH$
alkyne	C_nH_{2n-2}	$CH\equiv CH$, $CH_3-C\equiv CH$
Alcohol	$C_nH_{2n+2}O$	CH_3-OH, CH_3-CH_2-OH, $CH_3-CH_2-CH_2-OH$
ether	C_nH_{2n+O}	CH_3-O-CH_3, $CH_3-O-CH_2-CH_3$
Halide	$C_nH_{2n+1}X$	
aldehyde or	$C_nH_{2n}O$	$H-\overset{\overset{O}{\|}}{C}-H$, $CH_3-CH=O$, CH_3-CH_2-CHO
Ketone	$C_nH_{2n}O$	$CH_3-\overset{\overset{O}{\|}}{C}-CH_3$
acid/ester	$C_nH_{2n}O_2$	$H-\overset{\overset{O}{\|}}{C}-OH$, CH_3-COOH

Characterstics of Homologues :-

1. Have same general formula i.e. same method of preparation.
2. Have same type of atom.
3. have same functional gp. hence similiar chemical property.
4. Have different physical property.
5. 2 successive members differ by CH_2- gp i.e. a weight of 14 unit.

$R-CH_3$ $\xrightarrow{-CH_2}$ $R-H$ (Lower Homologue)

$R-CH_3$ $\xrightarrow{+CH_2}$ $R-CH_2-CH_3$ (Higher Homologue)

Ques Write lower & higher homologue of isopentyl alcohol.

$$CH_3-\underset{\underset{CH_3}{|}}{CH}-CH_2-CH_2-OH$$

Higher $CH_3-\underset{\underset{CH_3}{|}}{CH}-CH_2-CH_2-CH_2-OH$

Lower $CH_3-\underset{\underset{CH_3}{|}}{CH}-CH_2-OH$ isobutyl alcohol

Ques Which of the following is pair of Homologues?

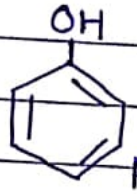
1. $CH_3-CH_2-CH_2-NH_2$ & $CH_3-\underset{\underset{CH_3}{|}}{CH}-NH_2$ X

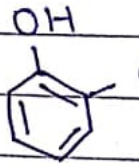
2. CH_3-CH_2-OH & $CH_3-CH_2-O-CH_3$ ✗

3. $H-\underset{\underset{O}{\|}}{C}-H$ & CH_3-CHO ✓

4. $CH_3-CH=CH_2$ & $CH_3-CH_2-\underset{\underset{H}{|}}{C}=CH_2$ ✓

5. CH_3-CH_2-Cl & CH_3-CH_2-Br ✗

6.  Phenol &  alcoholic ✗

7.  &  ✓

8. $CH_3-CH=CH_2$ &  ✗

9. $CH_3-CH_2-NH_2$ & $CH_3-CH_2-NH-CH_3$ ✗

10. $CH_3-\underset{\underset{O}{\|}}{C}-CH_3$ & $CH_3-CH_2-CH_2-CH=O$ ✗

11. $CH_3-CH_2-CH_2-OH$ & CH_3-O-CH_3 ✗

12.  &  ✓

13. $HCOOH$ & CH_3COOH ✓

14. $CH_3-\underset{\underset{O}{\|}}{C}-OH$ & $CH_3-O-\underset{\underset{O}{\|}}{C}-CH_3$ ✗

15. $CH_3-\overset{\overset{O}{\|}}{C}-CH_2-CH_3$ & $CH_3-\overset{\overset{O}{\|}}{C}-CH_3$ ✓

16. $CH_3-\underset{\underset{CH_3}{|}}{CH}-CH_2-Cl$ & $CH_3-CH_2-CH_2-Cl$ ✗

Nomenclature :-

~~Too~~ Common Name | Derived Name | IUPAC System

1. Common Name OR TRIVAL NAMING :-

- based on source of origin.

Structure	Source	Common Name
CH_4	Marshy area	Marsh Gas
CH_3-OH	distillation of wood	wood spirit
$HCOOH$	formica (Red Ant)	formic acid
CH_3-COOH	acetum (Vinegar)	acetic acid
$CH_3-\underset{\underset{OH}{|}}{CH}-COOH$	Milk	Lactic acid
C_3H_7-COOH	Butter	butyric acid
$HOOC-CH_2-\underset{\underset{OH}{|}}{CH}-COOH$	Apple	Malic acid
$HOOC-\underset{\underset{OH}{|}}{CH}-\underset{\underset{OH}{|}}{CH}-COOH$	~~tar~~ tamarind	tataric acid

$HOOC-CH_2-\underset{\underset{OH}{|}}{CH}-CH_2-COOH$ Citrus fruit Citric acid

(R) — (G) fun^c gp
hydrocarbon

Common Name
Functional Group

① Non-Terminating (with^{out} C containing func gp)
eg $-OH, -NH_2, -O-, -\underset{\underset{O}{\|}}{C}-$

② terminating (C containing func^c gp)
$-COOH$, $-CHO$

1. Non terminating

	Suffix
$-OH$	alcohol
$-SH$	thio alcohol
$-X$	Halide
$-NH_2$	amine
$-\underset{H}{N}-$	amine
$-\underset{|}{N}-$	amine
$-O-$	ether
$-S-$	thio ether
$-\underset{\underset{O}{\|}}{C}-$	Ketone

Polyvalent func gp:-
1. If same hydrocarbon gp is attached then di, tri, tetra etc used.
2. If diff. hydrocarbon are attached then write them in alphabatical order

Note: di, tri, tetra, primary, secondary, etc. are not considered in alphabatical order while terms iso, neo are considered in alphabetical order

CH_3-O-CH_3 — dimethyl ether

$CH_3-O-\underset{\displaystyle CH_3}{\underset{|}{CH}}-CH_3$ — isopropyl methyl ether

$CH_3-\overset{\displaystyle O}{\overset{||}{C}}-CH_3$ — dimethyl ketone

$CH_3-\overset{\displaystyle O}{\overset{||}{C}}-CH_2-CH_3$ — ethyl methyl ketone

CH_3-CH_2-OH — ethyl alcohol

$CH_3-\underset{\displaystyle CH_3}{\underset{|}{CH}}-Cl$ — isopropyl chloride

$CH_3-\underset{\displaystyle CH_3}{\underset{|}{CH}}-CH_2-NH_2$ — isobutyl amine

$CH_3-\underset{\displaystyle CH_3}{\underset{|}{\overset{\displaystyle CH_3}{\overset{|}{C}}}}-CH_2-OH$ — neo pentyl alcohol

$CH_2=CH-Cl$ — Vinyl chloride

$CH_2=CH-CH_2-OH$ — Allyl chloride

$CH_3-NH-CH_3$ — dimethyl amine

$CH_3-\underset{\displaystyle CH-CH_3}{\underset{|}{\overset{\displaystyle CH_3}{\overset{|}{N}}}}-CH_2-CH_3$ — ethyl isopropyl methyl amine

Terminating func^{c} gp :-

Prefix :-

Total no. of C	Prefix form
1C	form
2C	acet
3C	propion
4C	butyr — n, iso
5C	valer

3C + 1db	C–C=C	acryl
4C + 1db	C–C=C–C	croton

Suffix :-

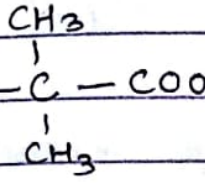
P. fg	Suffix
–COOH	ic acid
$-\overset{O}{\overset{\|\|}{C}}-O-R$	alkyl --- ate
$-\overset{O}{\overset{\|\|}{C}}-X$	yl halide
–CHO	aldehyde
$-\overset{O}{\overset{\|\|}{C}}-O-\overset{O}{\overset{\|\|}{C}}-$	ic anhydride
–CN	o nitrile
–NC	o iso nitrile
$-\overset{O}{\overset{\|\|}{C}}-NH_2$	amide

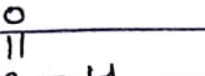
$H-\overset{O}{\overset{\|\|}{C}}-O-H$ formic acid

CH_3-COOH acetic acid

$CH_3-\underset{CH_3}{\underset{|}{CH}}-COOH$ isobutyric acid

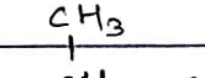
$CH_3-CH_2-CH_2-COOH$ — n-butyric acid

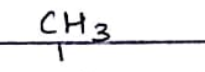

neo valeric acid


formaldehyde

$CH_3-CH=O$ — acetaldehyde

$CH_3-CH_2-CH=O$ — propionaldehyde


isobutyraldehyde

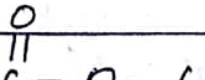

isovaler aldehyde

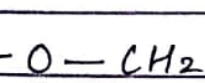
$CH_2=CH-CH=O$ — acryl aldehyde

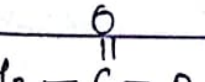
$CH_2=CH-COOH$ — acrylic acid

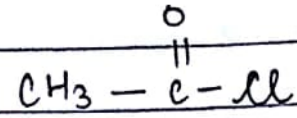
$CH_3-CH=CH-CHO$ — Croton aldehyde

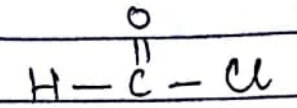
$CH_3-CH=CH-COOH$ — crotonic acid


methyl acetate

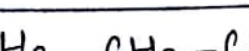

ethyl formate

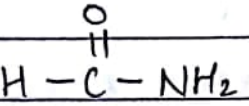

isopropyl ~~ethe~~ acetate

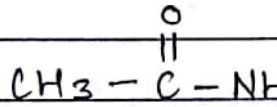

acetyl chloride

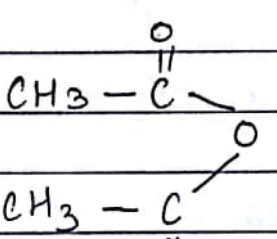

formyl chloride

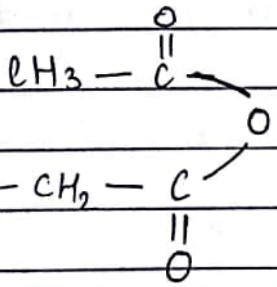
propionitrile


aceto isonitrile


formamide


acetamide


acetic anhydride


acetic propionic anhydride

DERIVED NAME :

(3) Based on famous Homolog

selection of parent chain
4°C > 3°C > 2°C > 1°C

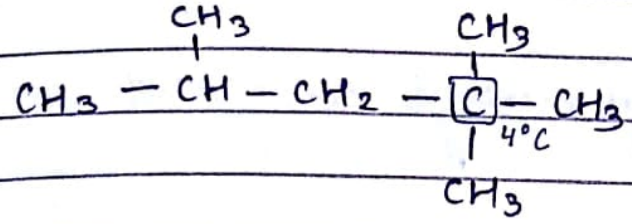
alkane :- CH_4 methane

$CH_3-\boxed{CH_3}$ methyl methane

$CH_3-\underset{CH_3}{CH_2}-\boxed{CH_2}-CH_3$ ethyl methyl methane

$CH_3-\underset{CH_3}{\boxed{CH}}-CH_2-CH_3$ ethyl dimethyl methane

Ques Write derived name of iso-octane?



