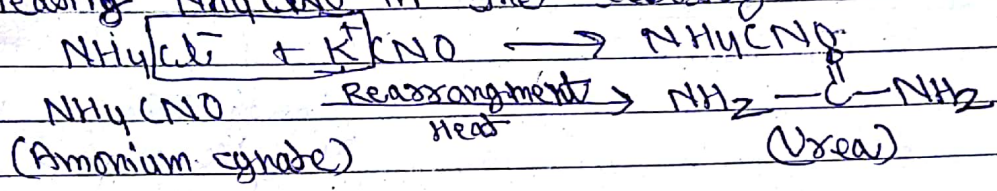


Note:- C has valency 4 and 1 covalent bond, an atom with covalency 4 is called

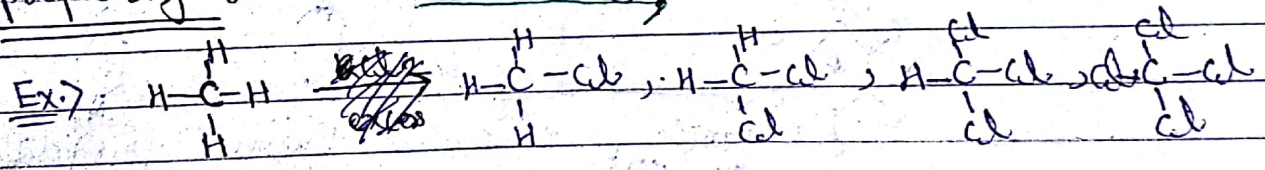
Vital Force theory :- (Brewster) :- ^{scientist} Organic comp's are those comp's which are ^{synthesise} synthesised inside living body and can't be synthesised outside the living body or laboratory.

Failure of vital force theory (sc. F. Wohler) :- F. Wohler gave VFT by preparing urea (an organic comp.) by heating NH_4CNO in the laboratory.

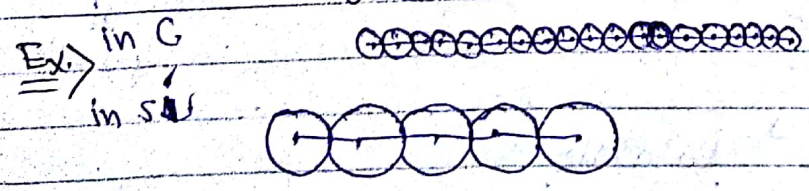


New concept of organic compounds :- Carbon and its derivatives are the organic comp. which may be synthesised inside the living body and outside the living body.

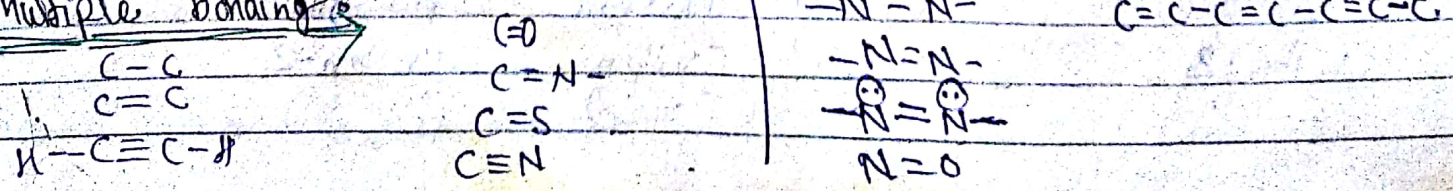
property :- (1) covalency

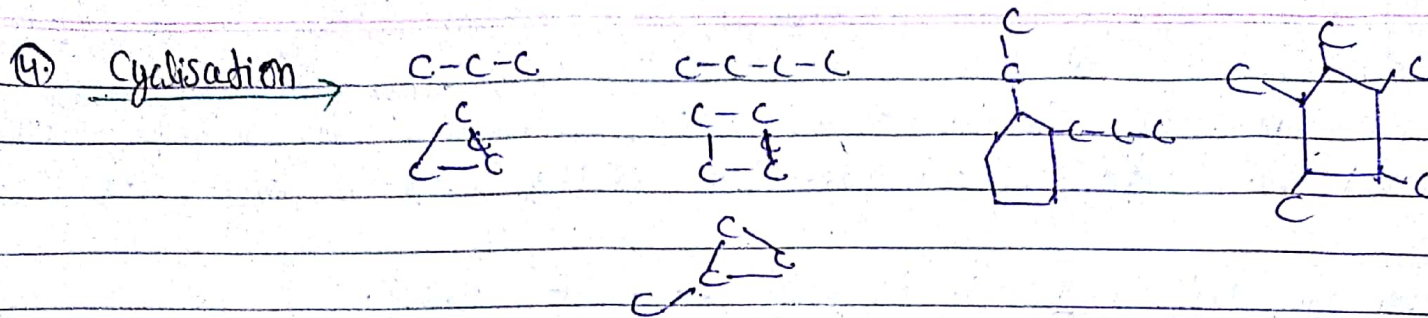


(2) Catenation :- self linking or self combining property of an atom.

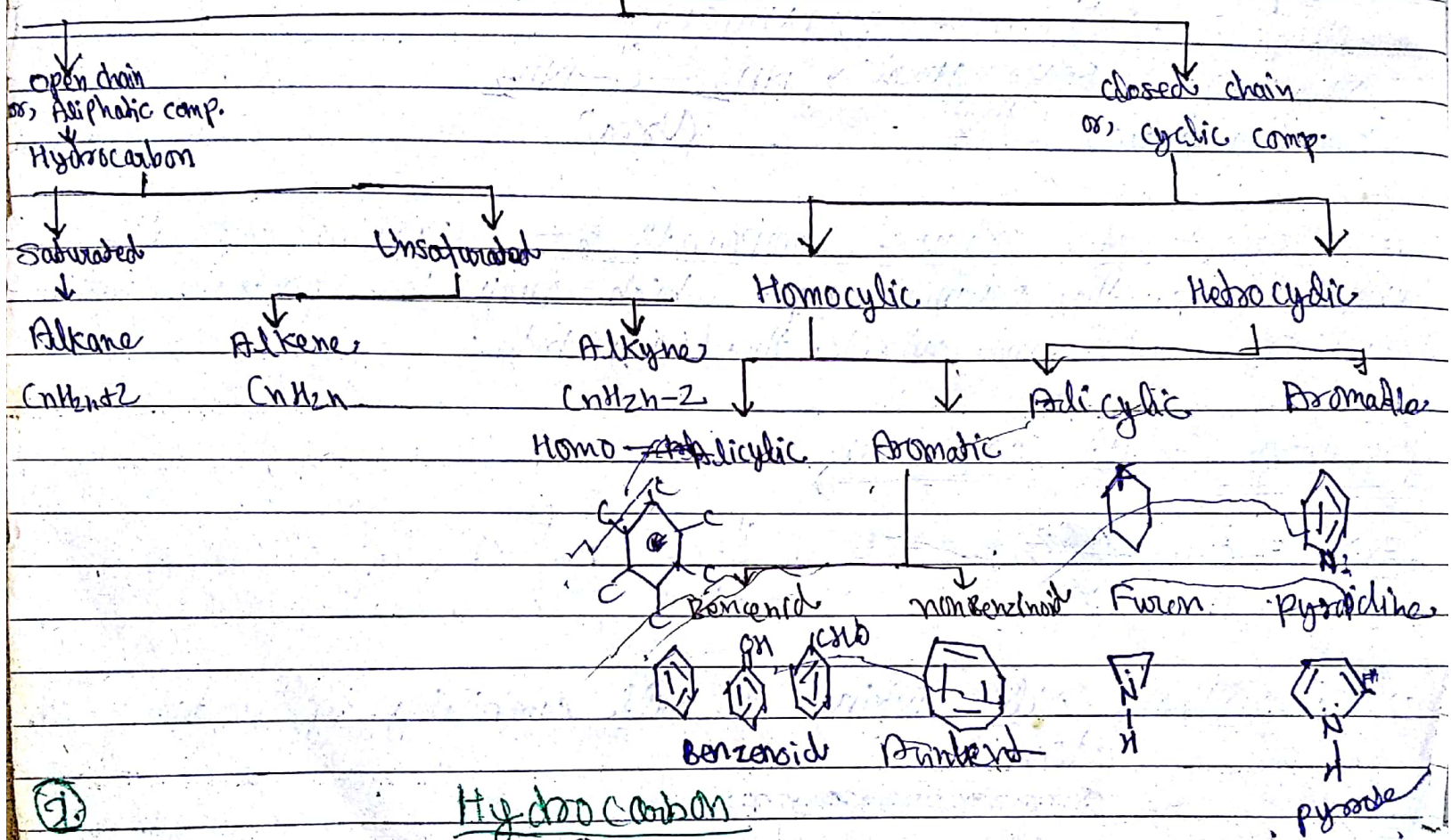


Multiple bonding

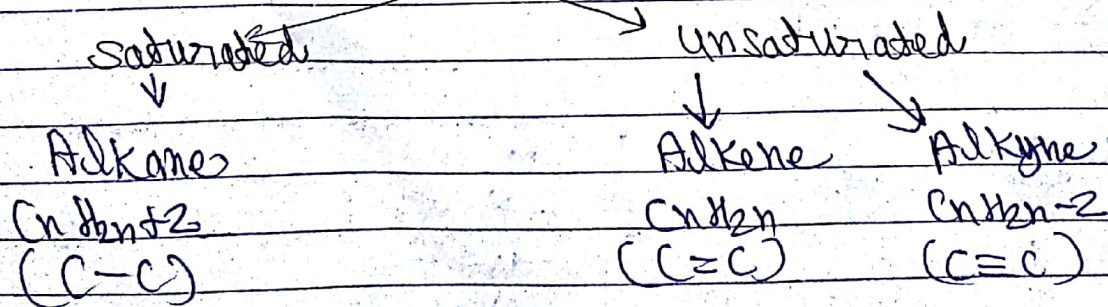




⑤ Organic compound



⑥ Hydrocarbon



All. = no. of carbon

n=1 → Meth	7 → Hept
2 → eth	8 → Oct
3 → prop	9 → Non
4 → but	10 → Dec
5 → pent	11 → Unde
6 → Hex	12 → dodec

* Alkane: $(C_n H_{2n+2})$ (C-C)

n=1 ⇒ CH ₄ → methane	n=3 ⇒ C ₃ H ₈ → propane
n=2 ⇒ C ₂ H ₆ → ethane	n=4 ⇒ C ₄ H ₁₀ → butane

* Alkene: $(C_n H_{2n})$ → (C=C)

n=2 ⇒ C ₂ H ₄ ⇒ ethene
n=3 ⇒ C ₃ H ₆ ⇒ propene

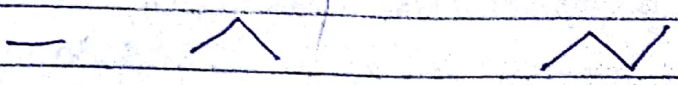
* Alkyne: $(C_n H_{2n-2})$ → (C≡C)

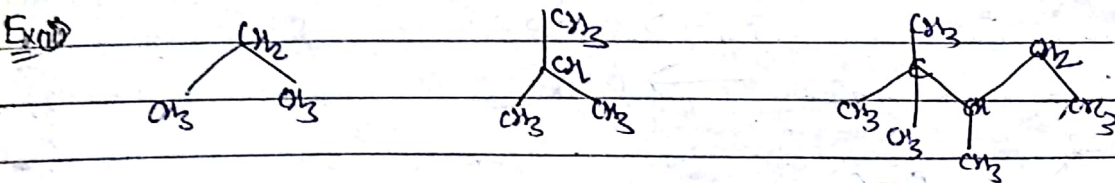
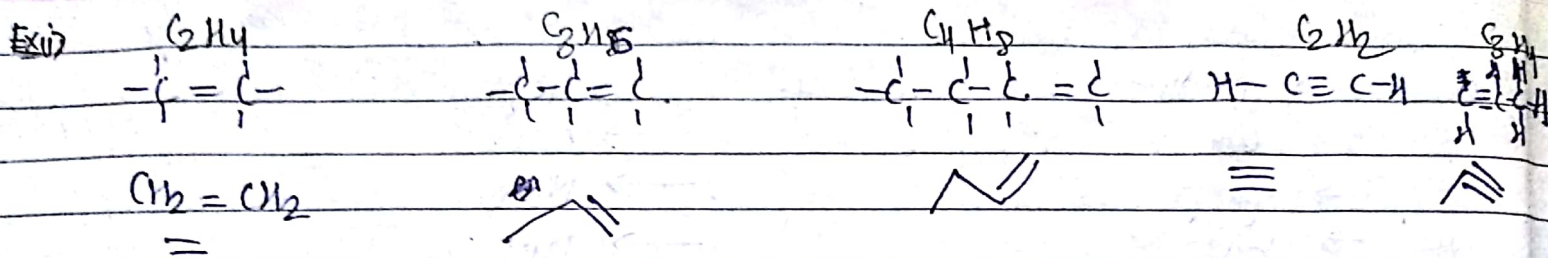
n=2 ⇒ C ₂ H ₂ ⇒ Ethyne
n=3 ⇒ C ₃ H ₄ ⇒ propyne

Formula: Methane Ethane propane Butane

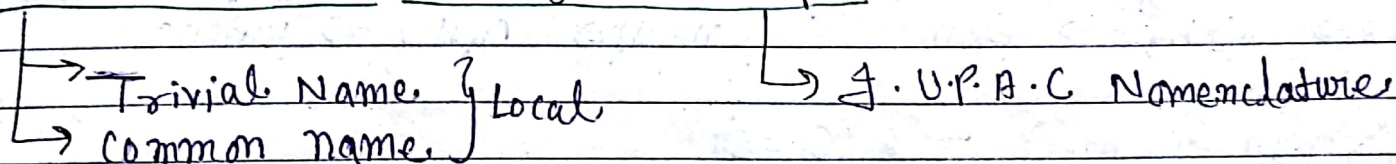
Molecular formula → CH ₄	C ₂ H ₆	C ₃ H ₈	C ₄ H ₁₀
Structural formula → $H-C(H)(H)-H$	$-C-C-$	$-C-C-C-$	$-C-C-C-C-$

Condensed formula → CH₄ C₂H₆ C₃H₈ C₄H₁₀

Bond line formula → 

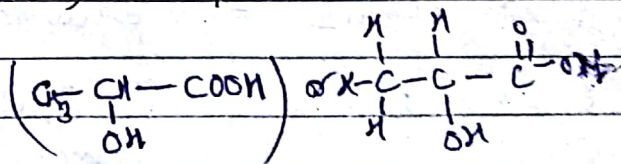



(3) Nomenclature of organic compounds -



CH_4 (Methane)	$HCOOH$	CH_3-OH	$\begin{array}{c} COOH \\ \\ COOH \end{array}$
\rightarrow Marsh gas	\rightarrow Formic acid	\rightarrow wood spirit	\rightarrow oxalic acid
\rightarrow Natural gas (CNG)	\rightarrow Formica	\rightarrow (Destructive distillation of wood)	\rightarrow Oxidus
\rightarrow Bio gas	\downarrow		
\rightarrow Fire damp.	Red Ant		