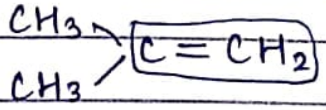
isobutyl trimethyl methane

alkene : $CH_2=CH_2$ ethylene

Fg > multiple bond
=> =

$CH_2=CH-CH_3$ methyl ethylene

$CH_3-CH=CH-CH_3$ sym. dimethyl ethylene



unsym. dimethyl ethylene

alkyne : $CH\equiv CH$ acetylene

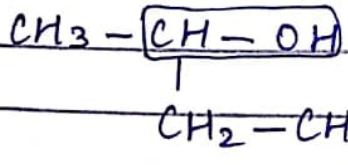
$CH_3-C\equiv CH$ methyl acetylene

$CH_3-C\equiv C-CH_2-CH_3$ ethyl methyl acetylene

$CH_2=CH-C\equiv CH$ Vinyl acetylene

alcohol : CH_3-OH Carbinol

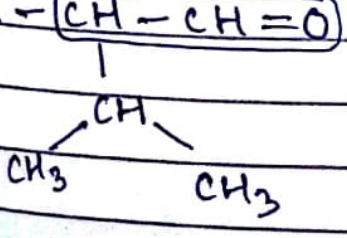
$CH_3-CH_2-CH_2-OH$ ethyl carbinol



ethyl methyl carbinol

aldehyde : $CH_3-CH=O$ acetaldehyde

$CH_3-CH_2-CH=O$ methyl acetaldehyde



isopropyl methyl acetaldehyde

acid : CH_3-COOH acetic acid

$CH_3-\underset{\displaystyle CH_2-CH_3}{\underset{|}{CH}}-COOH$ ethyl methyl acetic acid

$Cl_3C-COOH$ trichloro acetic acid

Ketone : $CH_3-\overset{\displaystyle O}{\overset{||}{C}}-CH_3$ acetone

$CH_3-CH_2-\overset{\displaystyle O}{\overset{||}{C}}-CH_2-CH_2-CH_3$

ethyl methyl acetone

IUPAC Naming

International Union of Pure & Applied Chemistry

Parts of IUPAC

1. Word Root
2. Prefix — Pri / Sec.
3. Suffix — Pri. / Sec.

Secondary Prefix	Primary Prefix	Word Root	Primary Suffix	Secondary Suffix
2°	1°	WR	1°	2°
substituent is written in alphabetical order - Cl chloro - NO_2 Nitro	used when compd is cyclic "cyclo"	represent total no. of C in Principal C chain 1C - meth 2C - eth 3C - Prop ⋮ 10C - dec 11C - undec 12 - dodec ⋮ 20 Eicos	tells whether compd. is saturated or unsaturated C - C ane 1C = C ene 1C ≡ C yne 2C = C diene 2C ≡ C diyne C=C & C≡C enyne	tells about Principal func gp. - COOH - CHO - OH - NH_2

PRIORITY ORDER of FUNCTIONAL GROUP :-

Functional Gp.	Prefix	Suffix
-(C)OOH (carboxylic acid)	x	oic acid
-COOH	carboxy	carboxylic acid
$-SO_3H$ (Sulphonic acid)	sulpho	sulphonic acid
$-(C)(=O)-O-(C)(=O)-$ (anhydride)	x	oic anhydride

Functional Group	Prefix	Suffix
-(C)OOR (ester)	x	alkyl --- oate
-COOR	alkoxycarbonyl or carbalkoxy	alkyl --- carboxylate
-(C)OX (acidhalide)	x	oyl halide
-COX	halo formyl	carbonyl halide
-(C)ONH$_2$ (amide)	x	amide
-CONH$_2$	carbamoyl	carboxamide
-(C)N (cyanide)	x	Nitrile
-CN	cyano	carbonitrile
-N$\rightleftarrows$(C) (isocyanide)	x	isonitrile
-NC	isocyano/ carbyl amino	carbyl amine
-(C)HO (aldehyde)	oxo	al
-CHO	formyl	carbaldehyde
-(C)- (Ketone) $\parallel$ O	Keto/oxo	one
-OH (alcohol)	hydroxy	ol
-SH (thio alcohol)	mercapto	thiol
-NH$_2$ (amine)	amino	amine

Substituents	Prefix
-R	alkyl
-X	halo
-NH$_2$	amino
$-N^{=O}_{\rightarrow O}$	nitro
-O-N=O	nitrite
-N=O	nitroso

Substituents	Prefix
$-OCH_2CH_3$	ethoxy
$-CH_2-OH$	hydroxy methyl
$-CH_2-Cl$	Chloro methyl
$-NH-CH_3$	methyl amino
$-S-$	thio
$-S-R$	alkyl thio
$CH_3-\overset{\overset{O}{\|}}{C}-O-$	acetoxy/ethanoyloxy
$CH_3CH_2-\overset{\overset{O}{\|}}{C}-O-$	propanoyloxy
$C_6H_5-\overset{\overset{O}{\|}}{C}-O-$	benzoyloxy
$-OR$	alkoxy

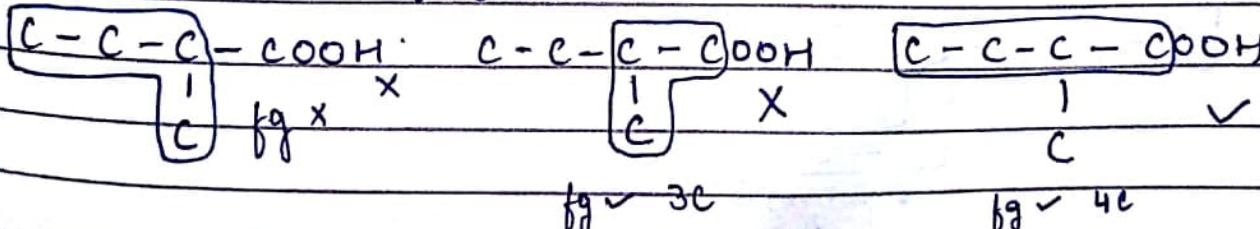
Rules of IUPAC :-

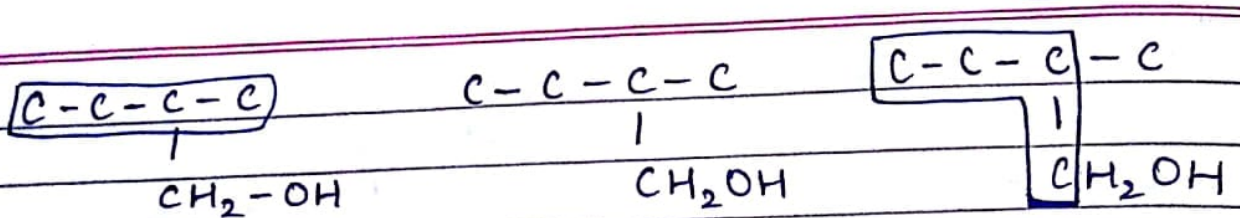
1. Selection of Principal Carbon chain
2. Numbering of P.C.C.

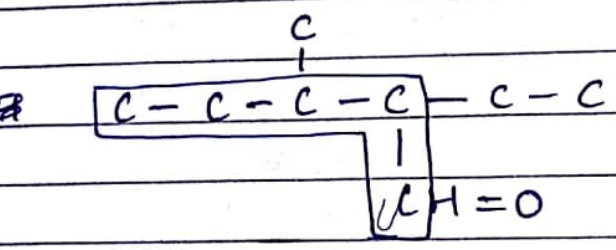
1. Selection of Principal Carbon Chain (PCC)

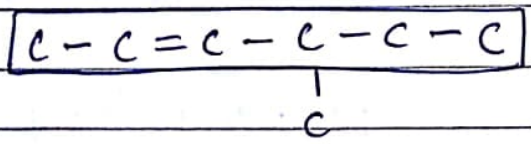
P.C.C. will be that ~~largest~~ longest continuous chain of carbon which have

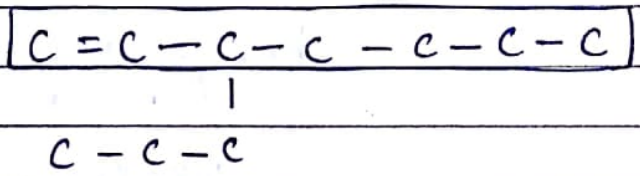
Pfg > Max^m no. of multiple bond > max. no. of C-atom

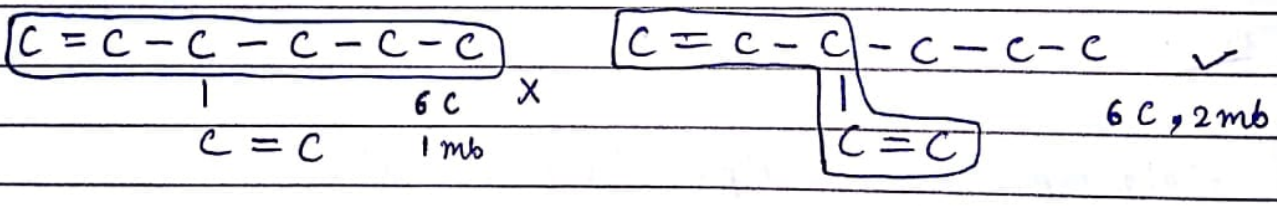


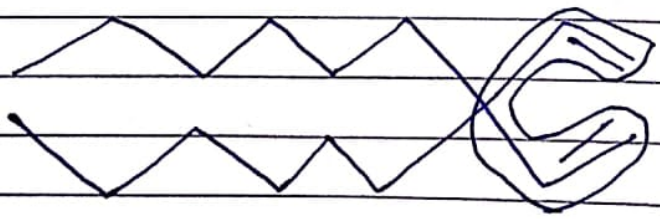




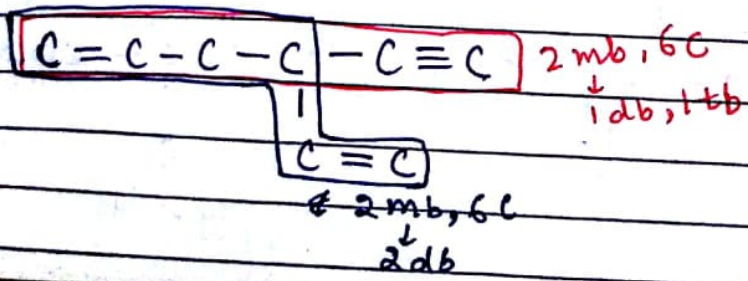


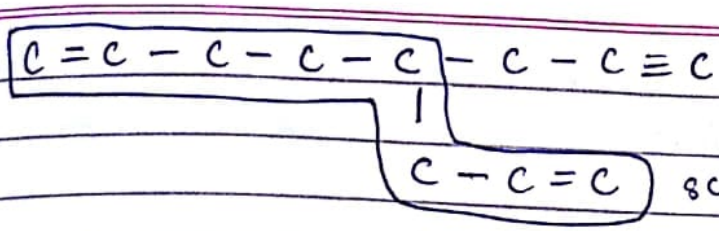




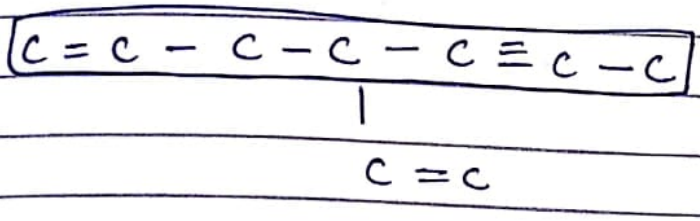


If symmetric molecule is not then priority is given to double & bond over triple bond

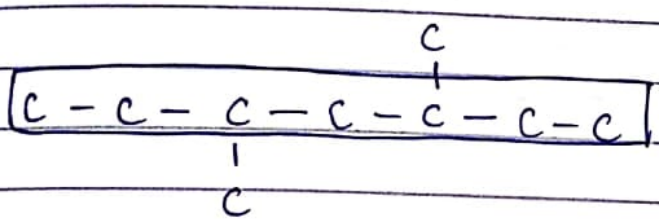




8C, 2 mb → 2 db

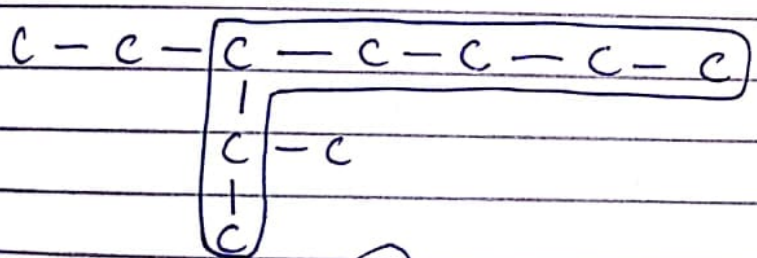


7C, 2 mb

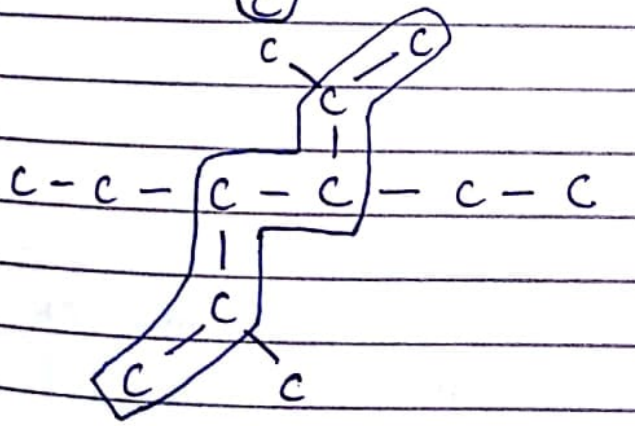


7C

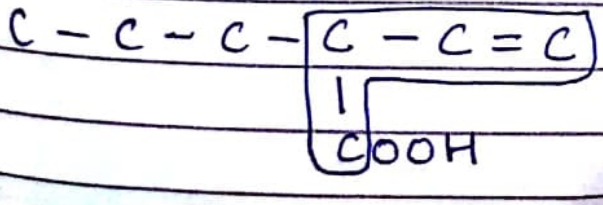
If 2 diff - P.C.C. have same no. of C in a molecule ~~tho~~ then that P.C.C. will be selected as P.C.C. which have max^m no. of sb (/ more branching)



2 sub, 7C

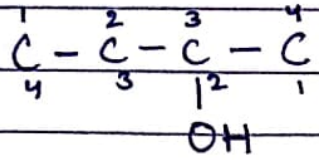


6C, 4 sub

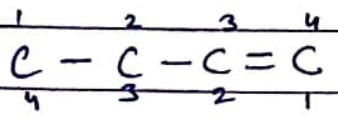


Numbering of P.C.C.:-

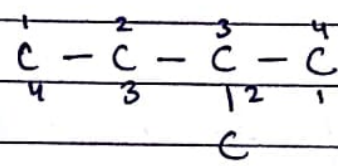
① Lowest no. will given to
Pfg > Mb > Sb



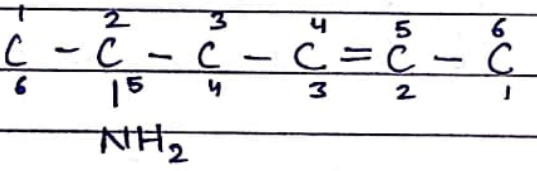
	fg
LHS	3
RHS	2



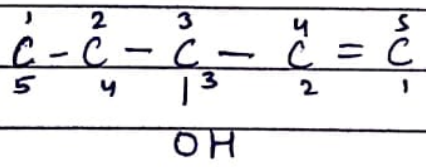
	mb
LHS	3
RHS	1



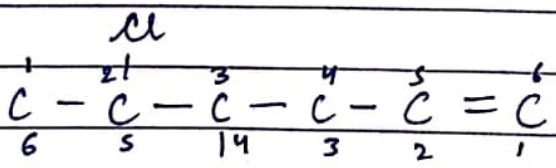
	Sb
L.H.S	3
RHS	2



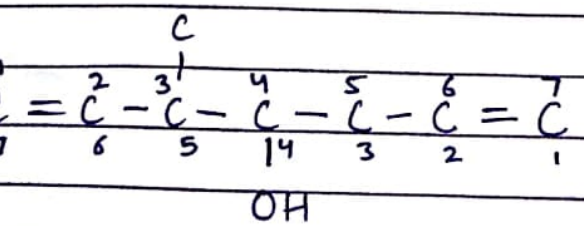
	fg
LHS	2
RHS	5



	fg	mb
LHS	3	4
RHS	3	1

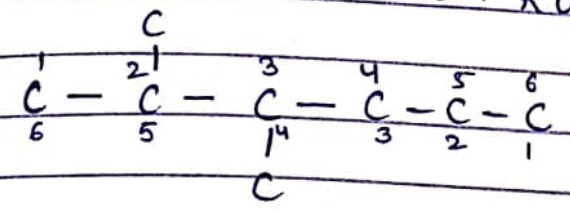


	fg
LHS	3
RHS	4

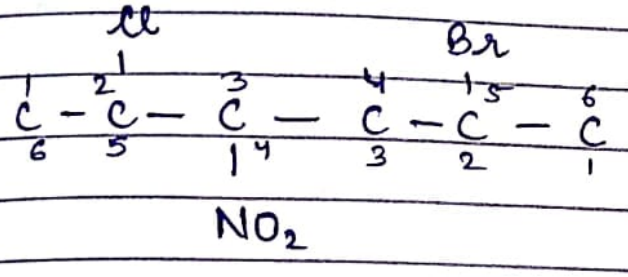


	fg	mb	sb
LHS	4	1, 6	3
RHS	4	1, 6	5

(2) If more than one substituent are +nt then least ~~low~~ locant no. Rule is followed.

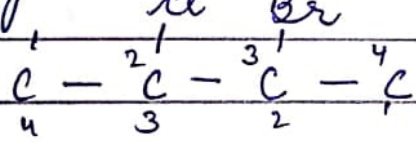


Sb
LHS 2,3 ✓
RHS 4,5

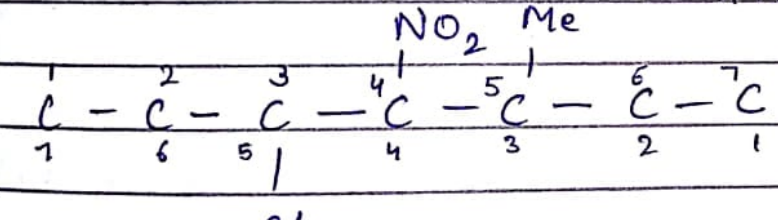


Sb
LHS 2, 3, 5 ✓
RHS 2, 4, 5

3. If 2 or more substituent are +nt at same posⁿ from either side then no. ~~of~~ starts from the side where alphabetically prefered substituent get lowest no.

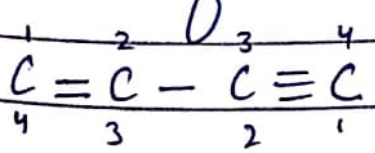


LHS 2, 3
RHS or 2, 3 ✓



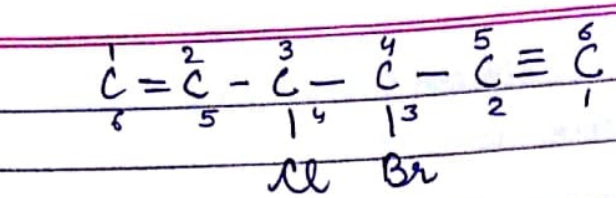
LHS 3 4 5 ✓
RHS 3 4 5

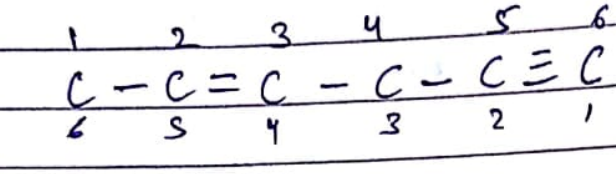
4. If db & tb are at similiar posⁿ from either side of PCC then priority is given to db ⁿ



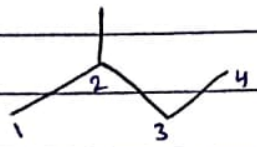
db tb
L·HS 1, 3 ✓
RHS 1, 3

PAGE NO.: 37
DATE: / /

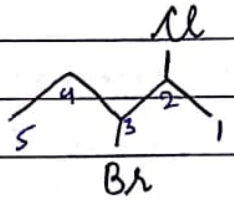




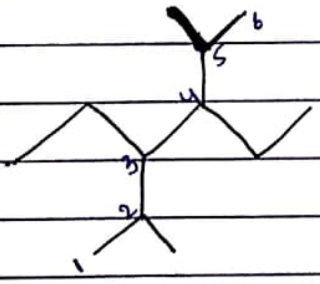
Ques



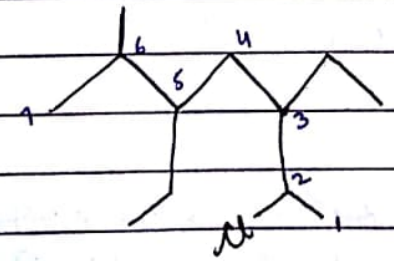
2-Methyl Butane



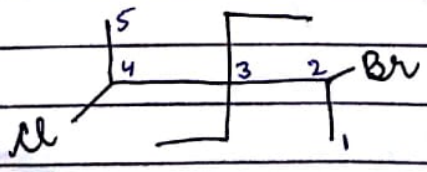
3-Bromo-2 Chloro Pentane



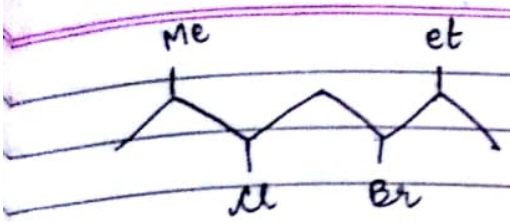
3,4 diethyl-2,5 dimethyl Hexane

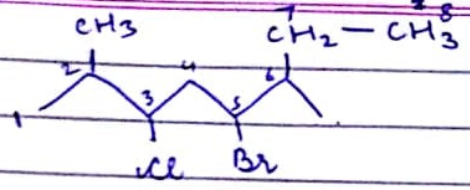


2-Chloro-3,5-diethyl-6-methyl heptane

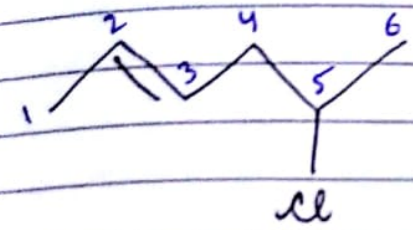


2-Bromo-4-Chloro-3,3-diethyl Pentane

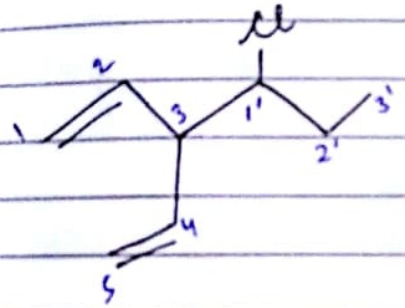




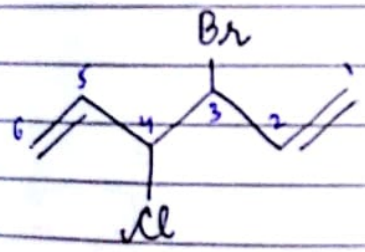
5-Bromo-3-Chloro, 2,6-dimethyl Octane



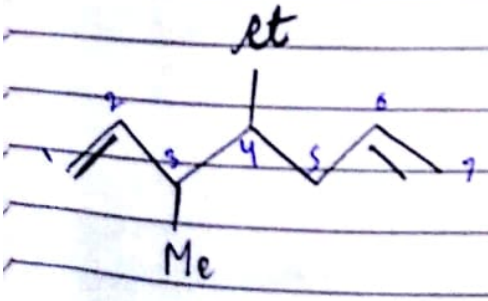
5-Chloro-~~Hex~~ Hex-2-ene



3(1'-Chloro Propyl) ~~Pent~~ Pent-1,4-diene



3-Bromo-4-Chloro hex-1,5 diene



4-ethyl-3-methyl-hept-1,6-di ene

$\overset{1}{CH_2}=\overset{2}{CH}-\overset{3}{CH_2}-\overset{4}{C}\equiv\overset{5}{CH}$ Pent-1-ene-4-yne