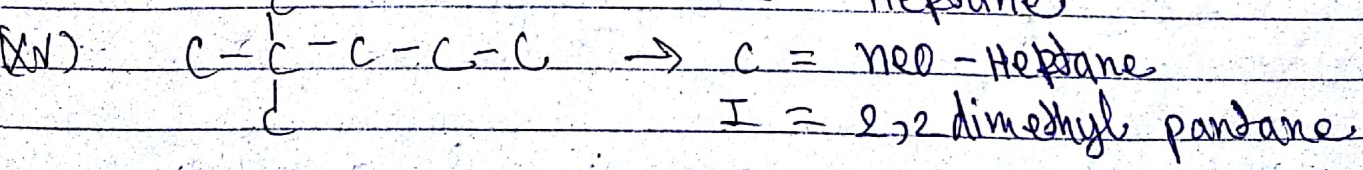
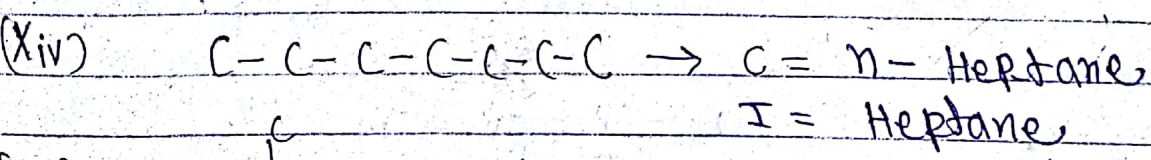
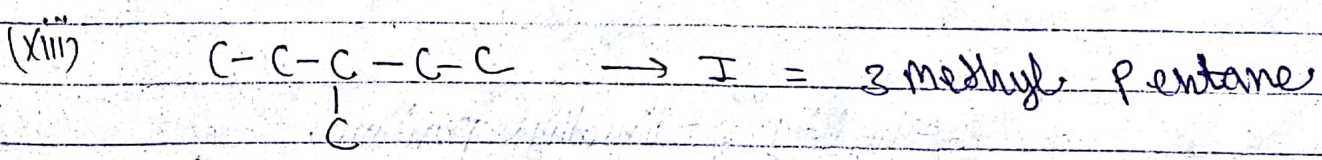
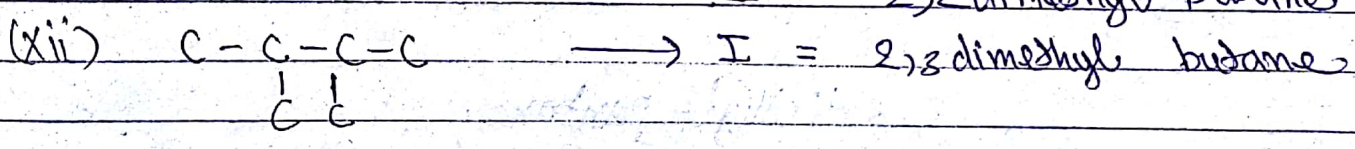
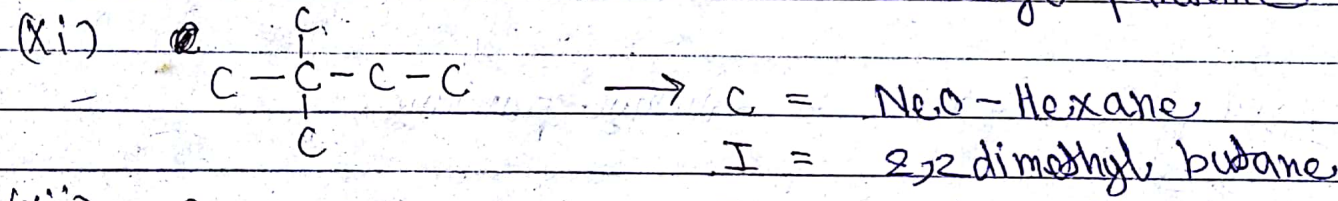
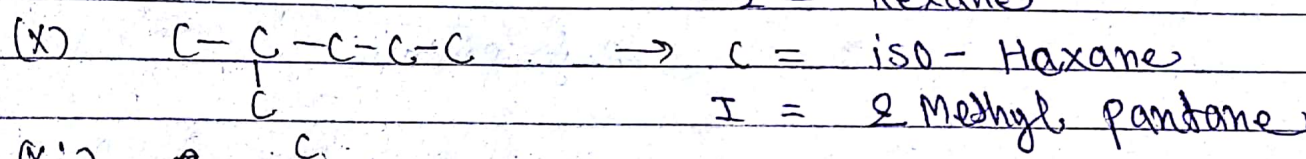
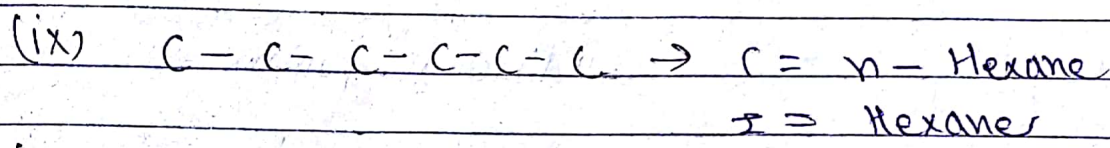
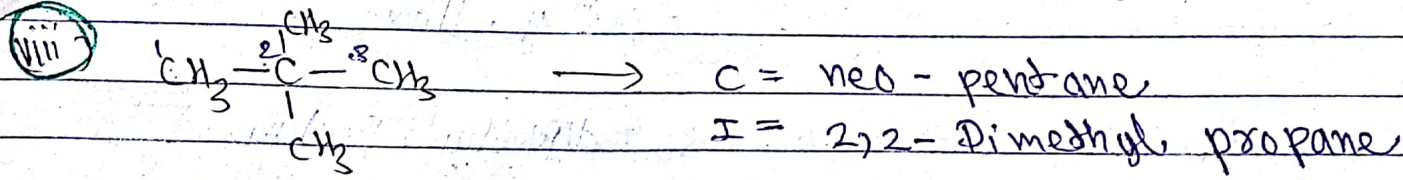
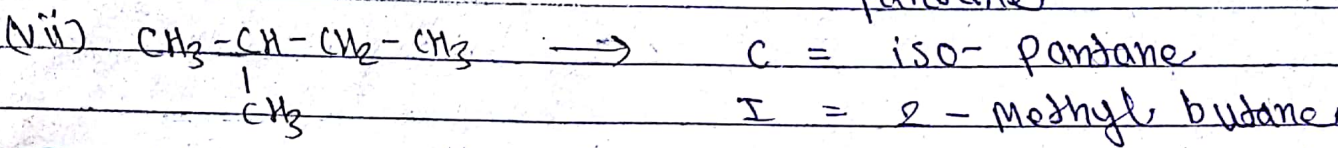
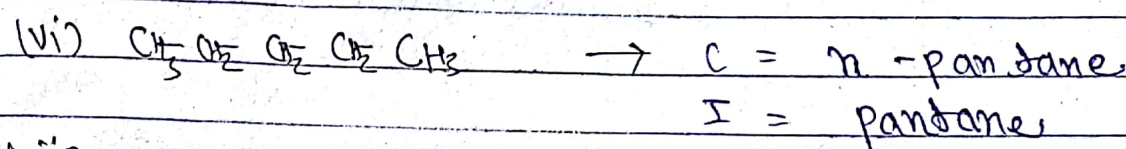
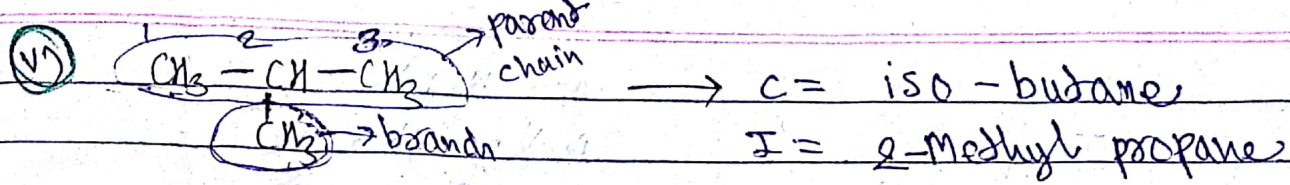
\rightarrow Lactic acid \rightarrow curd \rightarrow Lacto

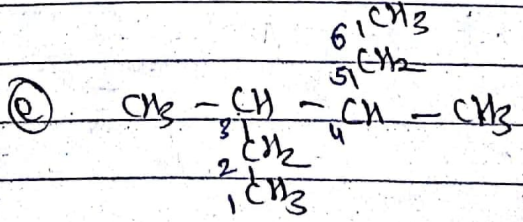
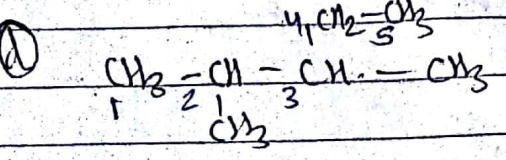
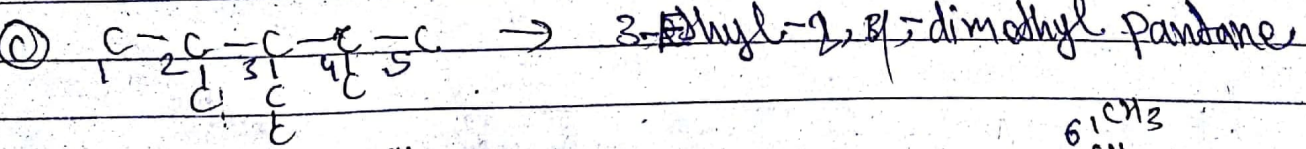
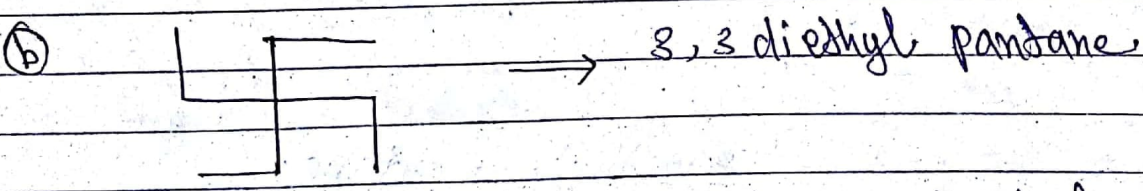
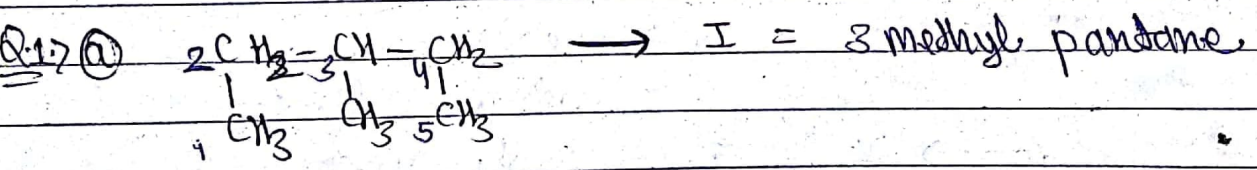
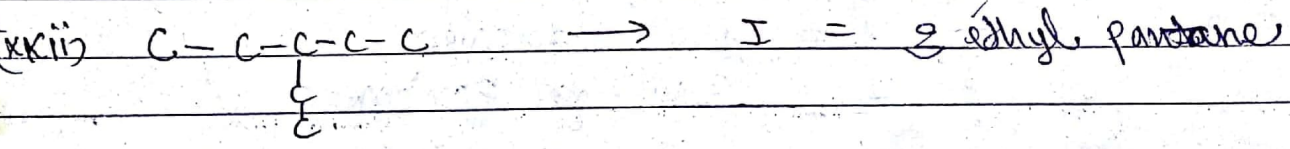
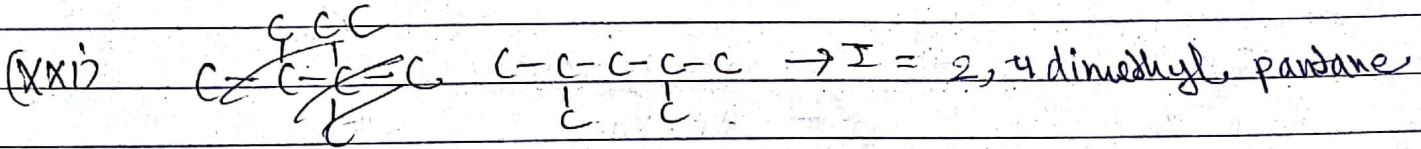
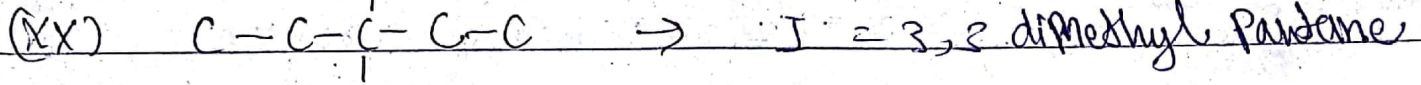
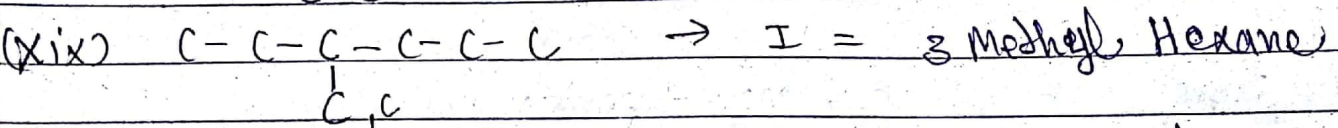
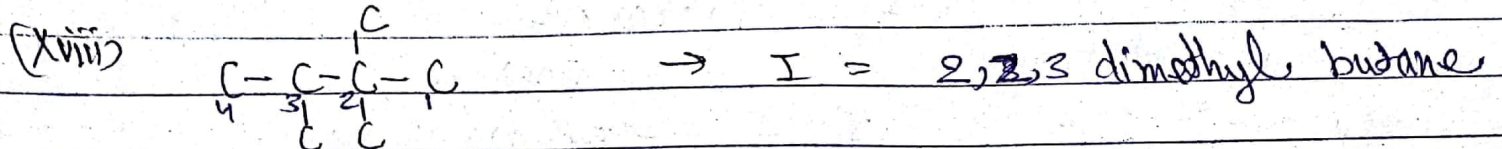
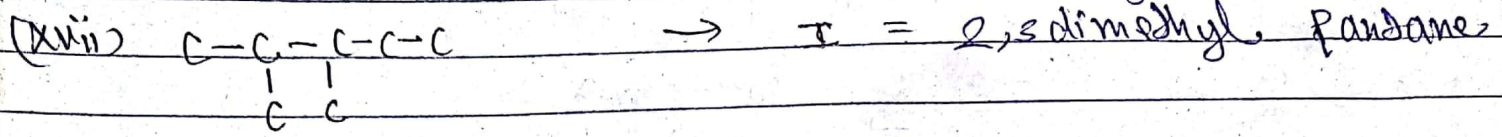
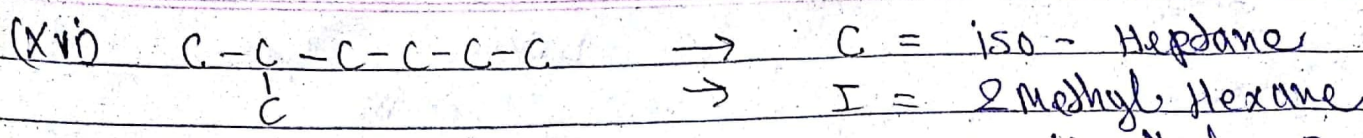


~~Imp~~  \rightarrow Phenol \rightarrow Phene \rightarrow Carboic acid \rightarrow phenic acid \rightarrow phenylic acid \rightarrow hydroxybenzene

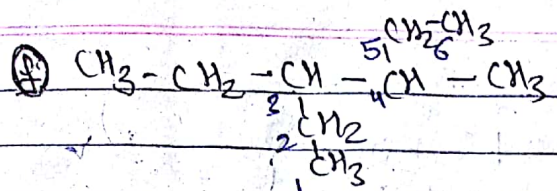
* Alkane, i.e. CH_4 - Methane
 (ii) CH_3-CH_3 - Ethane
 (iii) $CH_3-CH_2-CH_3$ - Propane
 (ix) $CH_3-CH_2-CH_2-CH_3$ - Butane
 I - Butane

→ Find C. no. abhi mathe 1

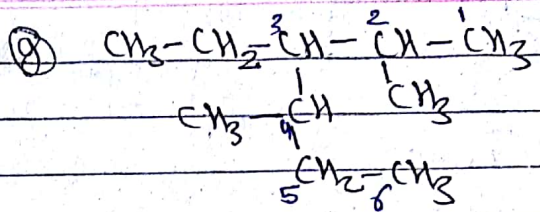




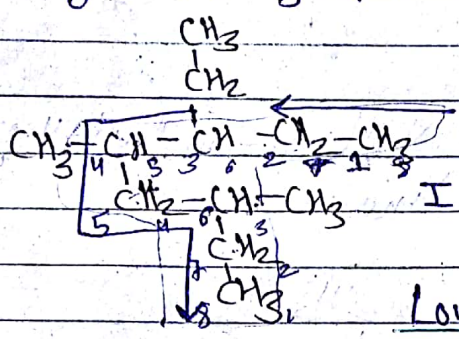
$\rightarrow 2,3\text{-dimethyl pentane} \rightarrow 3,4\text{-dimethyl pentane Hexane}$



→ 3-ethyl-4-methylhexane

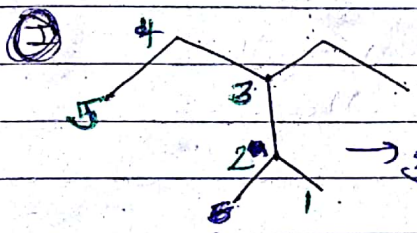


→ 3-ethyl-2,4-dimethylhexane

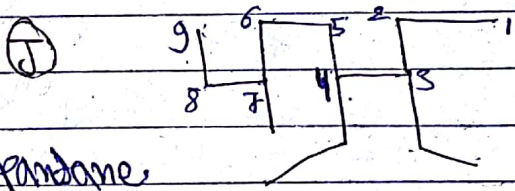


I → 3-ethyl-4,6-dimethyloctane

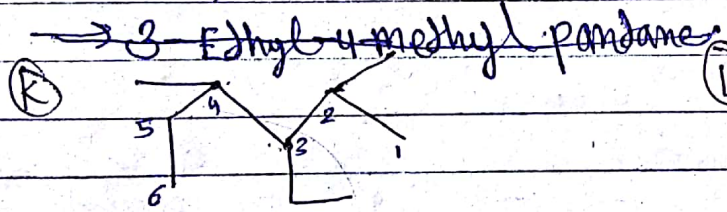
Lowest sum Rule $3+4+6=13$ vs $3+5+6=14$



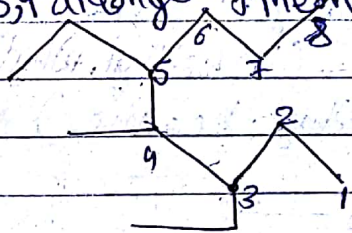
→ 3-ethyl-2-methylpentane



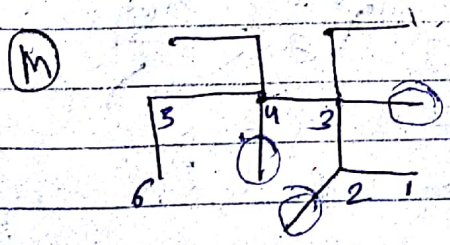
→ 3,4-diethyl-7-methylnonane



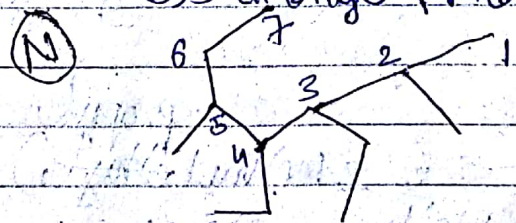
→ 3-ethyl-4-methylpentane



→ 3,5-diethyl-4-methyloctane

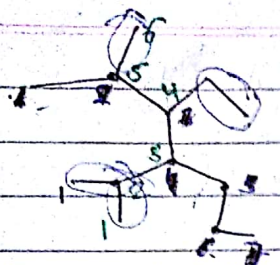


→ 3,4-diethyl-2,4-dimethylhexane



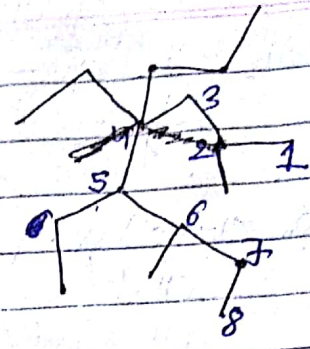
→ 3,4-diethyl-2,5-dimethylheptane

(M)



अनुसंधान के
द्वारा नो. प्रयोग
- 11/10/1

(N)

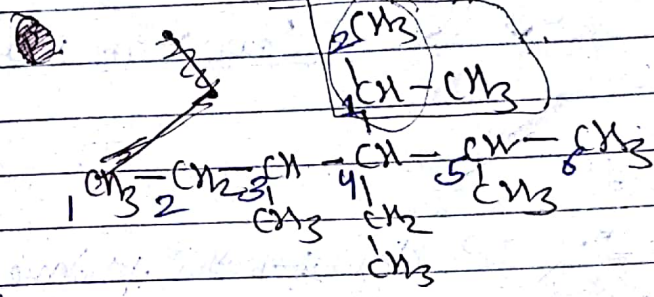
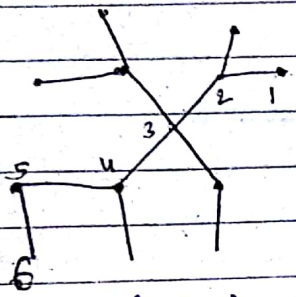


→ 3,4,4-trimethyl-2-methyl
Heptane

→ 4-ethyl-2,5-dimethyl-3-propyl hexane

→ 4,5-dimethyl-6-methyl-4-propyl
Octane

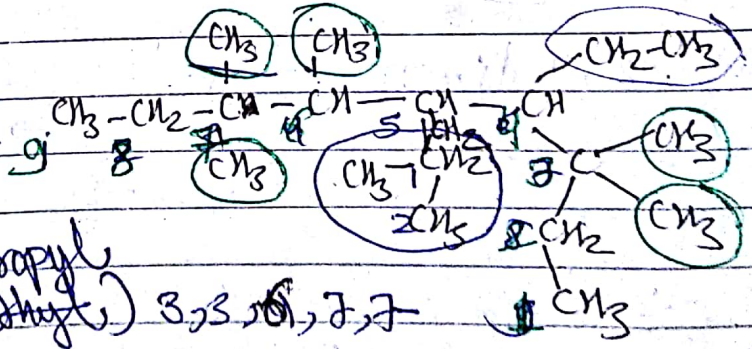
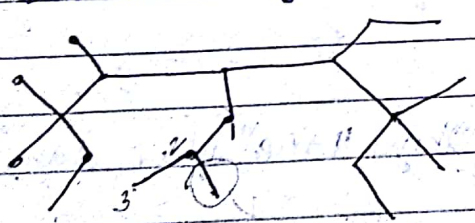
(O)



→ 3-ethyl-3-(isopropyl)-3-ethyl-2,4-dimethyl hexane

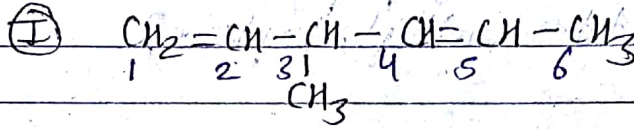
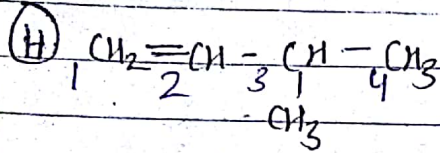
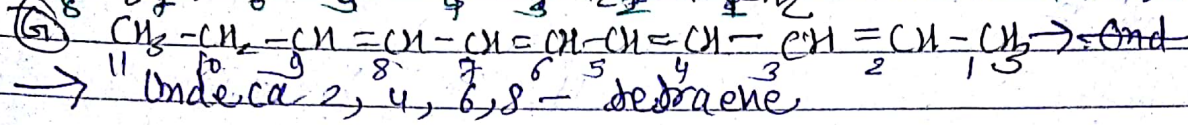
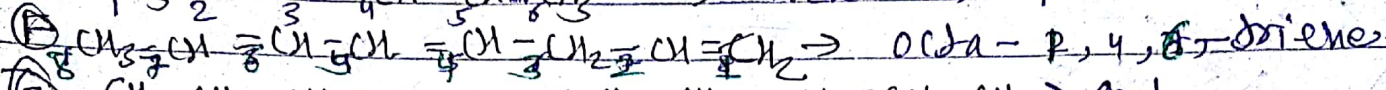
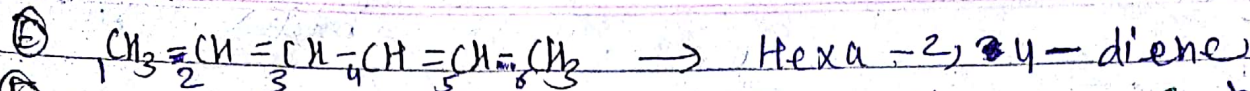
→ 3-ethyl-3-(1-methyl ethyl)-2,4-dimethyl hexane

(P)



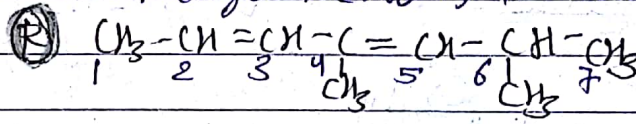
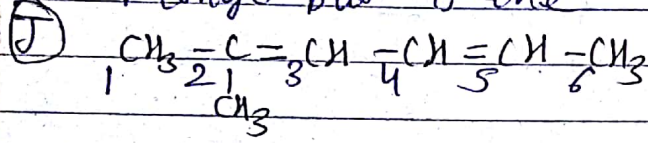
→ 4-methyl-5-(2-methyl propyl)-3,3,6,7,7-pentamethyl nonane

→ 5-(2-methyl propyl)-4-ethyl-3,3,6,7,7-pentamethyl nonane



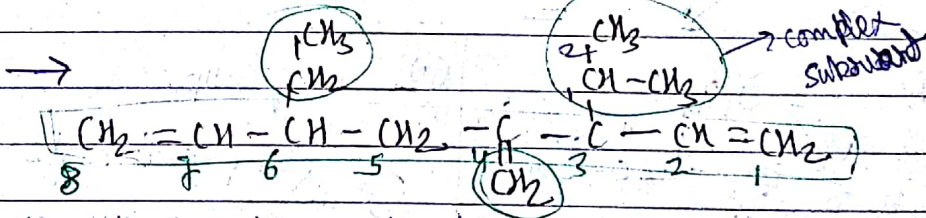
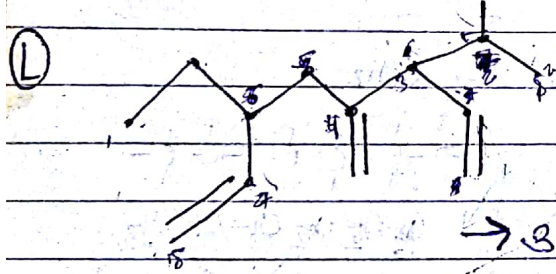
\rightarrow 3 Methyl but-1-ene

\rightarrow 3 Methyl Hexa-1,4-diene



\rightarrow 2 Methyl Hexa-2,4-diene

\rightarrow 4,6 dimethyl hepta-2,4-diene



\rightarrow 3 (1 methyl, Ethyl) - 6 ethyl, -4 methyloct-2,7-diene

Notes

- $\text{CH}=\text{CH}_2 \rightarrow$ Ethenyl (vinyl group)

- $\text{CH}=\text{CH} \rightarrow$ Ethenyl

= $\text{CH}_2 \rightarrow$ Methenyl

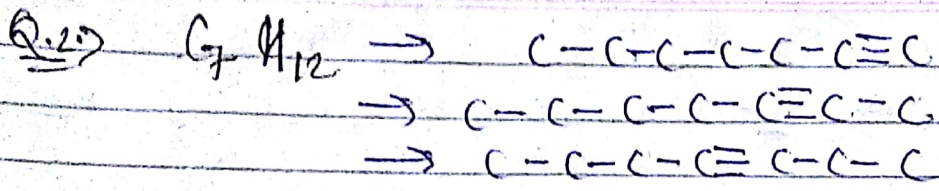
$\equiv \text{CH} \rightarrow$ Methyl

- $\text{CH}_2-\text{C}\equiv\text{CH} \rightarrow$ prop-2-yn-1-yl

- $\text{C}\equiv\text{C}-\text{CH}_3 \rightarrow$ propynyl

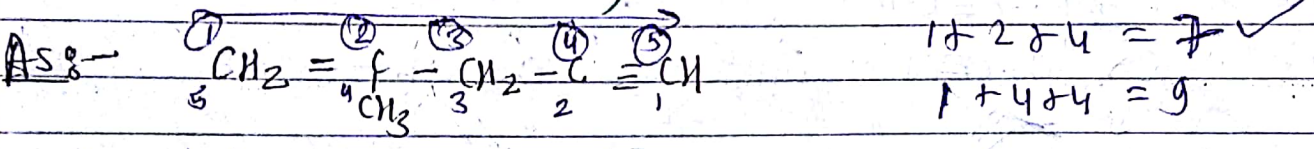
- $\text{CH}_2-\text{CH}=\text{CH}_2 \rightarrow$ prop-2-en-1-yl

- $\text{CH}=\text{CH}_2 \rightarrow$ propenyl

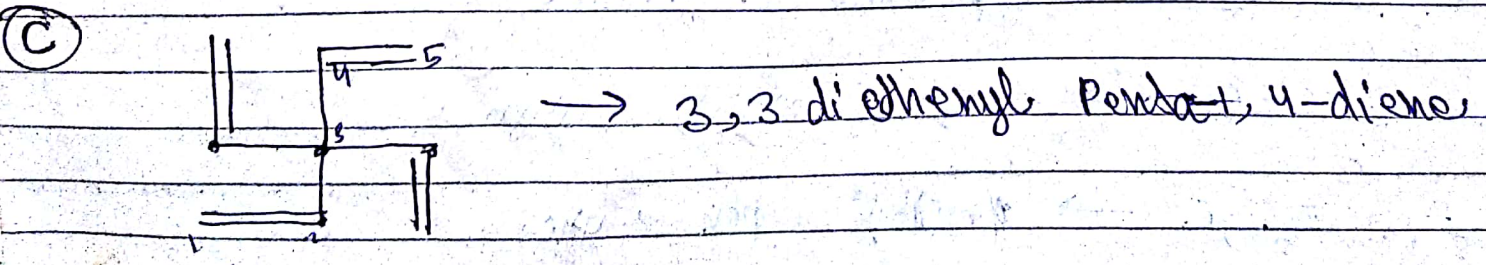
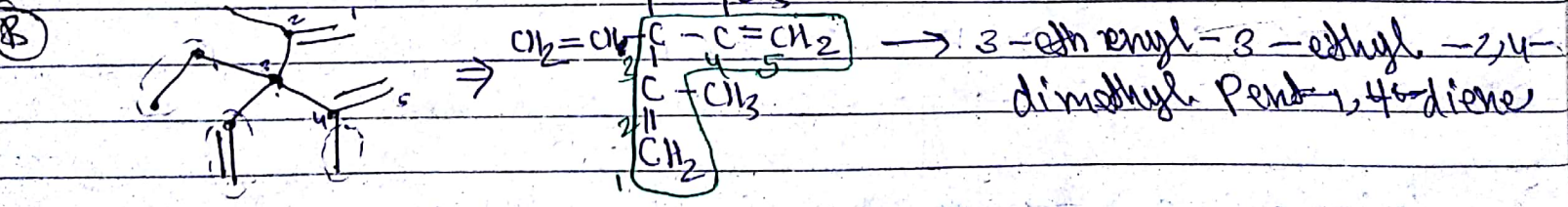
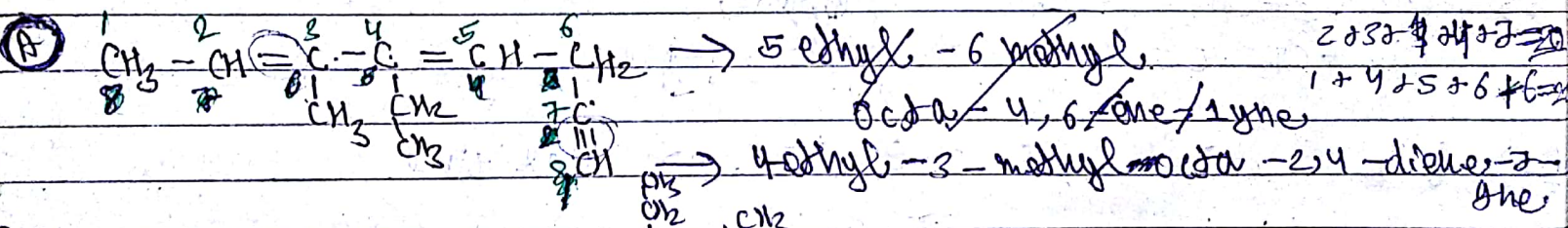