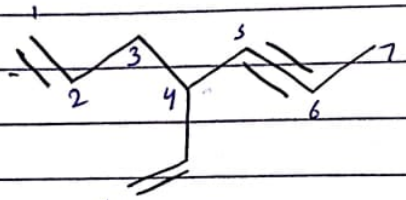
AIPMT-10

$CH_3 - \overset{6}{CH} = \overset{5}{CH} - \overset{4}{CH_2} - \overset{3}{C} \equiv \overset{2}{CH}$

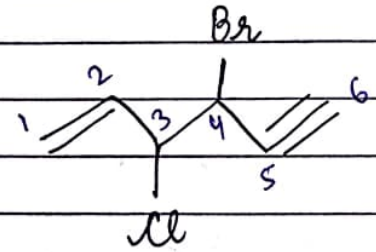
Hex - 4 - ene - 1 - yne

$CH_3 - \overset{3}{C}(=CH_2) - \overset{2}{C}(=CH_2) - CH_2 - CH_3$

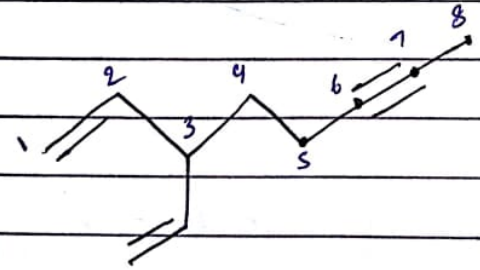
2 - Ethyl - 3 - Methyl But~~a~~ - 1, 3 - diene



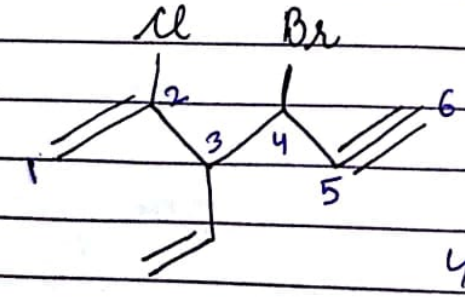
4 - ~~Eth~~ Vinyl hept - 1 ene - 5 - yne
or
4 - ethenyl hept - 1 - en - 5 - yne



4 - Bromo - 3 - Chloro - Hex - 1 - ene 5 - yne



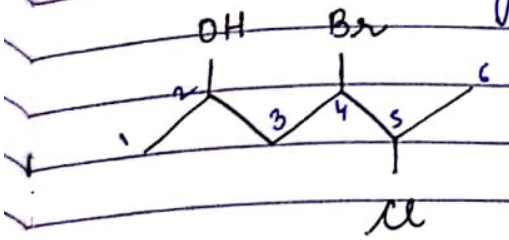
3 ethenyl - Oct - 1 ene - 6 - yne



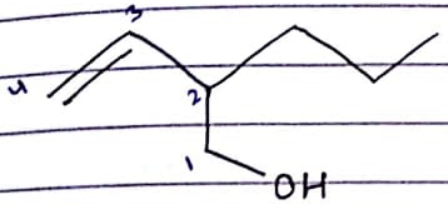
4 - Bromo - 2 - Chloro - 3 Ethenyl - Hex - 1 ene - 5 yne

4 - Bromo - 2 - Chloro - 3 - Vinyl Hex - 1 ene - 5 - yne

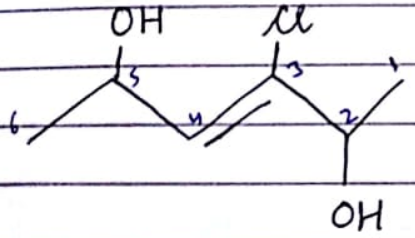
	Prefix	Suffix
-OH	Hydroxy	ol



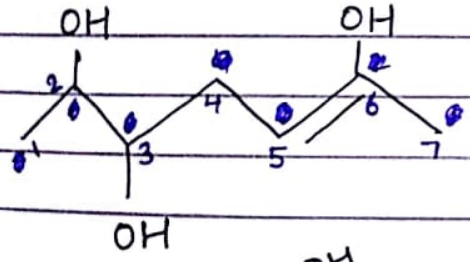
4 - Bromo - 5 - Chloro - 2 - Hexanol



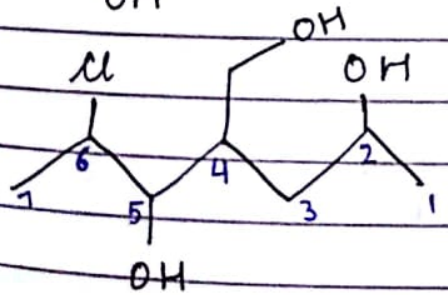
2 - Propyl - But - 3 - ene - 1 - ol



3 - Chloro - Hex - 3 - ene - 2, 5 - diol

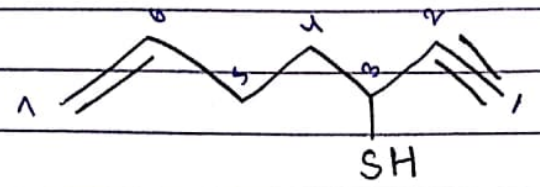


hept - 5 - ene - 2, 3, 6 - triol



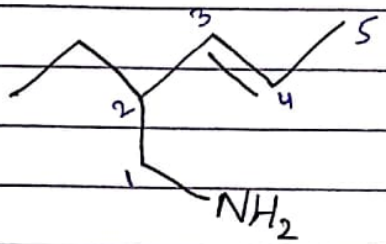
6 Chloro - 4 hydroxy Methyl hept - 2, 5 - diol

	Prefix	Suffix
$-SH$	mercapto	thiol

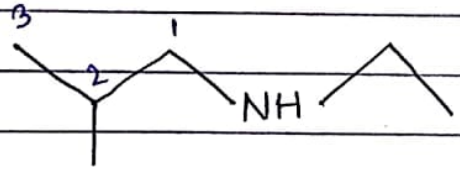


Hept - 6 - ene - 1 - yne - 3 - thiol

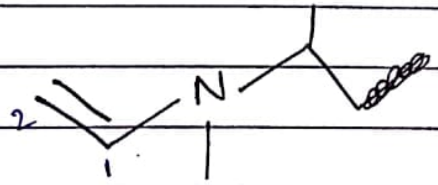
	Prefix	suffix
$-NH_2$	amino	amine



2 - Ethyl - Pent - 3 - ene - 1 - amine

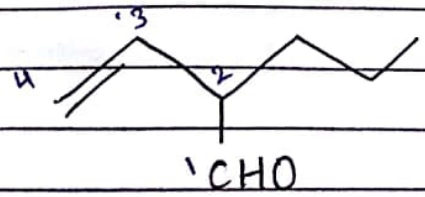


N - ethyl - 2 - methyl Propan amine

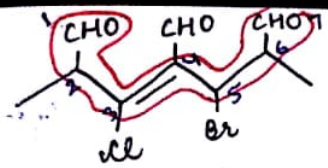


N - isopropyl - N - methyl ethene amine

	Prefix	Suffix
$-CHO$	oxo / formyl	al / carbaldehyde

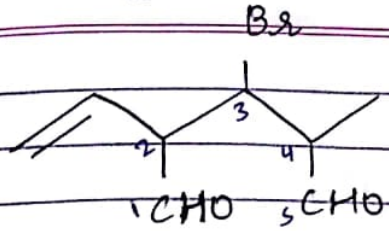


2 - Propyl But - 3 - en al



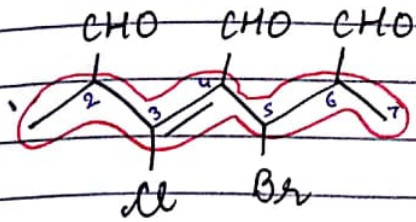
5-Bromo - 3 Chloro . 4 - formyl - 2,6 - dimethyl hept an - 1, 7 - dial

PAGE NO.: 43
DATE: / /

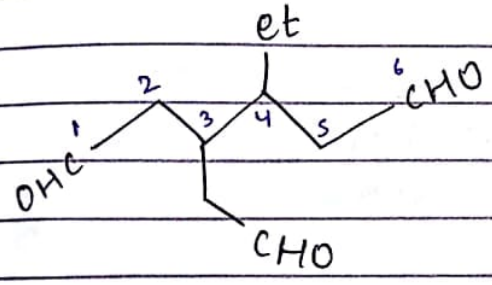


3 - Bromo - 2 - ethenyl - 4 - methyl Pentan - 1, 5 - dial

Note : When more than 2 carbon containing same func gp are attached to main ring chain then exclude all of them.

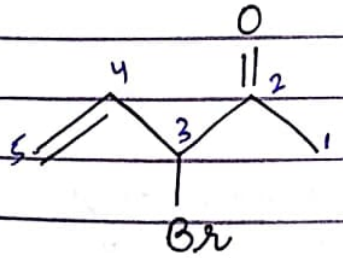


5 - Bromo - 3 - Chloro hept - 3 - ene - 2,4,6 - tricarbaldehyde

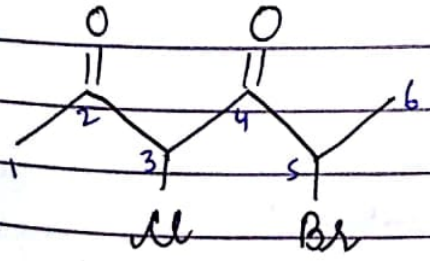


4 - ethyl - 3 - formyl methyl - hexan - 1, 6 - dial

	Prefix	Suffix
$-\overset{\overset{O}{\|\|}}{C}-$	oxo / keto	one



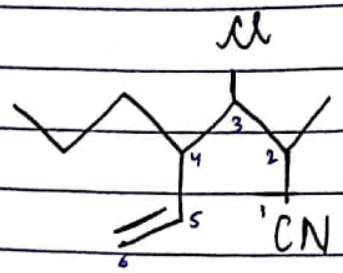
3 - Bromo - Pent - 4 - en - 2 - one



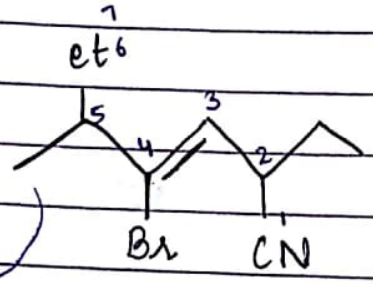
5 Bromo - 3 - Chloro hexan - 2,4 - dione

PAGE NO.: 44
DATE: / /

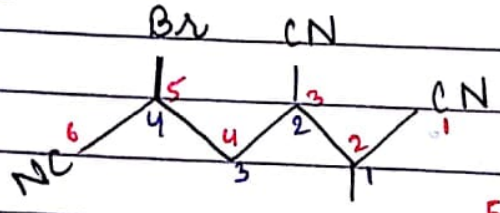
$-CN$ $-(C)N$ $-CN$ x cyano nitrile carbonitrile



3-Chloro-2-Methyl-4-Propyl-hex-5-ene nitrile



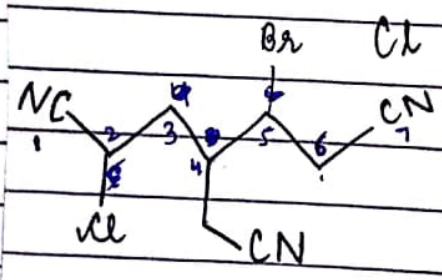
4-Bromo-2-ethyl-5-Methyl hept-3-ene nitrile



Pfg
CHS 1,3,4
RHS 1,2,4

4-Bromo-1-Chloro butane-1,2,4-carbo nitrile

5-Bromo-2-Chloro-3-cyano hexane-1,6-dinitrile



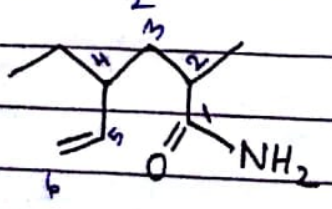
5-bromo-2-Chloro-4-cyano methyl heptane-1,7-dinitrile