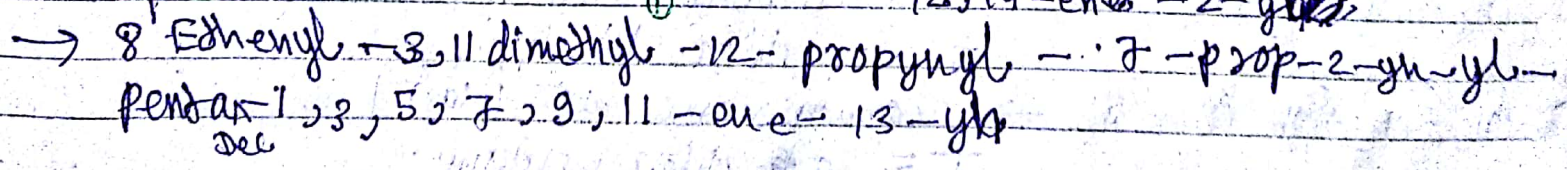
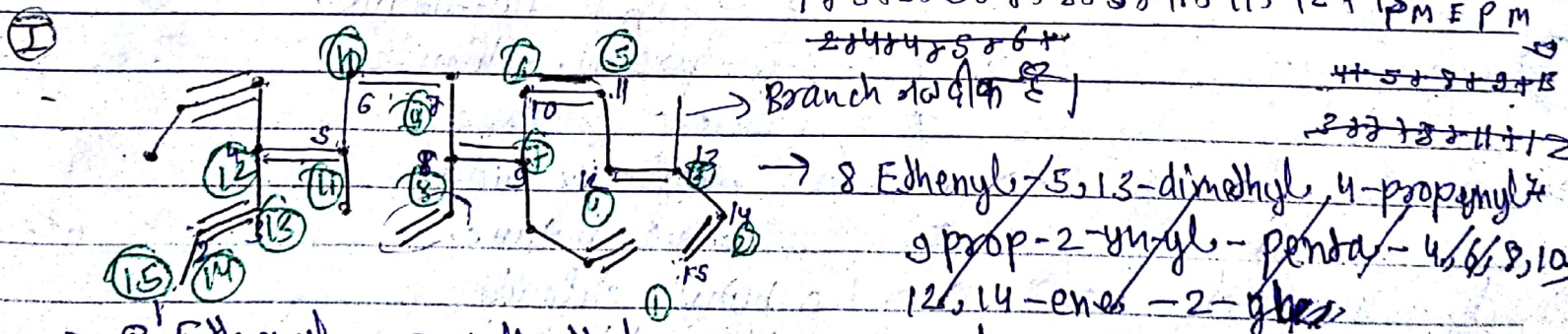
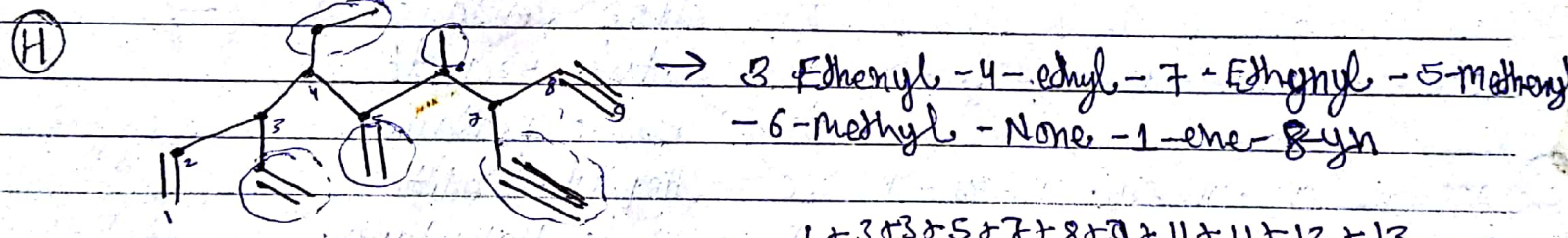
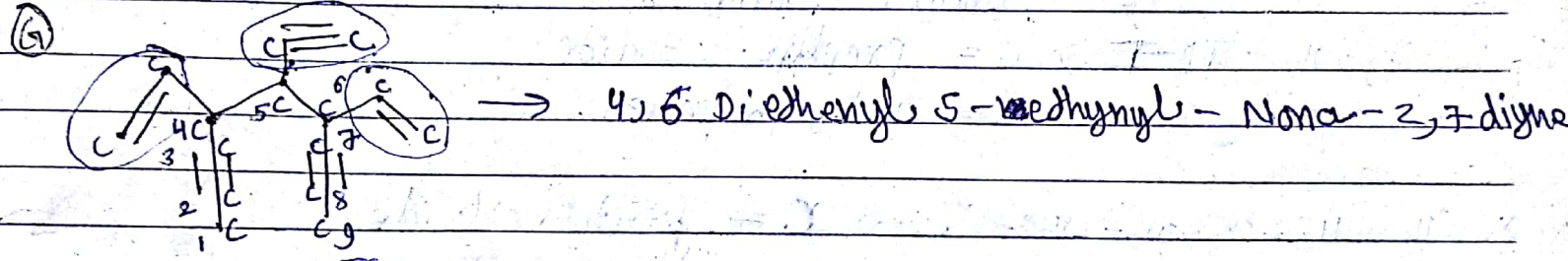
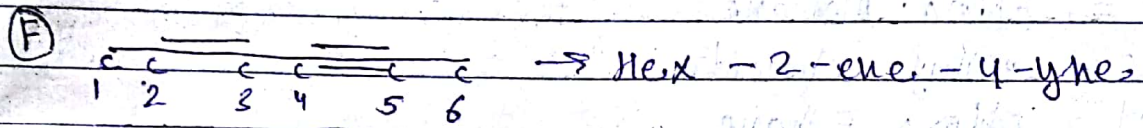
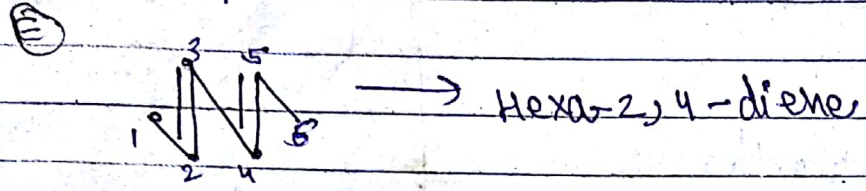
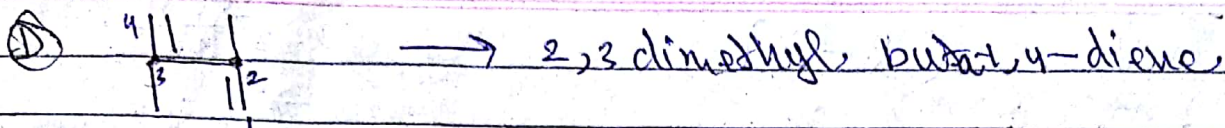
- (A) $^1CH_3 = ^2CH = ^3C \equiv ^4CH \rightarrow$ but-1-en-3-yne
- (B) $^5CH_3 - ^4CH = ^3CH - ^2C \equiv ^1CH \rightarrow$ pent-3-en-1-yne
- (C) $^5CH_3 - ^4C \equiv ^3C - ^2CH = ^1CH_2 \rightarrow$ pent-1-ene-3-yne
- (D) $^1CH_2 = ^2CH = ^3CH - ^4C \equiv ^5C - ^6CH_2 \rightarrow$ Hex-2-ene-4-yne

Lowest sum rule



2-methyl pent-1-ene-4-yne





$O^{-2}, Cl^{-}, N^{-3}, Br^{-}, S^{-2}$

Halogen	Halide
X_2	X^{-}
$F_2 \rightarrow$ Fluorine	$F^{-} \rightarrow$ Fluoride
Cl_2	$Cl^{-} \rightarrow$ Chloride
Br_2	$Br^{-} \rightarrow$ Bromide
I_2	$I^{-} \rightarrow$ Iodide

IV Alkyl Halide :- (R-X)

\rightarrow I.U.P.A.C. \rightarrow Halo Alkane

Cl	\rightarrow Chloro
Br	\rightarrow Bromo
F	\rightarrow Fluoro
I	\rightarrow Iodo

Ex. (i) $CH_3-Cl \rightarrow C =$ methyl chloride
 $I =$ chloro methane

(ii) $CH_3-CH_2-Cl \rightarrow C =$ Ethyl chloride
 $I =$ chloro ethane

(iii) $CH_3-CH_2-CH_2-I \rightarrow C =$ Propyl iodide
 $I =$ Iodo propane

Q. (A) $CH_3-CH_2-CH_2-CH_2-CH_2-Cl \rightarrow C =$ pentyl chloride
 $I =$ chloro pentane

(B) $Br-CH_2-CH_2-CH_2-CH_2-CH_2 \rightarrow C =$ pentyl bromide
 $I =$ bromo pentane

(C) $I-CH_2-CH_2-CH_2-CH_2-CH_2-CH_2-CH_2 \rightarrow C =$ Heptyl iodide
 $I =$ Iodo Heptane

(D) $CH_3-CH(Cl)-CH_3 \rightarrow C =$ iso-propyl chloride
 $I =$ 2-chloro propane

(E) $CH_3-CH_2-CH(Br)-CH_3 \rightarrow C =$ iso butyl bromide
 $I =$ 2-bromo butane

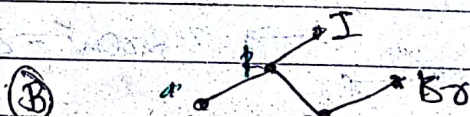
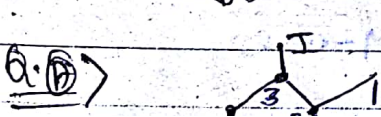
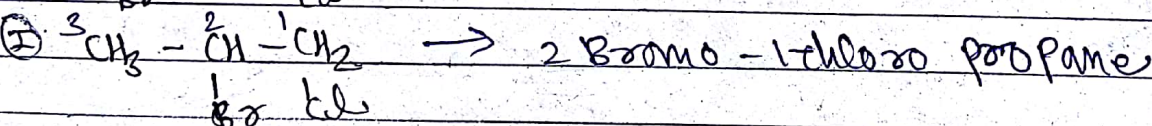
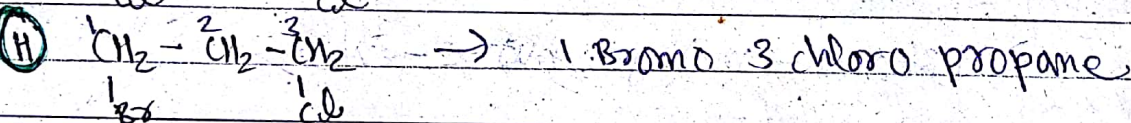
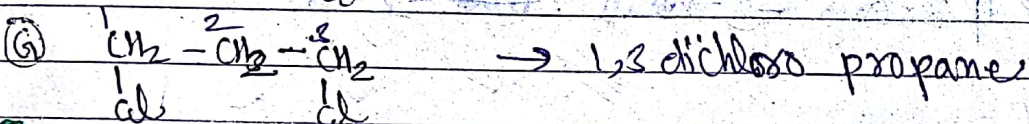
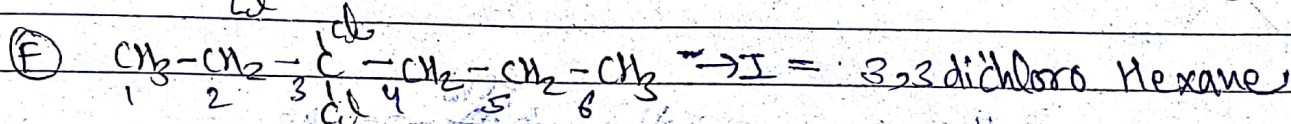
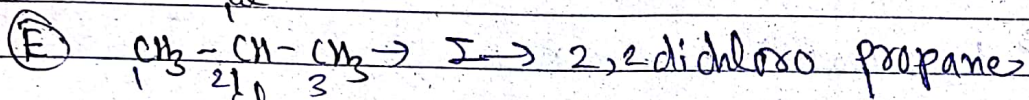
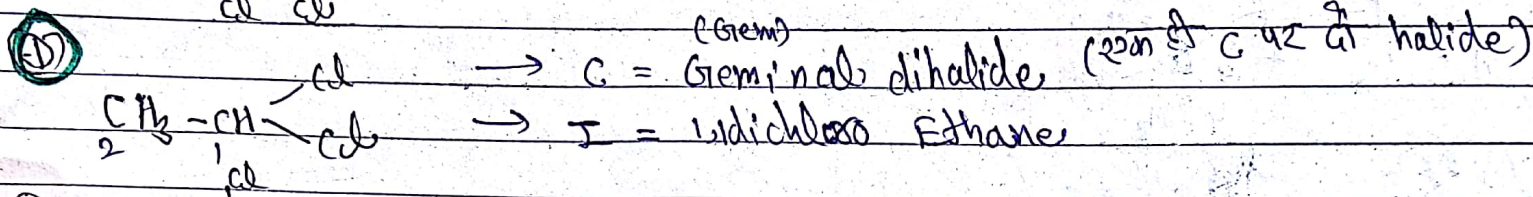
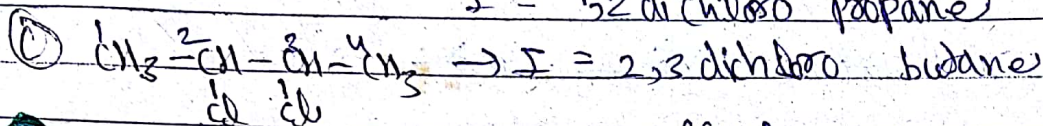
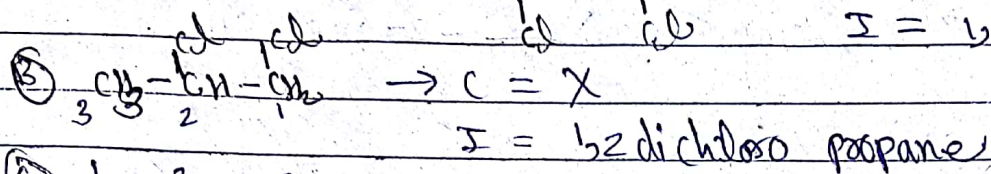
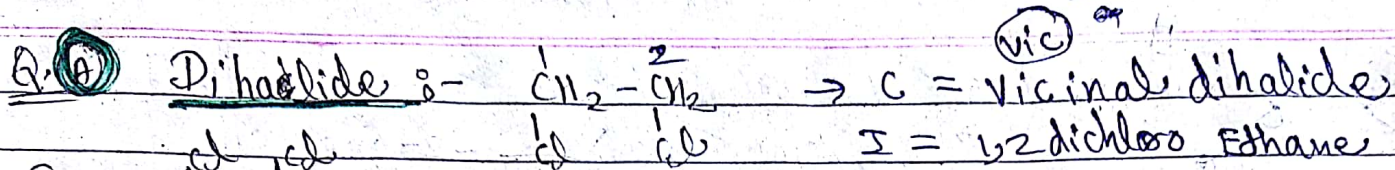
(F) $CH_3-CH_2-CH_2-CH_2-Cl \rightarrow C =$ iso butyl chloride
 $I =$ 2-chloro butane

(G) $CH_3-CH_2-CH_2-CH_2-CH_2-Cl \rightarrow$ common name, 1st part
 $I =$ 3-chloro pentane

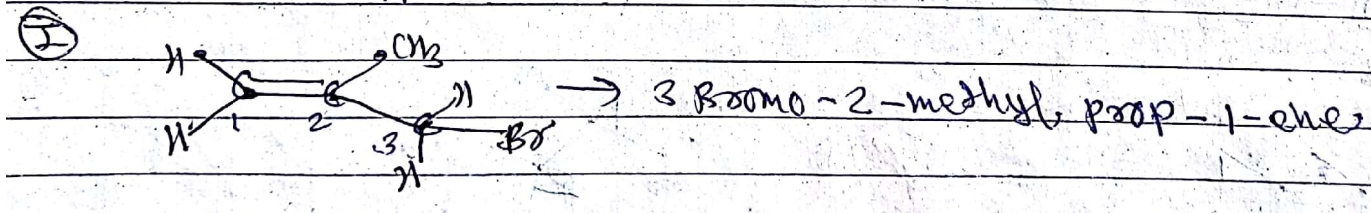
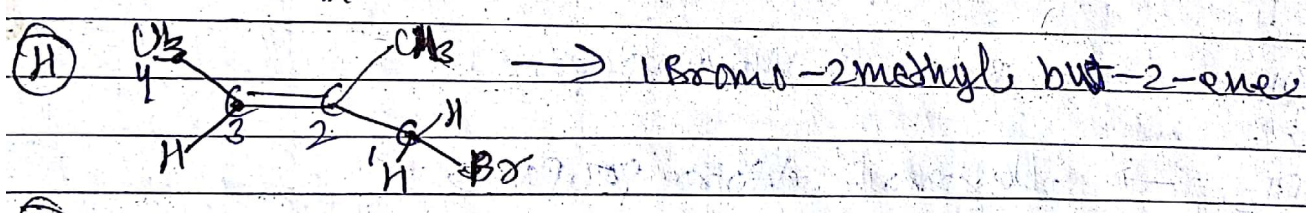
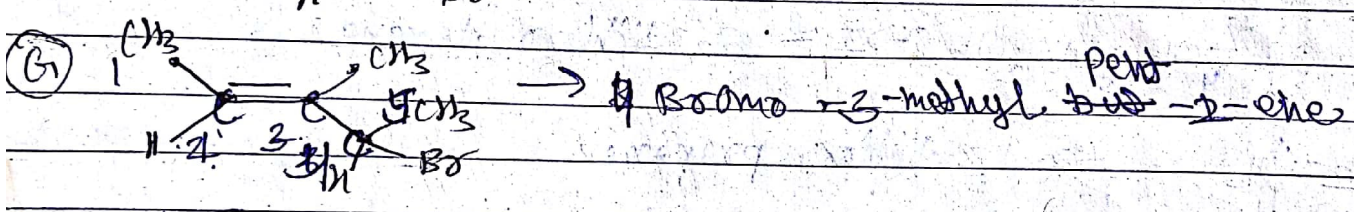
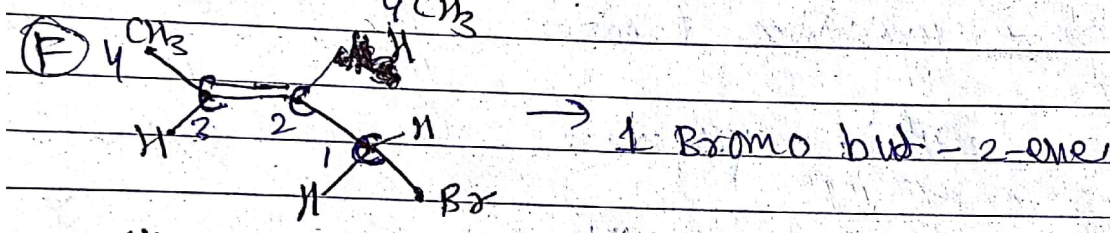
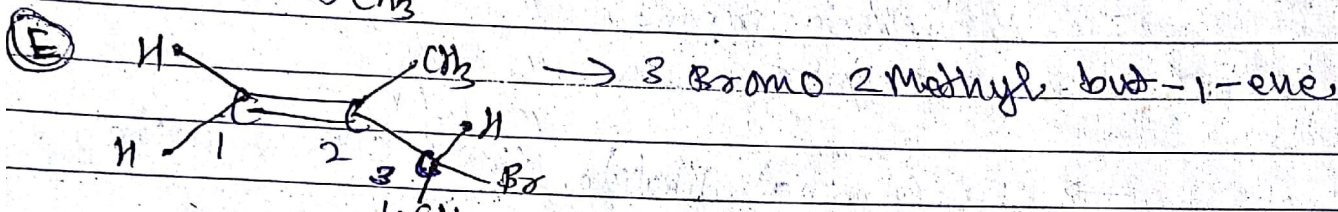
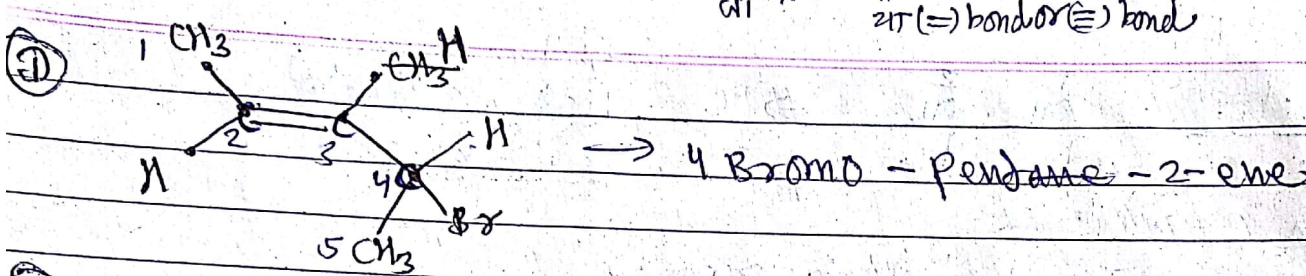
(H) $CH_3-CH_2-CH_2-CH_2-CH_2-I \rightarrow C =$ Iso hexyl iodide
 $I =$ 2-iodo hexane

(I) $CH_3-CH_2-CH_2-CH_2-CH_2-CH_2-CH_2-CH_2-I \rightarrow C =$ X
 $I =$ 4-iodo octane

Note: - In case of C के लिए।



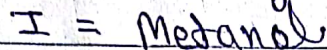
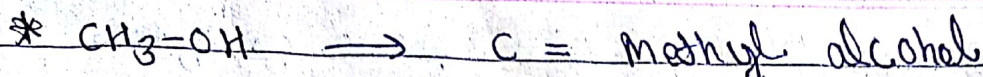
Note: - 21st 21st part is = 21st habit of 21st part -
 → 21st same position is 21st (=) bond or alkyl part
 → 21st 21st - 2. 21st is 21st, 21st, 21st is 21st No. alkyl
 21st (=) bond or (=) bond



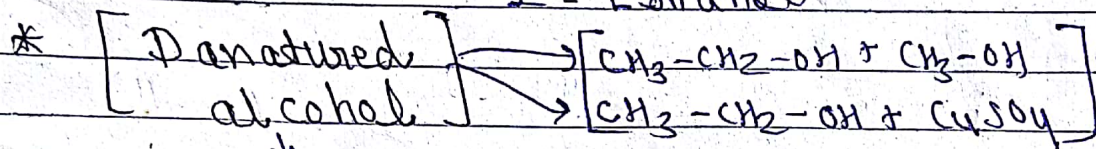
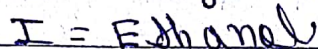
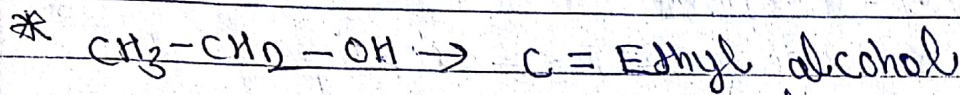
(V) Alcohols : → (R-OH)

C = Alkyl alcohol
 ↓
 Alkane + suffix = ol
 ↓
 I = Alkanol

suffix = ol
 prefix = hydroxy



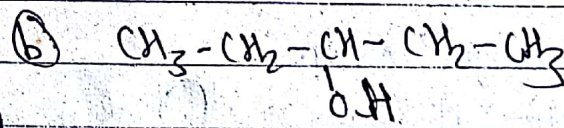
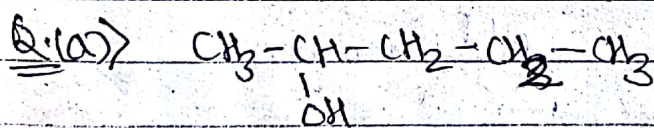
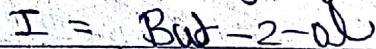
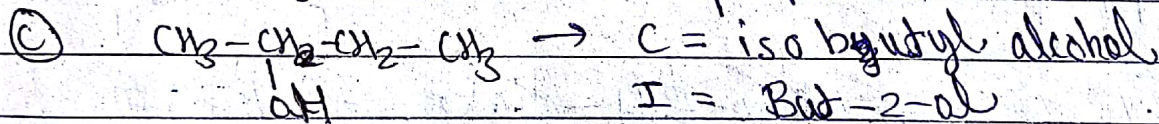
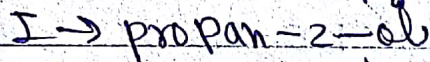
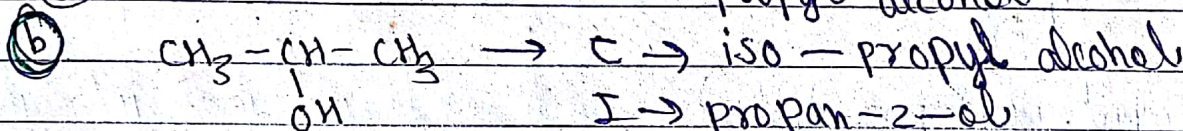
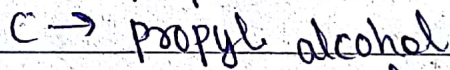
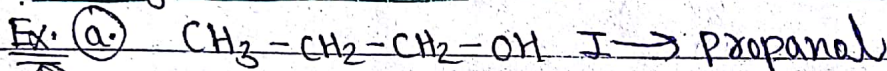
↳ poisonous alcohols



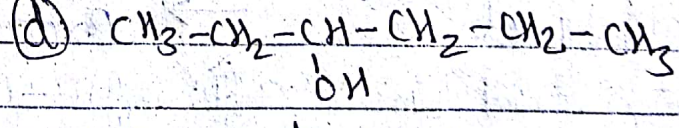
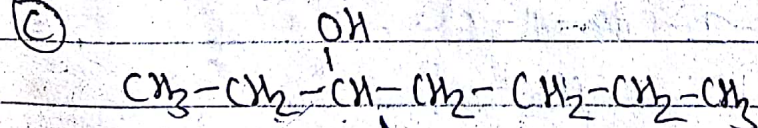
↓

unfit for drinking

Monohydric alcohol \rightarrow

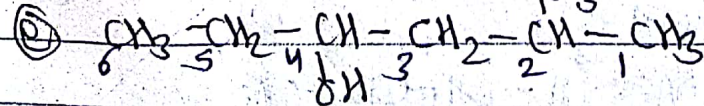


C = iso-pentyl alcohol, I = Pent-2-ol | I = Pent-3-ol

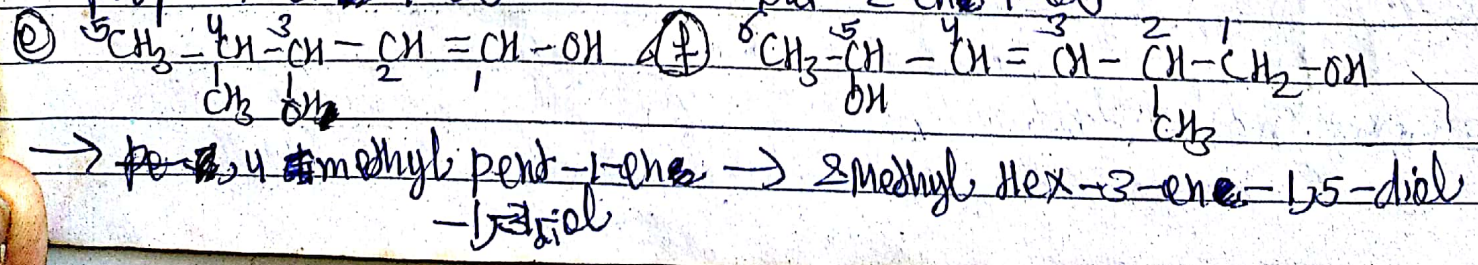
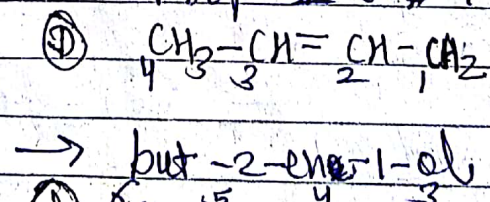
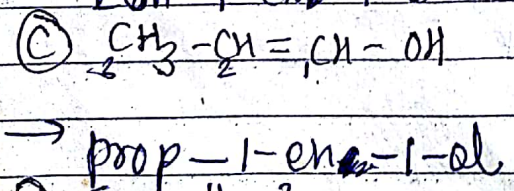
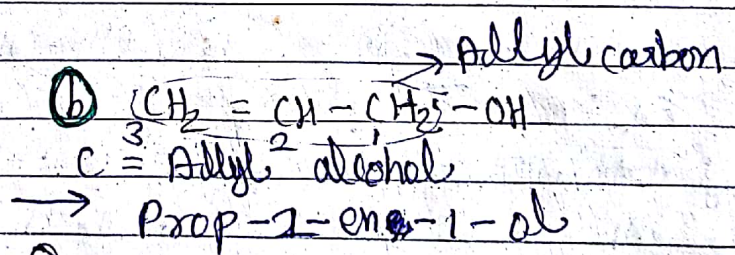
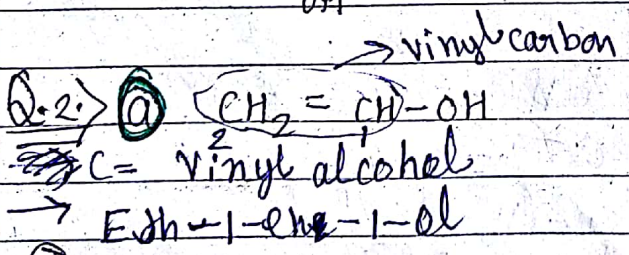
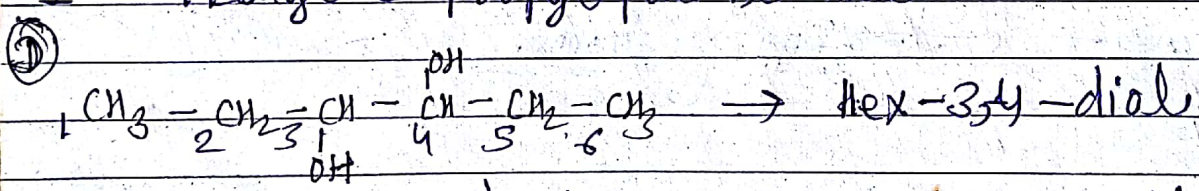
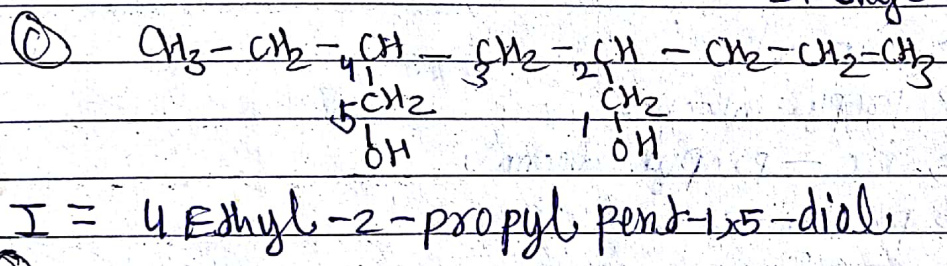
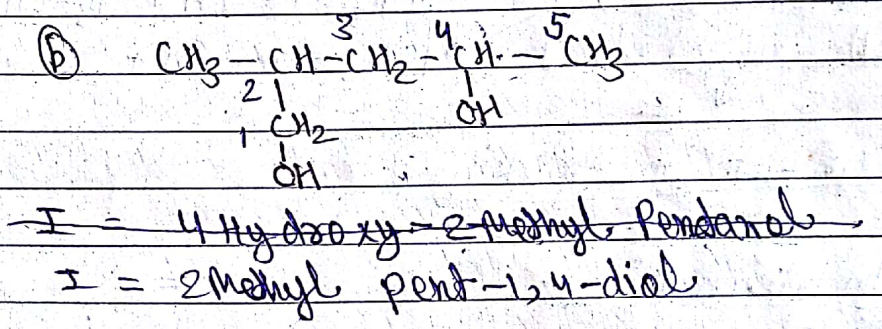
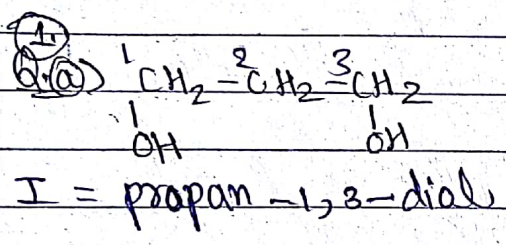
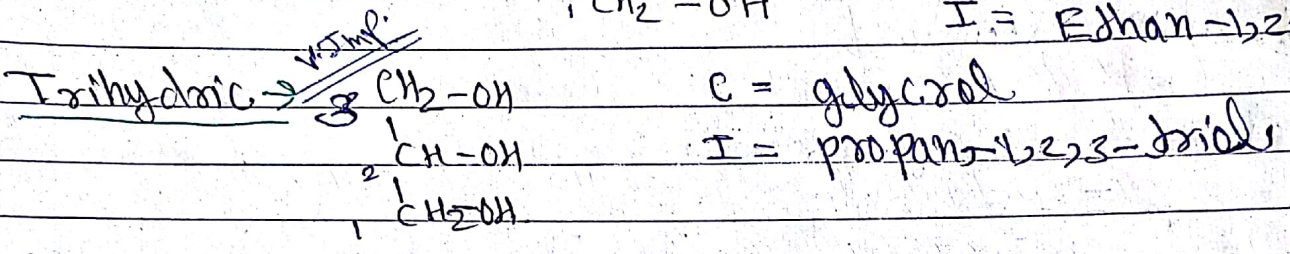
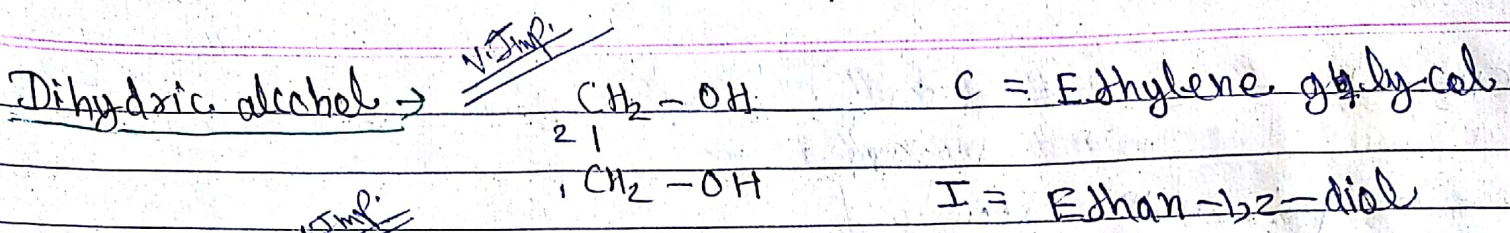