$-\overset{O}{\overset{\|}{C}}-NH_2$

$-(\overset{O}{\overset{\|}{C}})-NH_2$

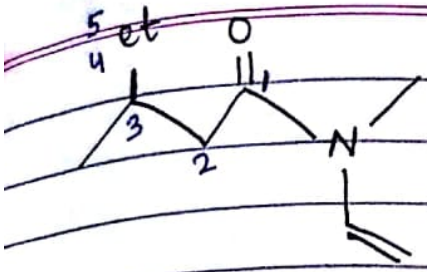
$-\overset{O}{\overset{\|}{C}}-NH_2$

x carbamoyl amide carboxamide

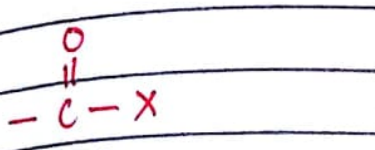


4-Ethyl-2-Methyl-hex-5-ene amide

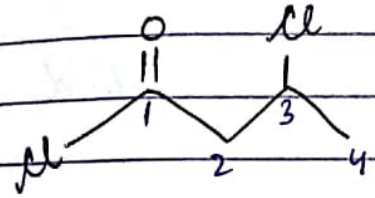
PAGE NO.: 45
DATE: / /



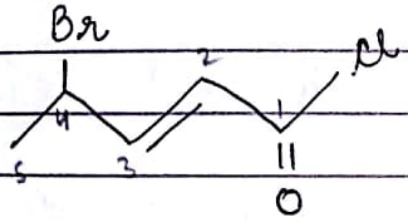
N - ethenyl - 3, N - dimethyl Pent -anamide



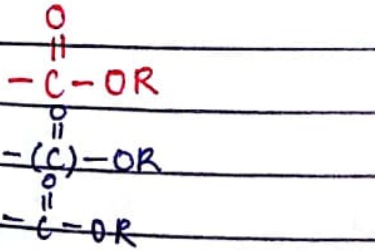
	X	oyl halide
	halo formyl	carbonyl halide



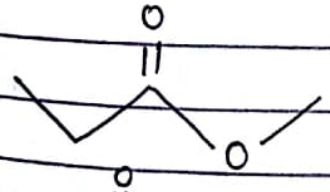
3 - Chloro - Butanoyl Chloride



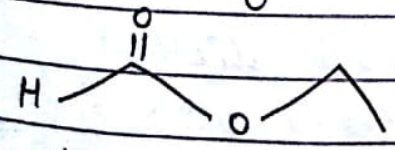
4 - Bromo - Pent - 2 - eneoyl Chloride



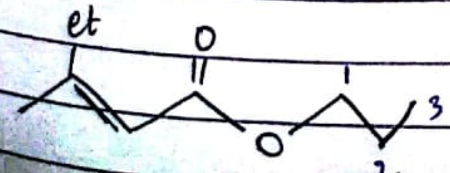
	X	alkyl --- oate
	alkoxy carbonyl	alkyl --- carboxylate



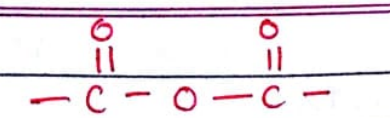
methyl - Propanoate

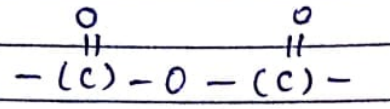


Ethyl - Methanoate

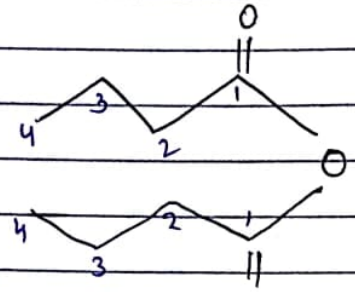


Propyl
~~Ethyl~~ - 3 - Methyl - Pent 2 - ene oate

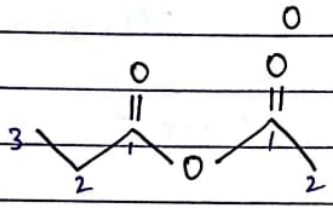




x oic anhydride

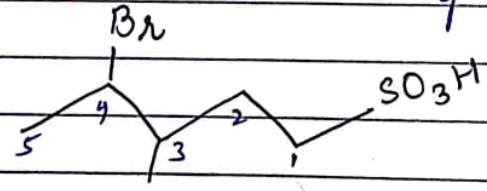


Butanoic anhydride



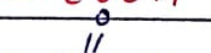
Ethanoic - Propanoic anhydride

$-SO_3H$ Sulpho sulphonic acid

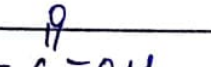


4-Bromo-3-Methyl Pentane sulphonic acid.

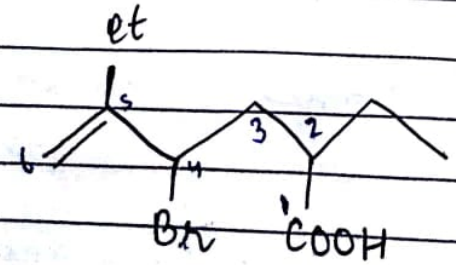
$-COOH$



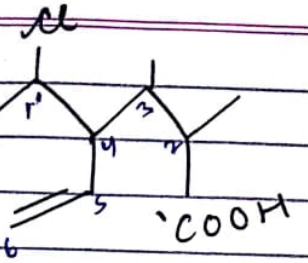
x oic acid



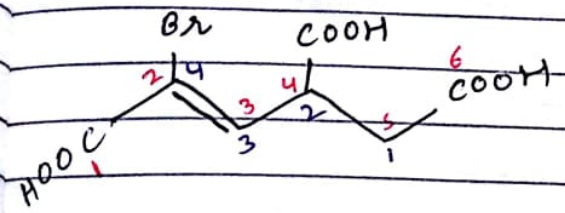
Carboxy carboxylic acid



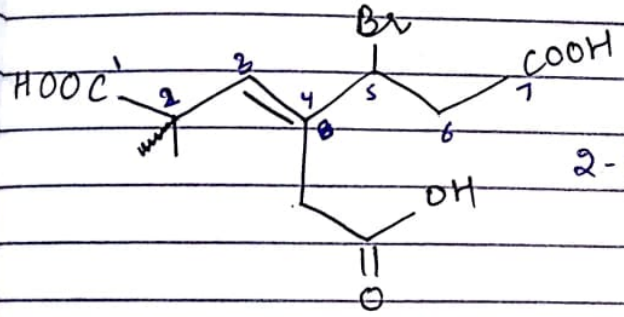
4-Bromo-2,5-diethyl hex-5-ene-oic acid



4 (1'-Chloro Ethyl) - 2, 3 - dimethyl hex - 5-ene - oic acid



~~2-bromo - 4-carboxy hex - 2 - ene - 1, 6 - dioic acid~~ But-3-ene -
4 - Bromo - 1, 2, 4 - ~~dicarboxy~~ dicarboxylic acid ~~But-3-ene~~

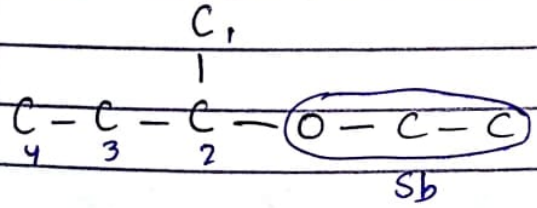


5 - Bromo - 4 - carboxy ~~ethyl~~ Methyl
2 - Methyl-Hept - 3 - ene - 1, 7 - dioic acid

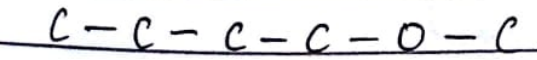
IUPAC Naming

$CH_3 - O - CH_3$ — methoxy methane

$CH_3 - CH_2 - O - CH_3$ — methoxy ethane



2 - ethoxy butane

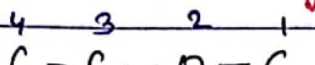


butoxy ethene

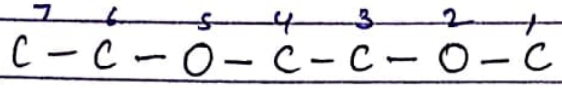
PAGE NO.: 48
DATE: / /

Oxa - System →

O को C atom से जोड़कर Least locant no. देनी है

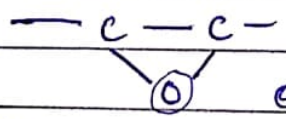


2-oxa-butane

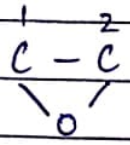


2,5-dioxa heptane

Cyclic ether

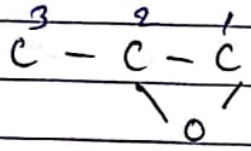


epoxy

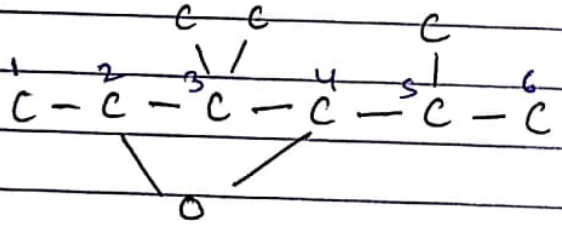


epoxy ethane

Or oxirane (trivial name)

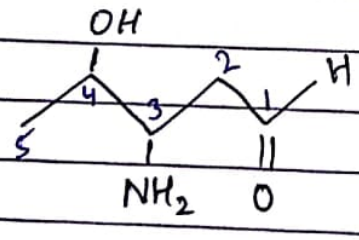


1,2-epoxy propane

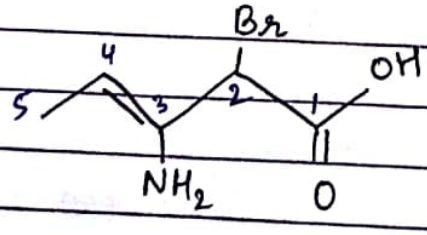


2,4-epoxy-3,3,5-trimethyl hexane

IUPAC Naming of Polyfunctional group :-

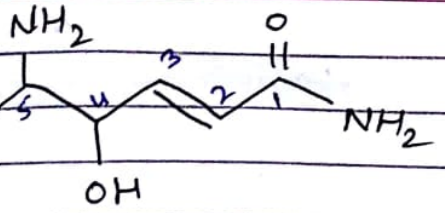


3-amino-4-hydroxy Pentanal

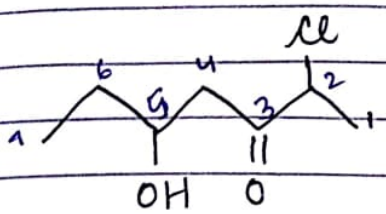


3-amino-2-Bromo-Pent-3-enoic acid

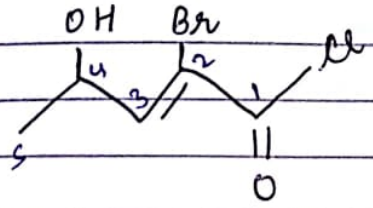
PAGE NO.: 79
DATE: / /



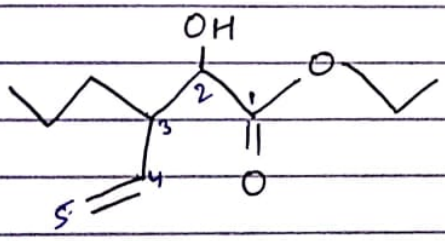
5 - amino - 4 - hydroxy hex - 2 - ene amide



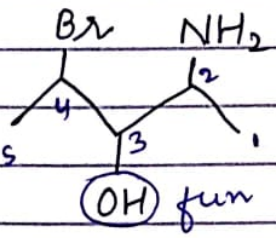
2 - Chloro - 5 - hydroxy heptan - 3 - one