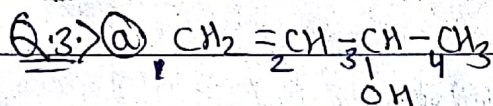
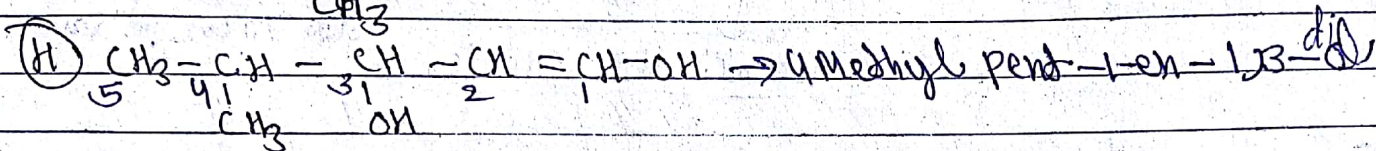
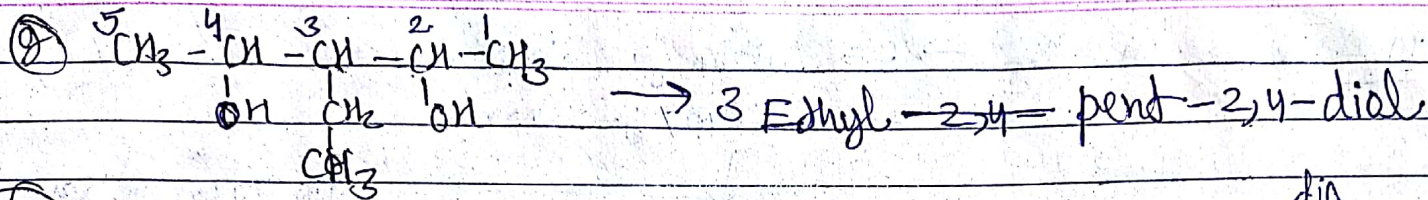
\rightarrow Hept-3-ol

\rightarrow Hex-3-ol

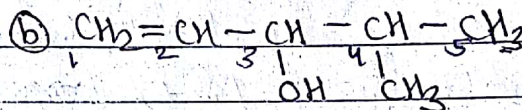


\rightarrow 2 methyl Hex-4-ol

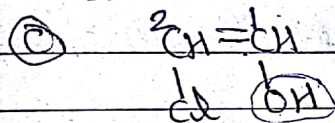




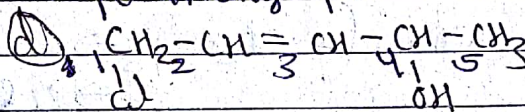
\rightarrow But-1-en-3-ol



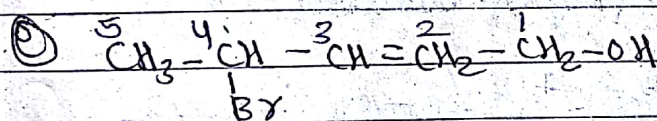
\rightarrow 4 methyl pent-1-en-3-ol



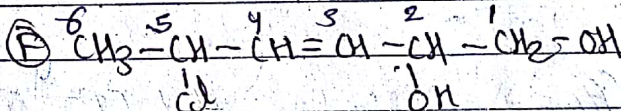
\rightarrow 2 chloro Eth-1-en-1-ol



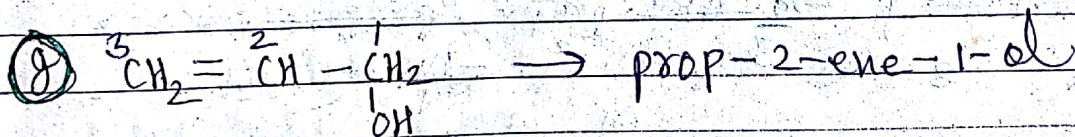
\rightarrow 1 chloro pent-2-en-4-ol



\rightarrow 4 Bromo pent-2-en-1-ol



\rightarrow 5 chloro Hex-3-en-1,2-diol



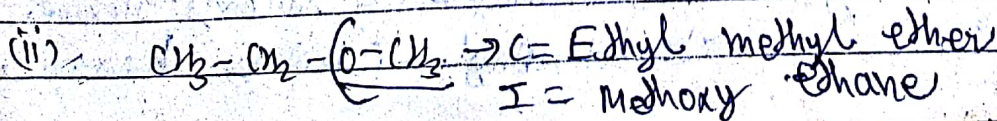
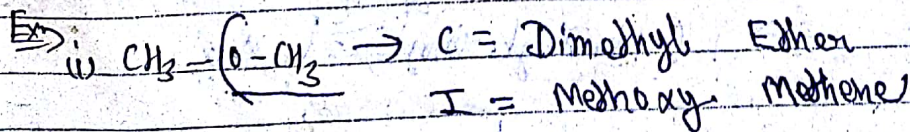
VI

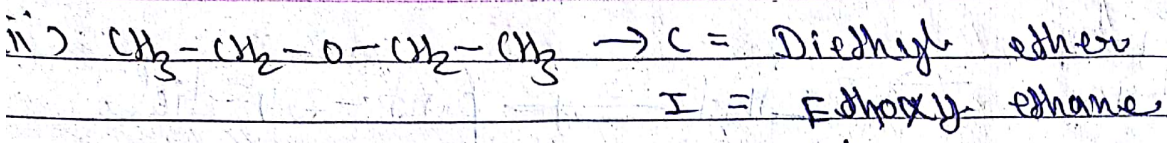
Ether (R-O-R)

C = Dialkyl Ether

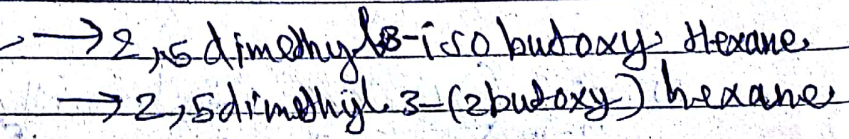
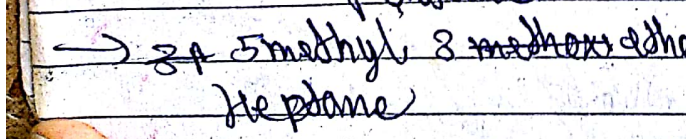
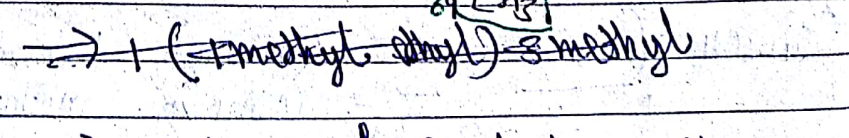
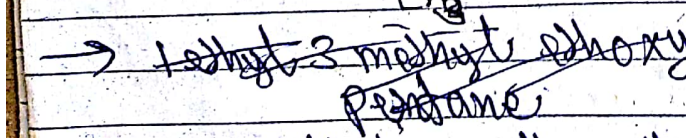
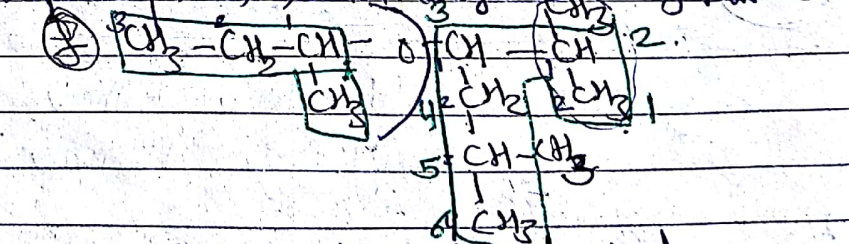
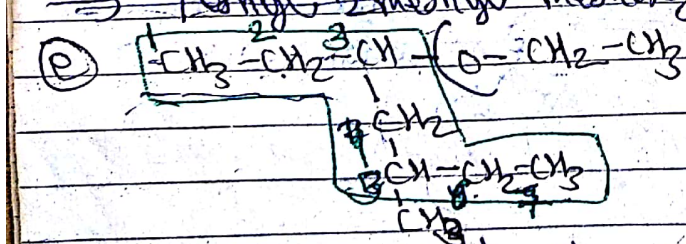
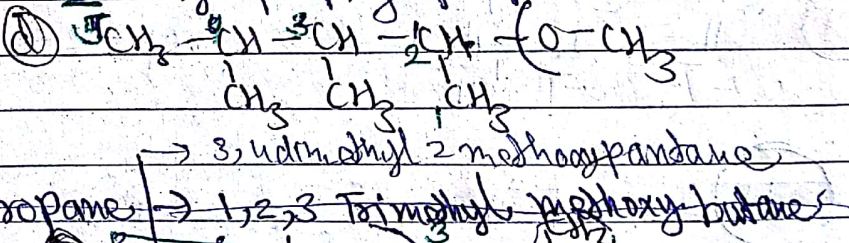
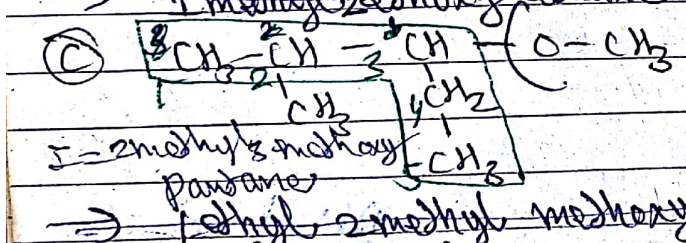
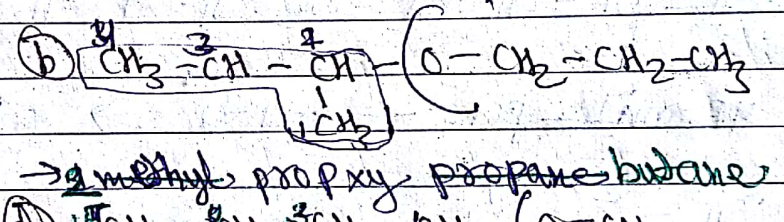
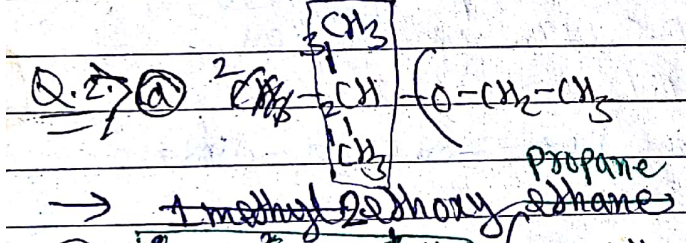
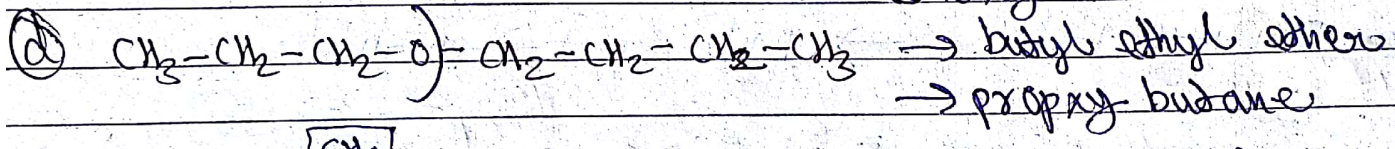
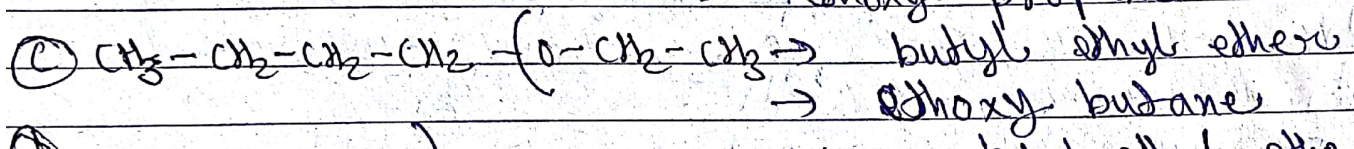
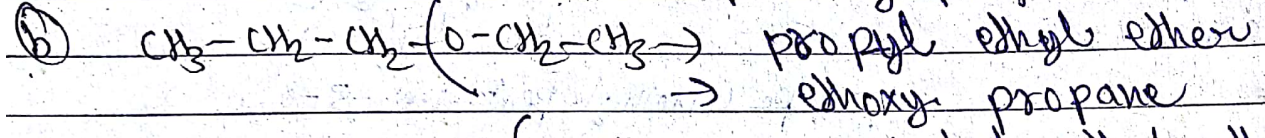
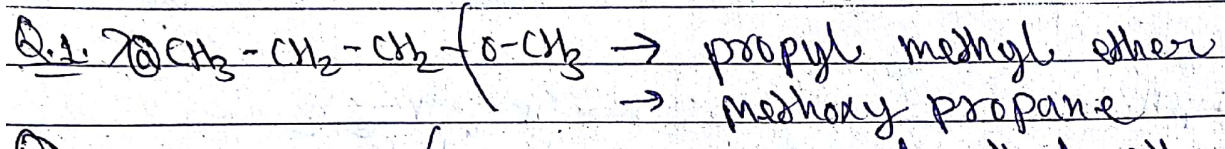
I = Alkoxy Alkane

Note: - OR Alkyl ether
 OR Alkyl ether
 OR Alkyl ether

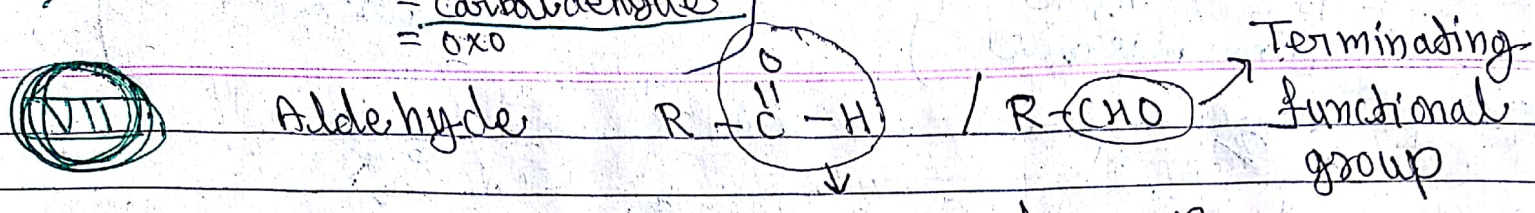




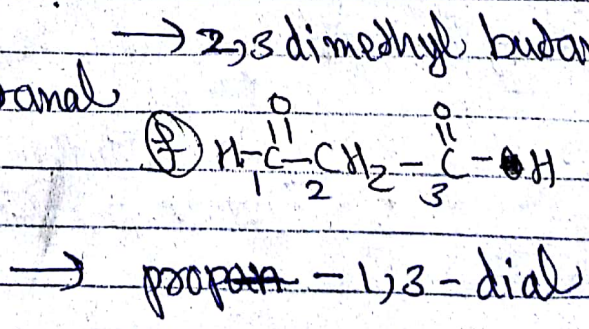
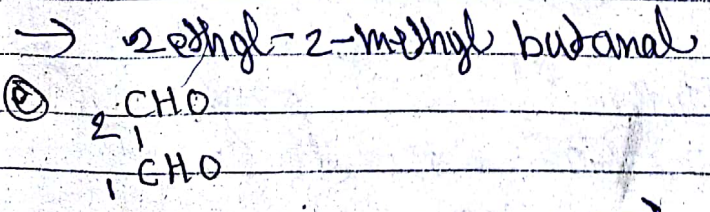
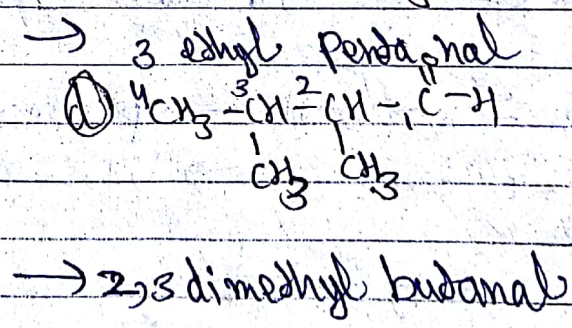
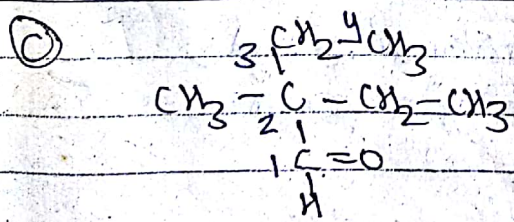
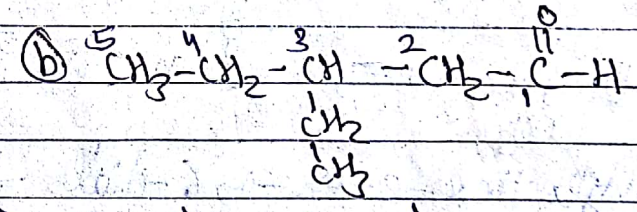
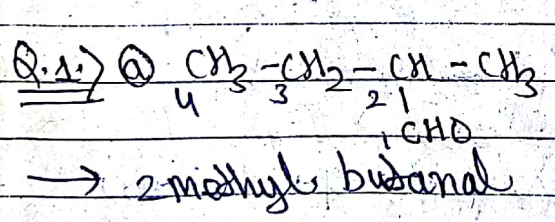
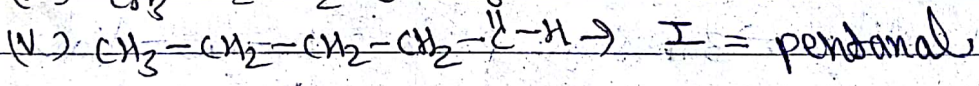
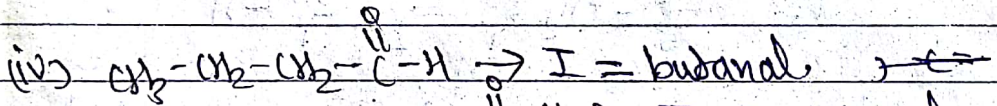
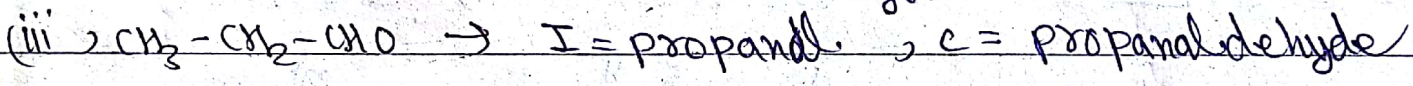
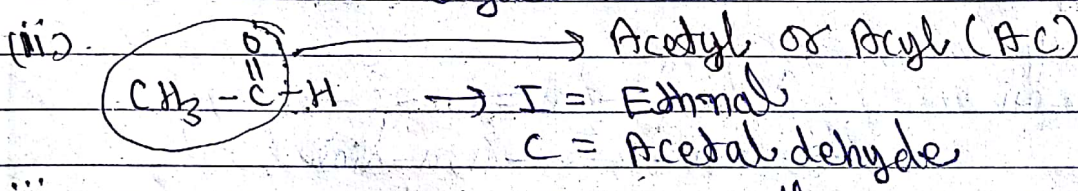
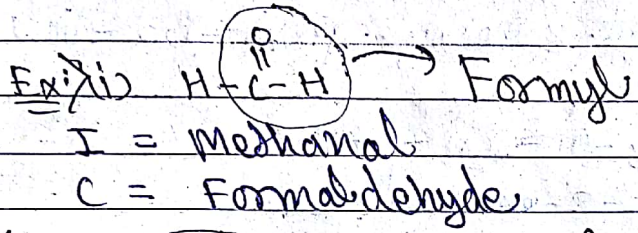
↓
Local Anesthetic diethyl ether