2 - Bromo - 4 - hydroxy Pent - 2 - eneoyl Chloride

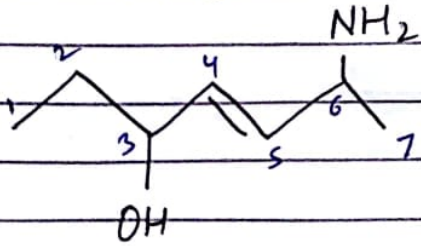


ethyl - 2 - hydroxy - 3 - Propyl Pent - 4 - eneoate

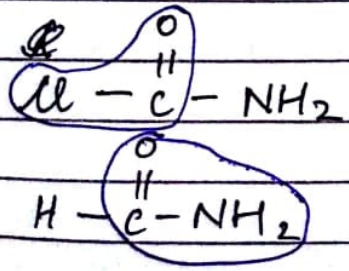


2 - amino - 4 - Bromo ~~- 3 - hydroxy~~ Pentan 3-ol ~~amine~~

(OH) fun

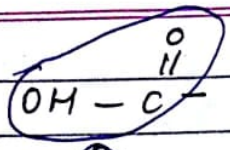


6 - amino hept - 4 - ene - 3 - ol

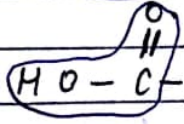


amino · methanoyl chloride

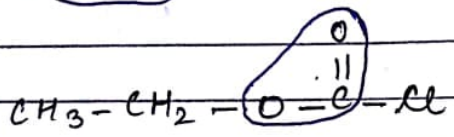
methanamide



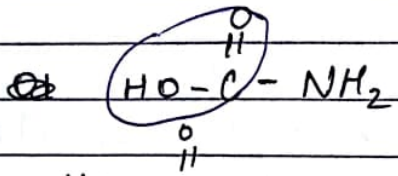
methoxy methanoic acid



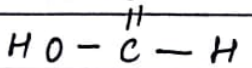
Chloro methanoic acid



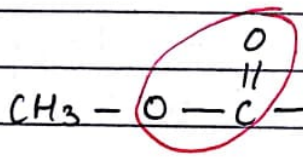
ethyl chloro methanoate



amino methanoic acid

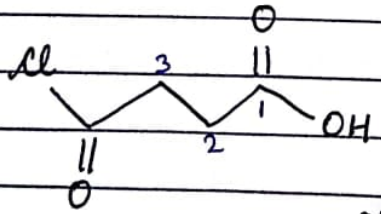


methanoic acid

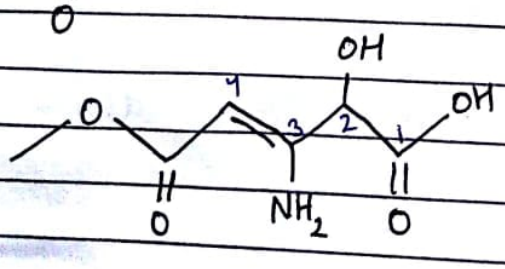


methyl amino methanoate

Note: In polyfunctional gp containing substance if C-containing func gp is +nt as substituents then its C is not considered in P.C.C. but in case of -CHO & $-\overset{\overset{O}{\|}}{C}-$ its C ~~is~~ can be included in chain.

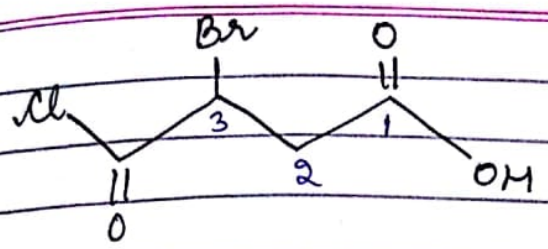


chloro formyl
3-~~methyl chloride~~ - Propanoic acid

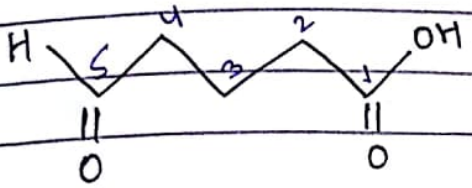


3-amino-2-hydroxy-4-methyl carbonyl-3-butenoic acid

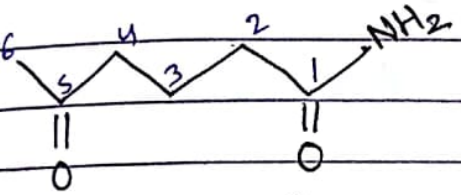
PAGE NO.: 51
DATE: / /



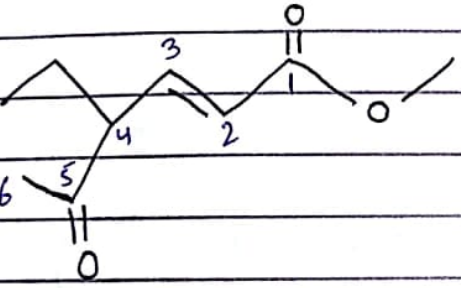
3 - Bromo - 3 - Chloroformyl Propanoic acid



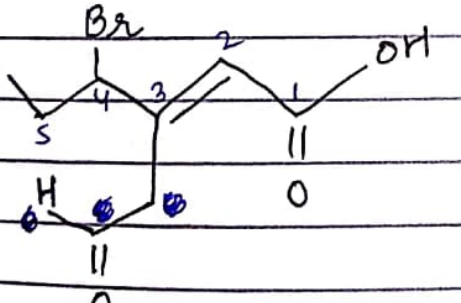
5 - Oxo - Pentanoic acid



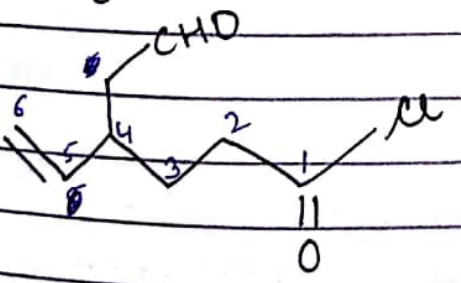
5 - Keto - Hexanamide



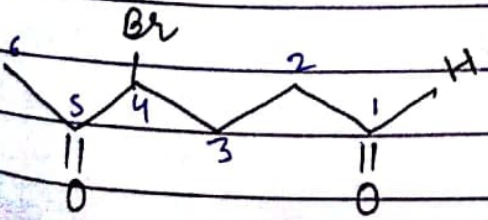
4 - ethyl - ~~methoxy carbonyl~~
5 - keto ~~Hexo - 2 - ene~~ methyl
Hex - 2 - ene oate



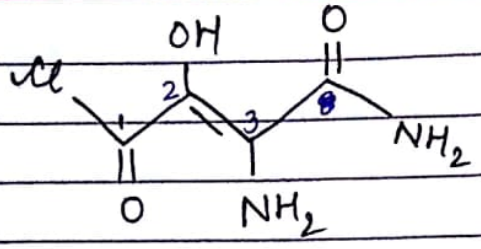
~~5 - Keto - 3 - Bromo Propyl~~
~~Hex - 2 - ene oic acid~~
4 - Bromo - 3 - methyl formyl hex - 2 - enoic acid



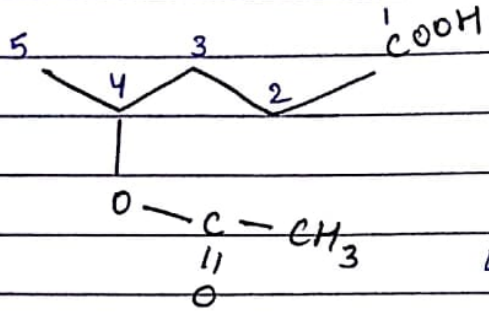
formyl methyl
4 - ~~Methyl formyl~~ - 5 - hexenoyl Chloride



4 - Bromo - 5 - keto Hexanal

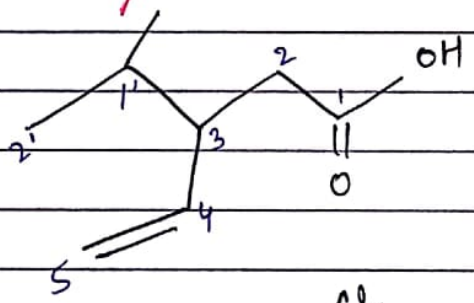


3 - amino - 3 - carbo amoyl
2 - hydroxy Prop
- 2 ene - ~~Chlor~~
oyl chloride

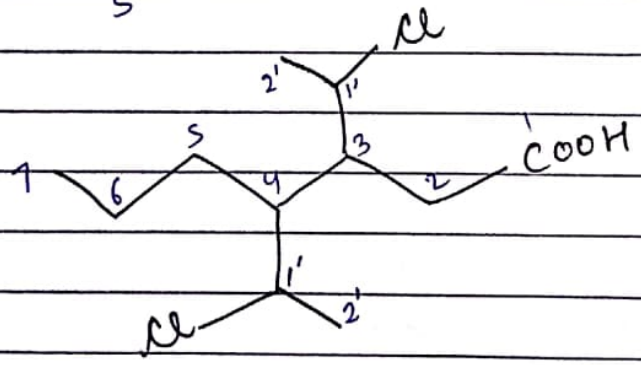


4 - ethanoyloxy pentanoic acid
or
4 - acetoxy Pentanoic acid

Complex Substitutent :-



3 isopropyl Pent - 4 - eneoic acid
or
3 (1'-methyl ethyl) pent - 4 - enoic acid

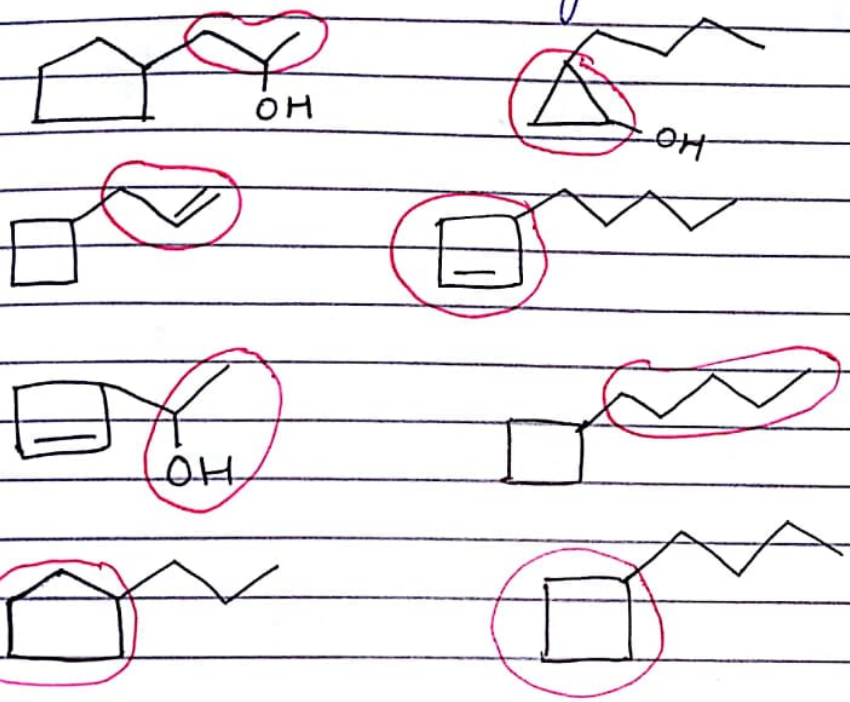


bis - 3,4 - (1'- Chloroethyl) heptanoic acid

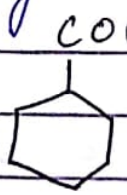
IUPAC name of Cyclic compounds :-

Rule ① If an org. comp. is consist of open chains & closed chain then that chain will be PCC which have -

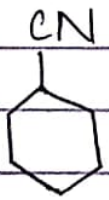
Pfg > max^m no. of m.b. > max^m no. of C-atom or longest chain > ring if open & closed chain have same no. of C-atom



Rule-2 If C-containing Pfg is directly attached to ring then it is consider as part of ring but its C is not included in P.C.C.