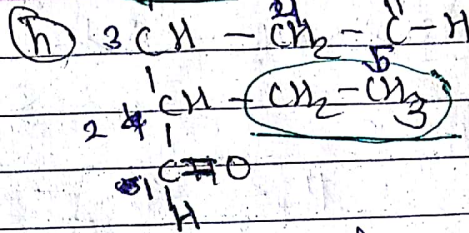
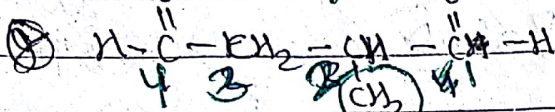
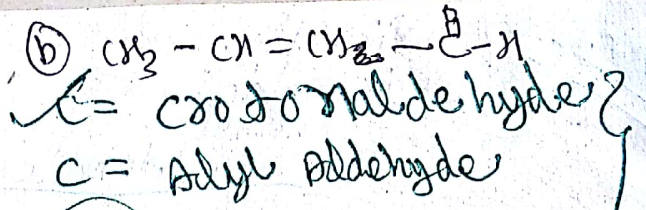
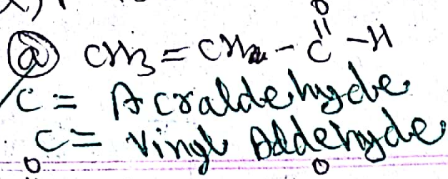
M. Imp. \Rightarrow suffix = al
 \Rightarrow prefix = aldo
 = formyl
 = carbaldehyde
 = oxo



Functional group
 $I = \text{Alkanal} \rightarrow \text{al} = \text{Alkanal}$



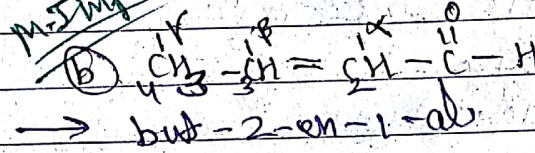
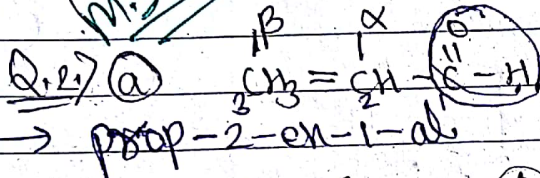
α, β unsaturated carbonyl compounds



→ 3 Methyl but-1,4-dial

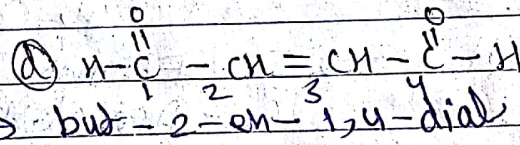
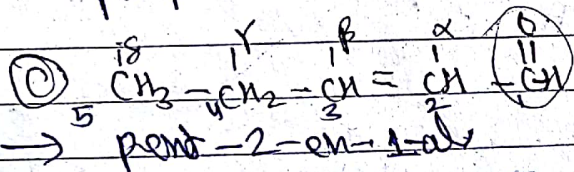
→ 2 ethyl 3 Methyl, pent-1,5-dial

M.I.M.P.



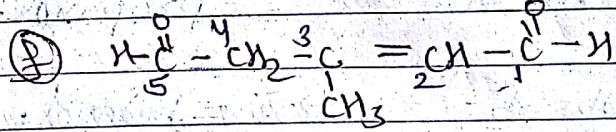
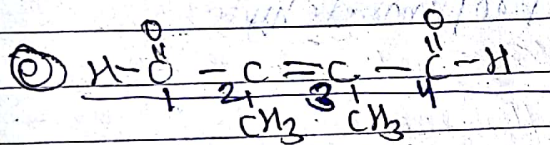
→ prop-2-en-1-al

→ but-2-en-1-al

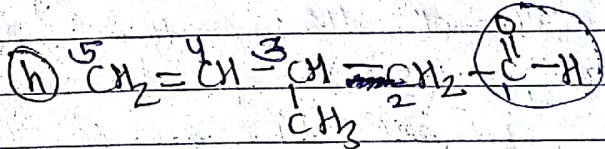
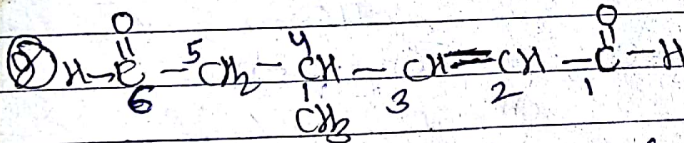


→ pent-2-en-1-al

→ but-2-en-1,4-dial

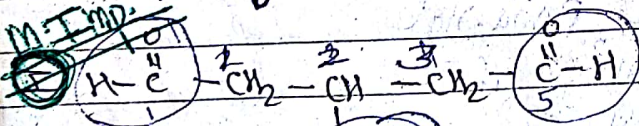


→ 2,3 dimethyl but-2-en-1,4-dial → 3 methyl pent-2-en-1,5-dial

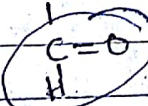


→ 4 methyl hex-2-en-1,6-dial → 3 methyl pent-4-en-1-al

M.I.M.P.

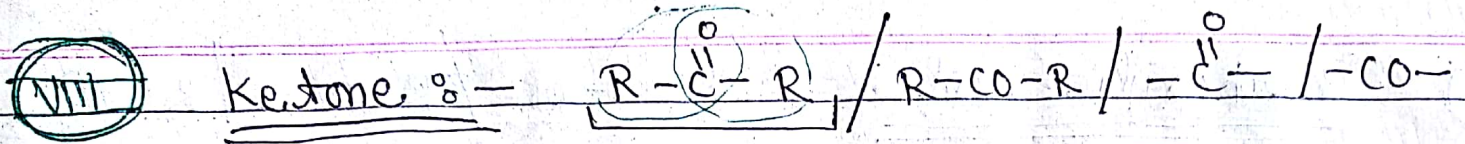


→ propan-1,2,3-tricarbaldehyde
 → 1,2,3-tricarbaldehyde propane



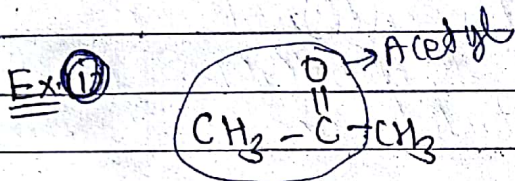
→ but-1,5-dial

Carbonyl group containing compound \rightarrow $\text{C}=\text{O}$
 Kato ueta carbon uze etat 3it etant & det of aldehyde etant 211



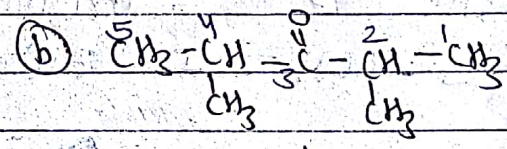
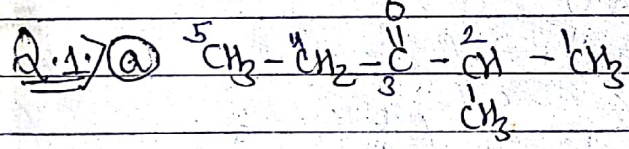
Suffix = one
 prefix = keto
OXO

I = Alkane = Alkanone
one



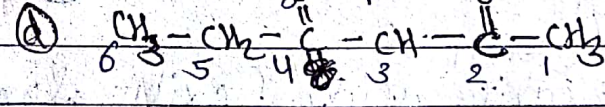
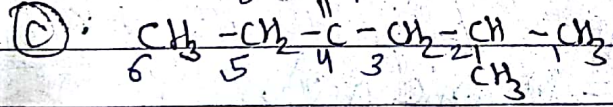
\rightarrow I = propanone
 \rightarrow C = Acetone
 \rightarrow C derived = Dimethyl ketone

- (ii) $\text{CH}_3-\text{CH}_2-\overset{\text{O}}{\parallel}{\text{C}}-\text{CH}_3$ \rightarrow I \rightarrow butanone, C = ethyl methyl ketone
- (iii) $\text{CH}_3-\text{CH}_2-\overset{\text{O}}{\parallel}{\text{C}}-\text{CH}_2-\text{CH}_3$ \rightarrow I = pent-2-one
- (iv) $\text{CH}_3-\text{CH}_2-\overset{\text{O}}{\parallel}{\text{C}}-\text{CH}_2-\text{CH}_3$ \rightarrow I = pent-3-one, C = Diethyl ketone



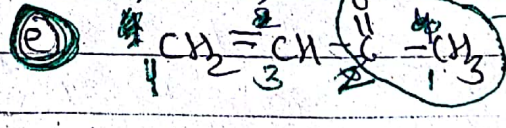
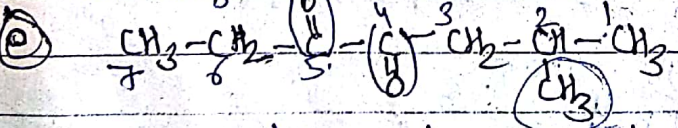
\rightarrow 2 methyl pent-3-one

\rightarrow 2,4 dimethyl pent-3-one



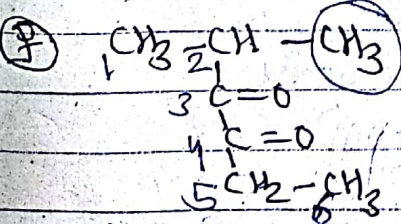
\rightarrow 2 methyl Hex-4-one

\rightarrow Hex-2,4-dione



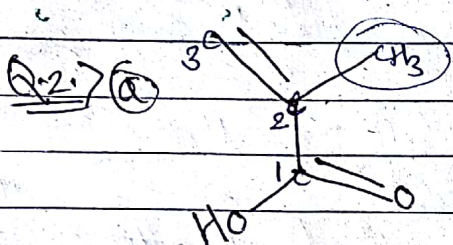
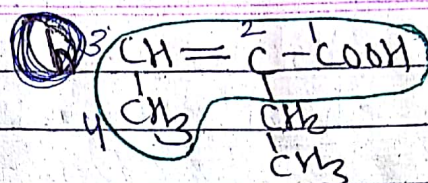
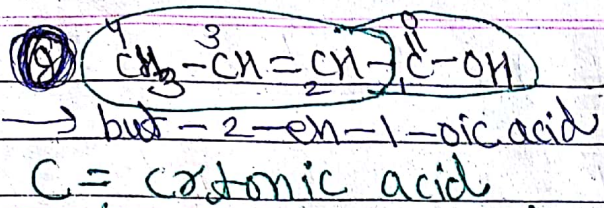
\rightarrow 2 methyl Hept-4,5-dione

\rightarrow but-1-en-3-one

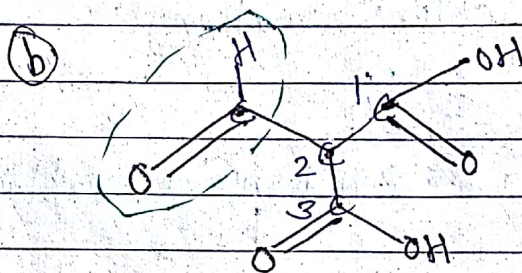


\rightarrow But-3-en-2-one

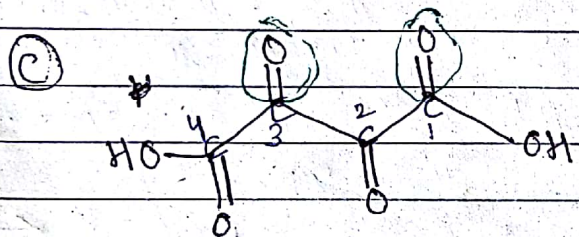
\rightarrow 2-methyl Hex-3,4-dione



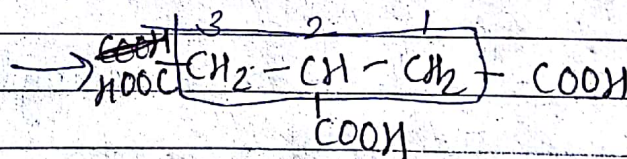
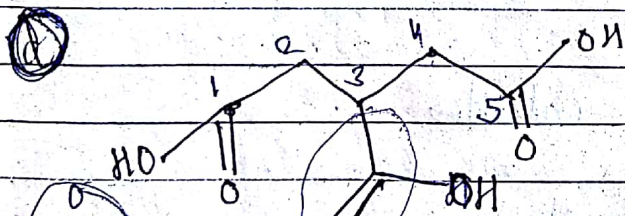
→ 2-methyl prop-2-en-1-oic acid



→ 2-formyl prop-1,3-dioic acid



→ 2,3-dioxo-but-1,4-dioic acid

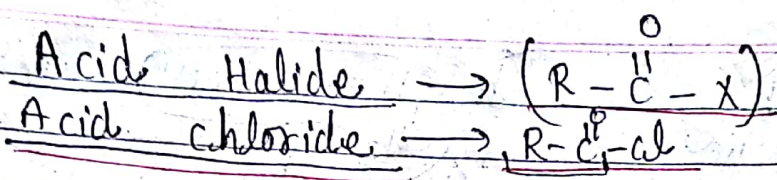


→ 3-formyl pent-1,5-dioic acid

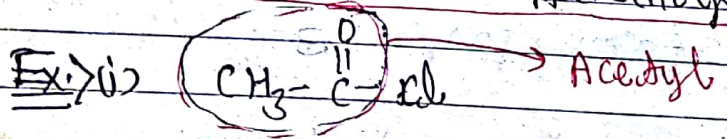
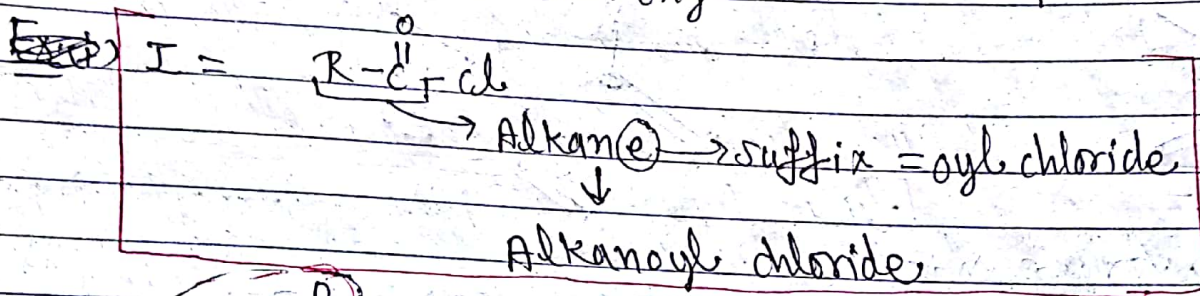
→ prop-1,2,3-tricarboxylic acid