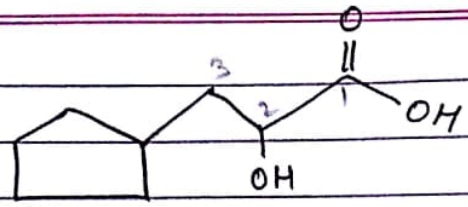
cyclo hexane carboxylic acid



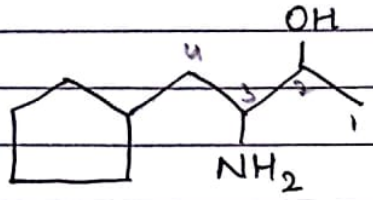
cyclo hexane carbonitrile

Numbering as usual

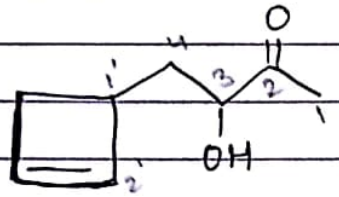
Note:- ① ring prefix → cyclo
② Word cyclo is considered in alphabetical order



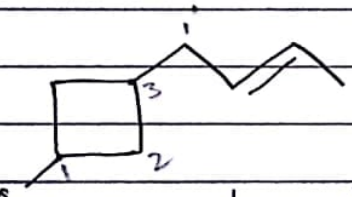
3-cyclo Pentyl-2-hydroxy Propanoic acid



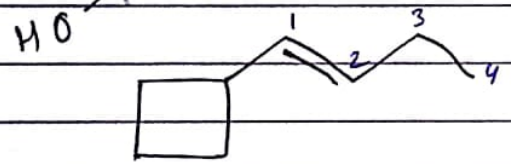
3-amino-4-cyclo Pentyl -Butan-2-ol



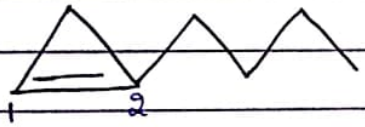
4 (2'-cyclobutenyl) -3-hydroxy -2-butanone



3 (2'-butenyl) butanol

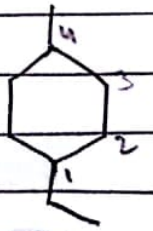


1-cyclo Butyl Buto-1-ene



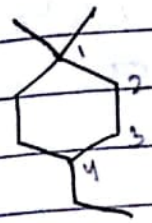
2-Butyl-cyclo Propene

1-Butyl-2-Chloro-cyclo Butane

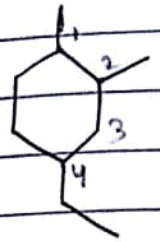


1-Ethyl-4-Methyl cyclo hexane

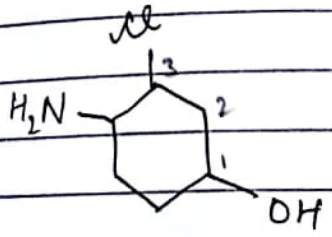
PAGE NO.: 55
DATE: / /



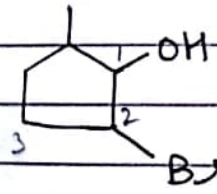
4-Ethyl-1,1-dimethyl cyclohexane.



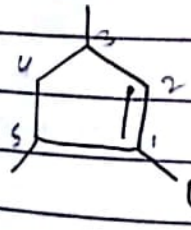
4-ethyl-1,2-dimethyl cyclohexane



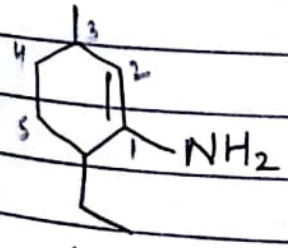
4-amino-3-chloro cyclohexanol



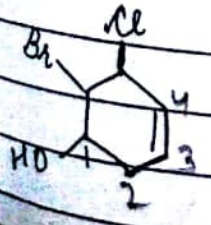
2-Bromo-5-methyl cyclo Pentanol



3,5-dimethyl cyclo Penteneol

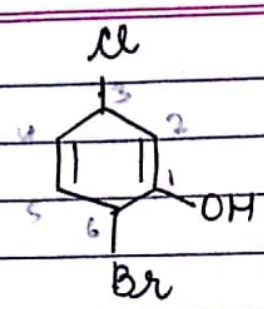


6-Ethyl-3-Methyl cyclo Hexeneamine

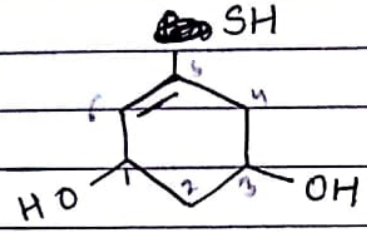


6-Bromo-5-Chloro-cyclohex-3-eneol

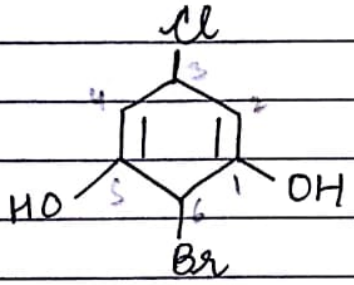
PAGE NO.: 56
DATE: / /

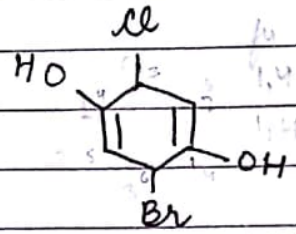


6 - Bromo - 3 - Chloro - cyclohex-1,4 - dienol



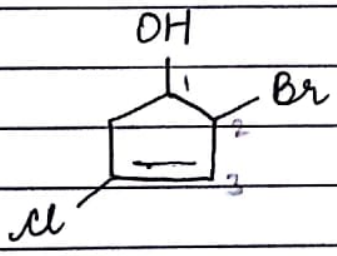
5- mercapto - cyclo - hex - 5 - ene - 1, 3 - diol



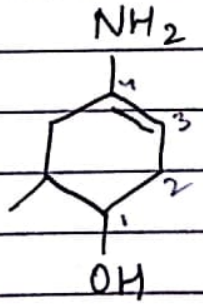


3- Bromo - 6 - Chloro cyclohex-1,4-diene - 1,4-diol

6-Bromo - 3- Chloro cyclohiex - 1,4-di-ene-1,5-diol

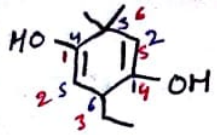


2- Bromo - 4 - Chloro - cyclo Pent-3-enol



4- amino - 6 - methyl - cyclo hex-3-enol

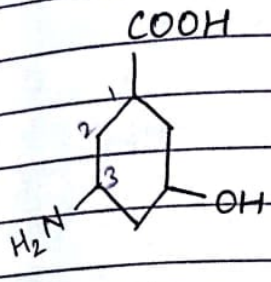
1 - carbonyl chloride - cyclohexane



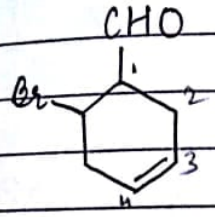
Pf	mb	sb
1,4	1,4	3,3,6 ✓
1,4	1,4	3,6,6

PAGE NO.: 57
DATE: / /

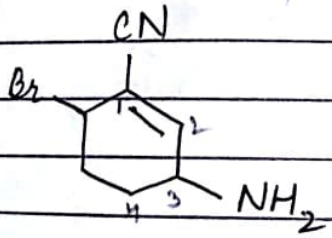
2- cyclo hexyl ethanal



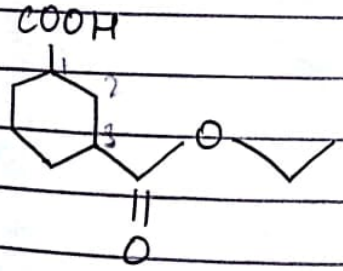
3 - amino - 5 - hydroxy - cyclohexane carboxylic acid



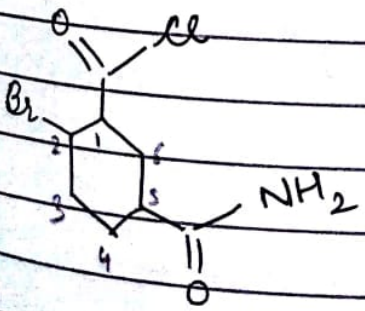
6 - Bromo - cyclohex - 3 - ene - carbaldehyde



3 - amino - 6 - Bromo - cyclohexene carbonitrile

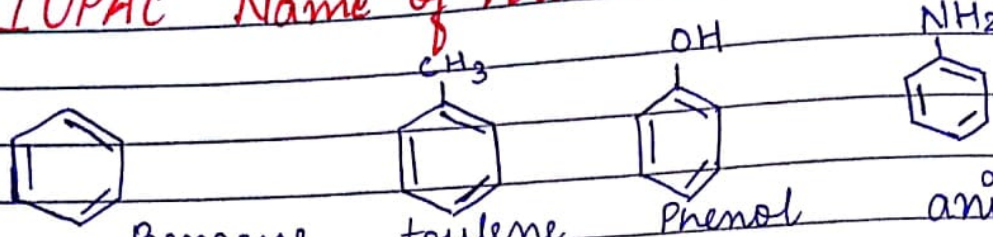


3 - ethoxy carbonyl - cyclo hexane carboxylic acid

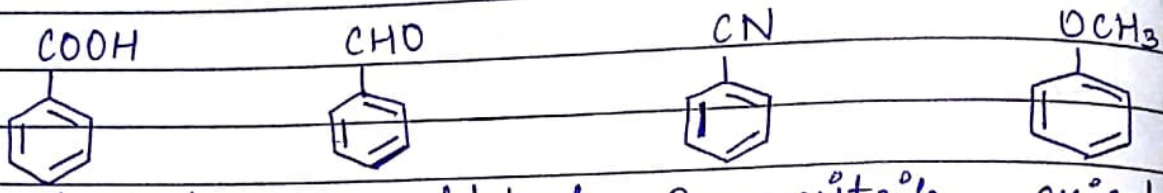


2 - Bromo - 5 - carbamoyl cyclo - hexane carbonyl halide

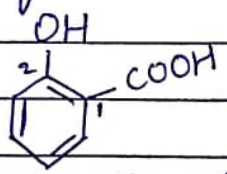
IUPAC Name of Aromatic



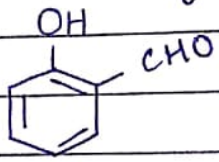
Benzene toulene Phenol aniline



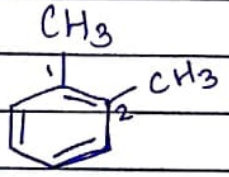
Benzoic acid Benzaldehyde Benzonitrile anisole



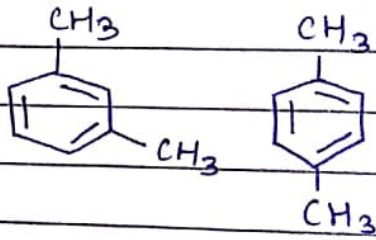
Salicyclic acid
2-hydroxy benzoic acid



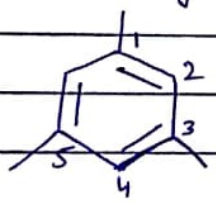
Salicyaldehyde
2-hydroxy benzaldehyde



1, 2-dimethyl benzene
O-Xylene



m-Xylene P-Xylene

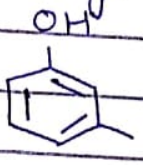


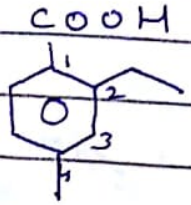
1, 3, 5-trimethyl benzene
or
mesitylene

Rule: ~~Its~~ If aromatic compd is consist of open and closed chain then PCC will be
① which have P func gp.
② Open chain if it have more than 2 carbon

(in case when fun^{c} gp not +nt)

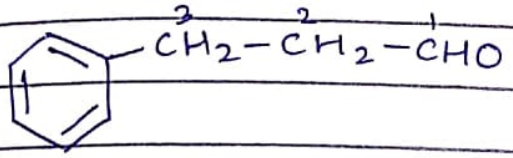
2-methyl Phenol
O-Cresol



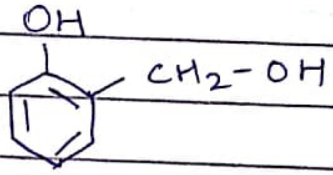


2-ethyl-4-methyl benzoic acid

2-phenyl ethanoic acid

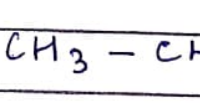


3-phenol propanal



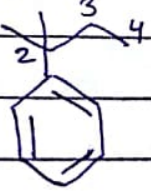
✓2-hydroxy methyl phenol
≡(2-Hydroxy phenyl)-methanol

phenolic gp
Phenol



alcoholic gp

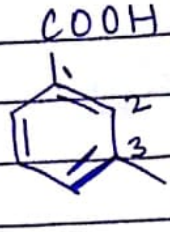
Alcoholic gp has more priority than phenolic gp.



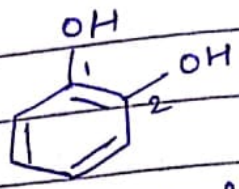
methyl benzene ethyl benzene 2-phenyl propane

1-Chloro-3 phenyl propane

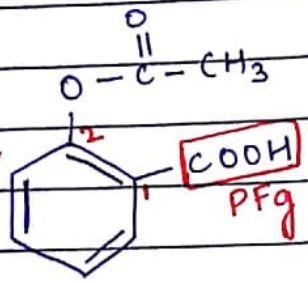
If more than 2 fun^{c} gp are +nt



benzene-1,3-dicarboxylic acid

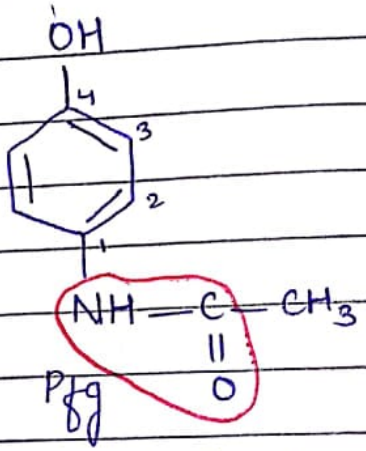


catechol
benzene-1,2-diol

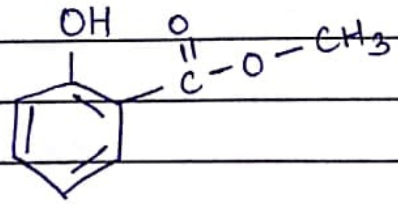


Aspirin

2-ethanoyl oxy benzoic acid
2-acetoxy benzoic acid
acetyl salicylic acid

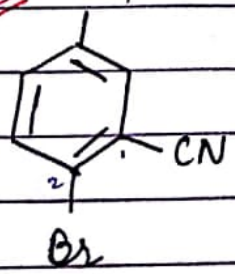


N-(4-hydroxy phenyl) ethanamide
Paracetamol



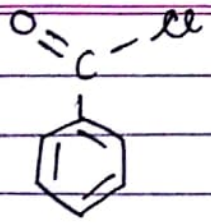
IUPAC
methyl-2-hydroxy Benzoate
common methyl salicylate
→ "oil of winter green"

IIT-09

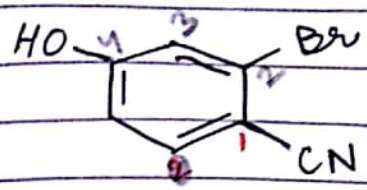


● 2-Bromo-5-hydroxy ~~phenyl cyano~~ benzo nitrile

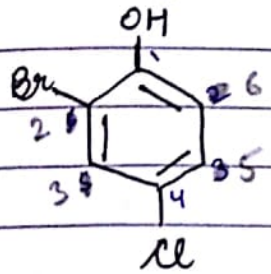
PAGE NO.: 6 /
DATE: / /



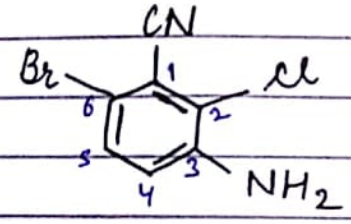
Benzene carbonyl chloride



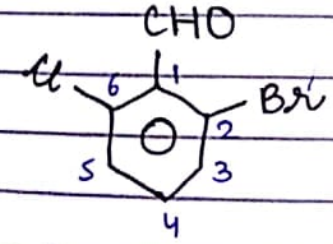
6 - Bromo - 4-hydroxy ~~benzene~~ benzo nitrile



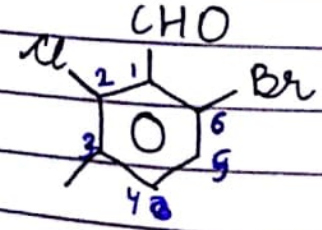
2 6 - Bromo - 4-Chloro - benzanol

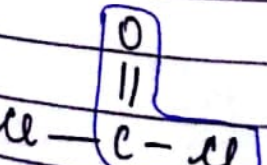


3 - amino - 6 - Bromo - 2 - Chloro ~~Cyclo~~ Benzo - nitrile.



2 - Bromo - 6 - Chloro - Benzaldehyde





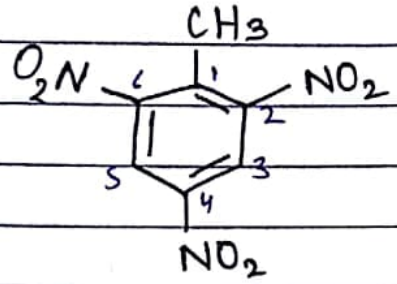
Phosgene gas
poisonous gas
Chloro - methanoyl chloride

PAGE NO.: 12
DATE: / /

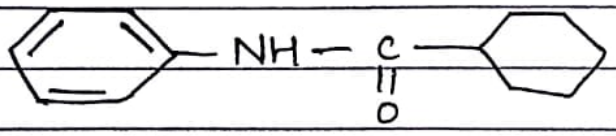
CCl_4 tetrachloro methane

$CHCl_3$ → chloroform / 1,1,1 - trichloro methane

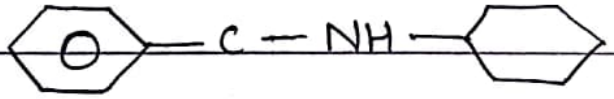
tear gas → $CCl_3.NO_2$
trichloro nitro methane



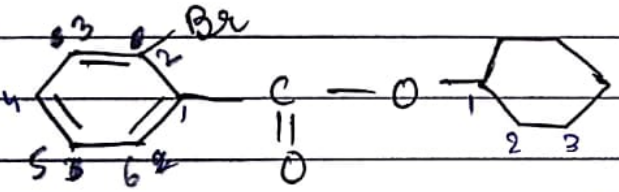
2,4,6 - trinitro toulene



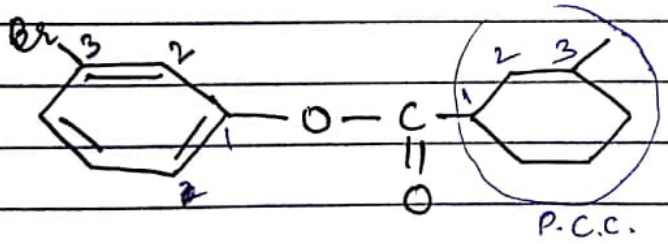
N - Phenyl ~~carbox~~ cyclo hexane carboxamide ~~hexane~~



N - cyclo - hexyl benzene carboxamide



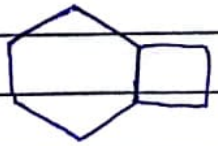
(4 - methyl cyclo hexyl) - 6 - Bromo Benzoate



(5 - bromo phenyl) - 3 methyl cyclo hexane carboxylate

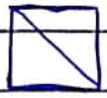
Bicyclo compounds

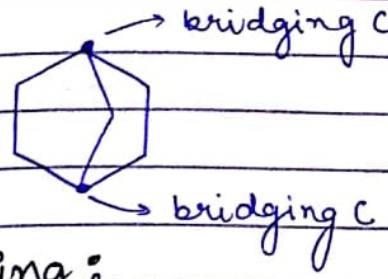
4 rings are fused



General formula → C_nH_{2n-2}

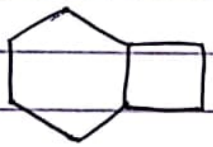
min C → 4C



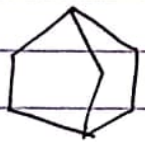


Naming:

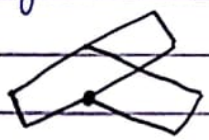
- Prefix → bicyclo
- no. of C-atoms in bridges are written in ↓ order in square bracket separated by full stop (.)
  [ ]
- word root → total no. of C-atom



bicyclo [4.2.0] Octane

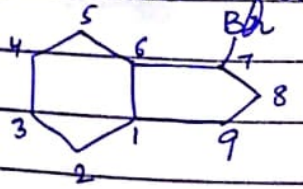


bicyclo [2.2.1] heptane

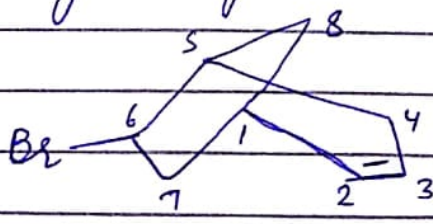


bicyclo [2.2.2] Octane

If subs. is +nt then numbering start from bridged carbon & then we move from large ring to small ring



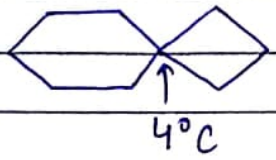
bicyclo [4.3.0] -7-Bromo nonane



bicyclo [3.2.1] 6-bromo-Oct-2-ene

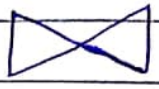
Spiro - compounds :-

When 2 rings are fused at common carbon (4°C)

eg 

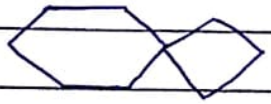
G.F. → C_nH_{2n-2}

min. no. of C → 5C

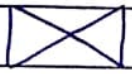


Naming :-

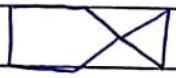
- Prefix spiro
- no. of C atoms in ring are written in ↑ order
- word root, suffix.



Spiro [3.5] nonane



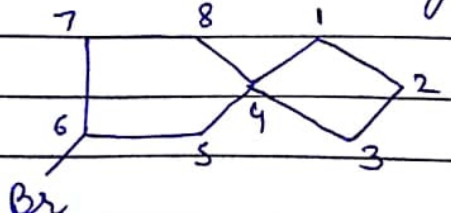
Spiro [2.2] Pentane



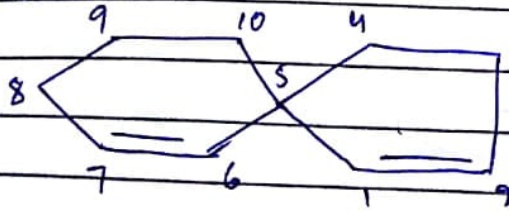
Spiro [2.4] heptane

* Sub is +nt

Numbering starts from adjacent carbon to common C & move through smaller ring to large ring

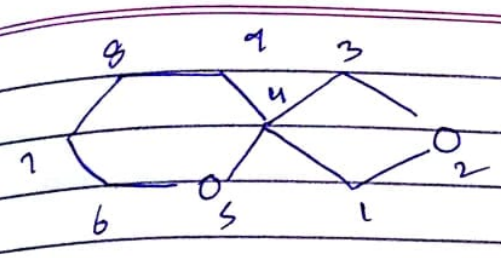


Spiro [3.4] - 6 - Bromo Octane



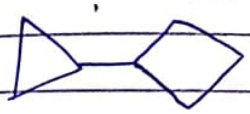
Spiro [4.5] dec - 1, 6 - diene

c)



spiro [3.5] 2,5 dioxa nonane

Extra different org i.e. no common carbon



1 - Chloro - 4 - Methyl benzene
4 - Chloro - toulene