Amine $\rightarrow R-NH_2$ $4^\circ \rightarrow$ quaternary
 Ammine $\rightarrow (NH_3)$ complex comp. h^+ $[Co(NH_3)_6]^{+3}$

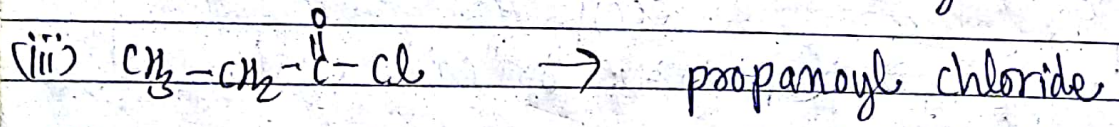
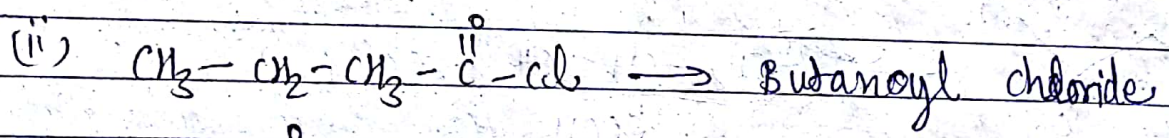
(X)



Suffix = oyl halide or oyl chloride
 Prefix = Carbonyl chloride
 = chloro carbonyl



E = Ethanoyl chloride
 C = Acetyl chloride

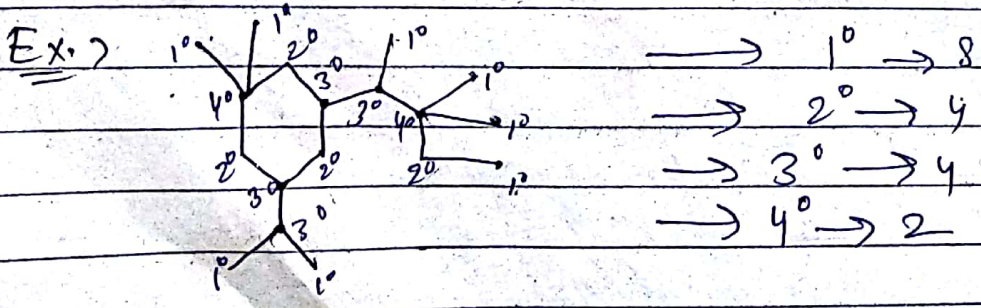


(XI)

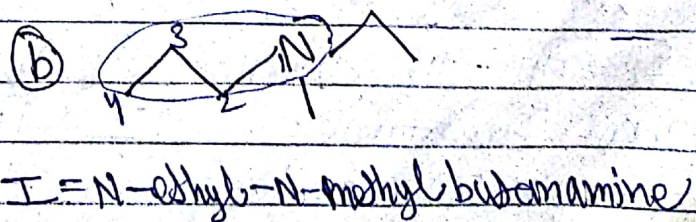
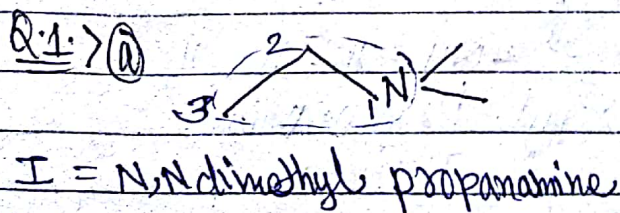
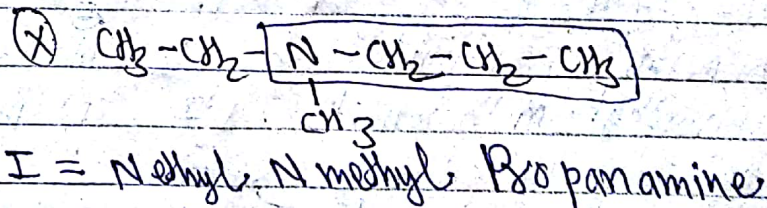
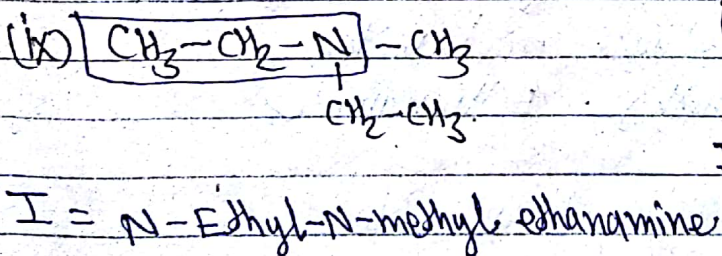
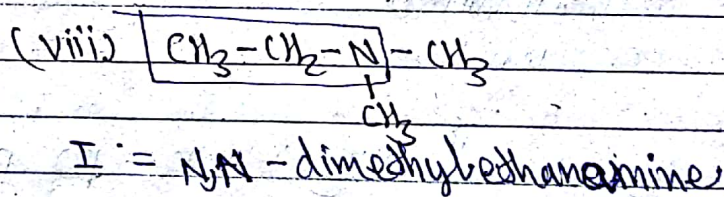
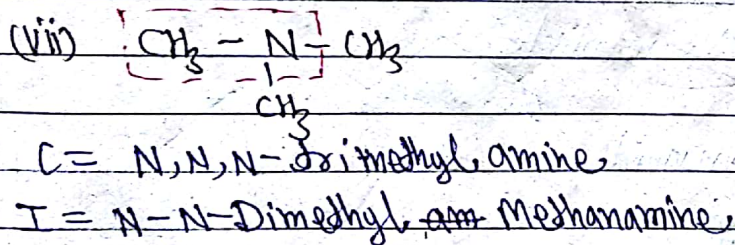
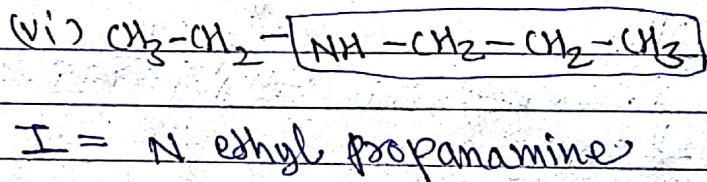
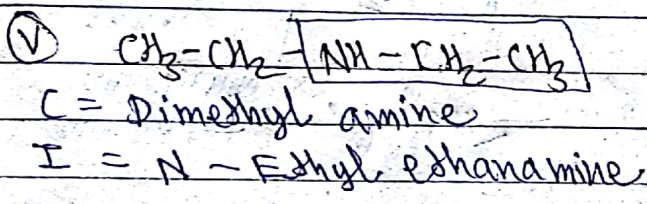
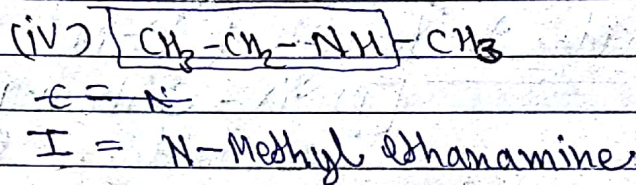
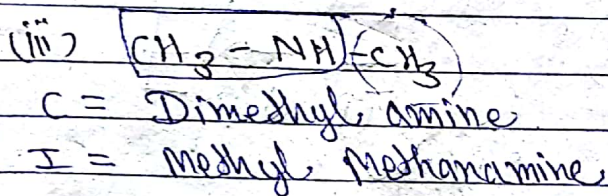
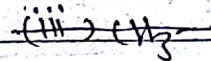
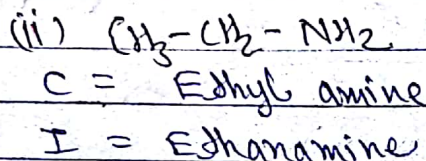
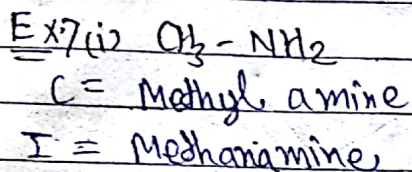
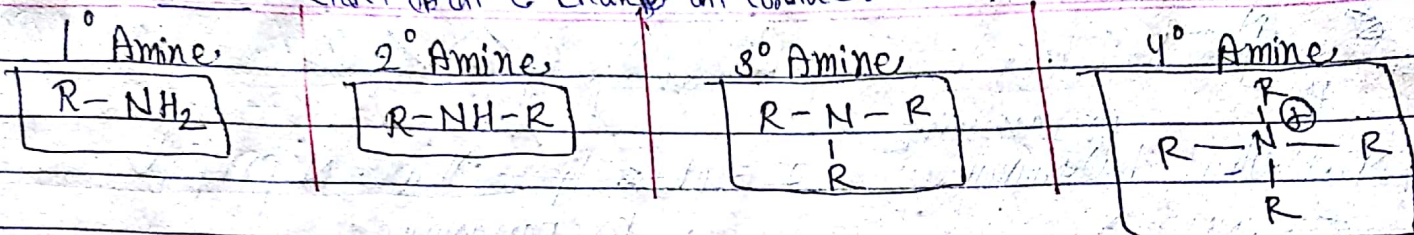
Amines :-

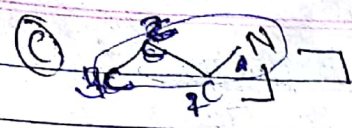
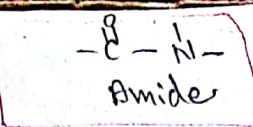
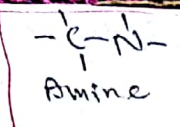
Suffix - Amine
 Prefix - Amino

$1^\circ, 2^\circ, 3^\circ, 4^\circ$ (quaternary) carbon \rightarrow

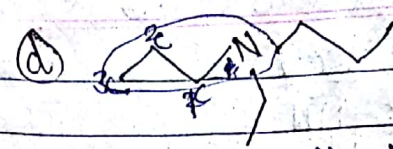


Note: - If diff C change is considered

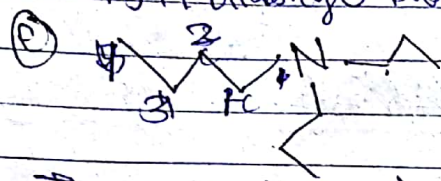




I = N,N diethyl propanamine



I = N,N diethyl N propyl propanamine



I = N,N dipropyl butanamine

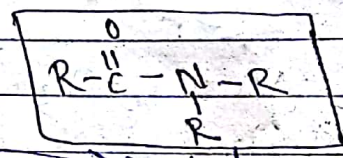
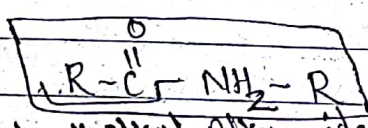
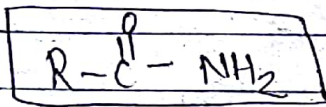
(XIII)

Acid Amide :- $(\text{R}-\overset{\text{O}}{\parallel}{\text{C}}-\text{NH}_2)$ / $[\overset{\text{O}}{\parallel}{\text{C}}-\text{N}]$
 [prefix = Carbo + alkyl] (Amides)

1° Amide

2° Amide

3° Amide



N,N dialkyl Alkylamide

Ex. (i) $\text{H}-\overset{\text{O}}{\parallel}{\text{C}}-\text{NH}_2$
 I = Methanamide
 C = Formamide

(ii) $\text{CH}_3-\overset{\text{O}}{\parallel}{\text{C}}-\text{NH}_2$
 C = Acetamide
 I = Ethanamide

(iii) $\text{CH}_3-\text{CH}_2-\overset{\text{O}}{\parallel}{\text{C}}-\text{NH}_2 \rightarrow$ I = propanamide

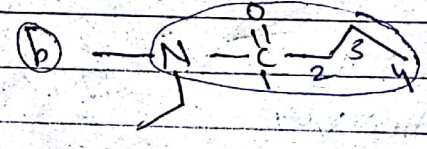
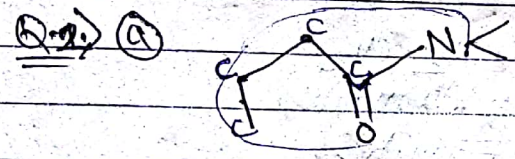
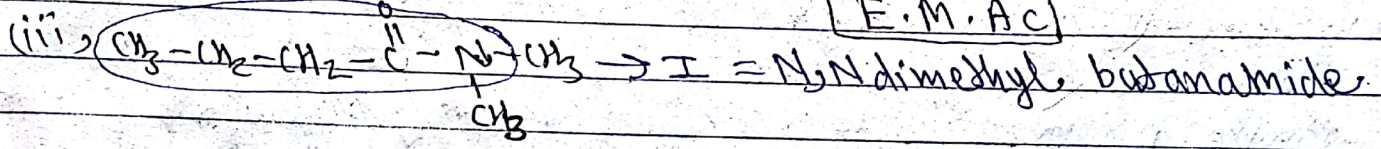
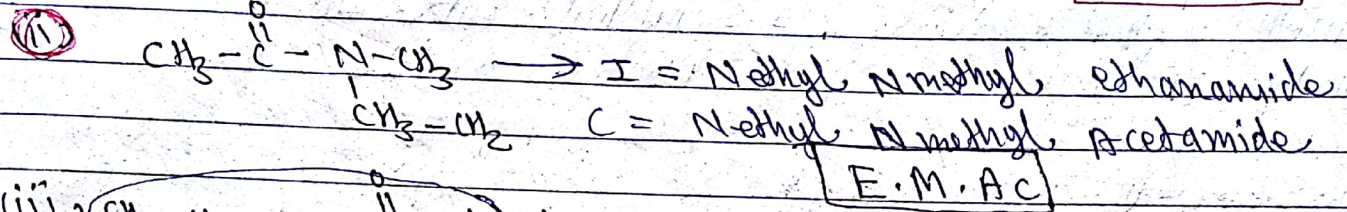
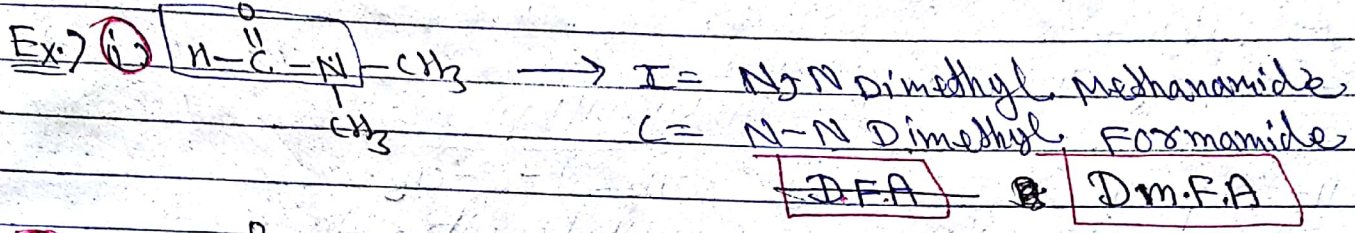
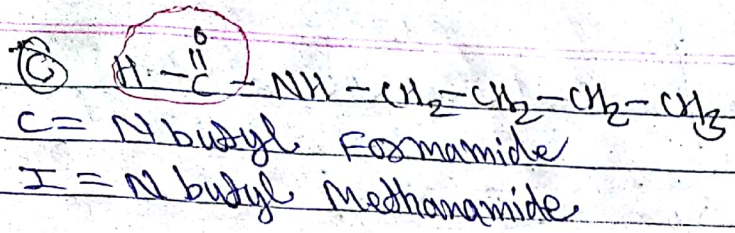
(iv) $\text{H}-\overset{\text{O}}{\parallel}{\text{C}}-\text{NH}-\text{CH}_3$
 C = N methyl Formamide
 I = N methyl Methanamide

(v) $\text{CH}_3-\overset{\text{O}}{\parallel}{\text{C}}-\text{NH}-\text{CH}_3$
 C = N methyl Acetamide
 I = N methyl Ethanamide

(vi) $\text{CH}_3-\overset{\text{O}}{\parallel}{\text{C}}-\text{NH}-\text{CH}_2-\text{CH}_3$
 C = N ethyl Acetamide, I = N ethyl ethanamide

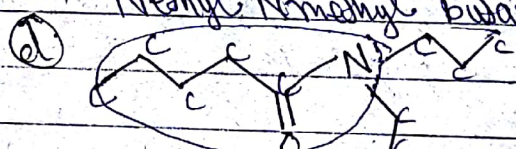
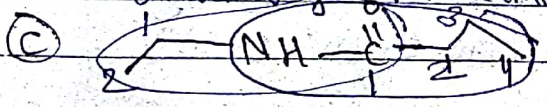
(a) $\text{CH}_3-\overset{\text{O}}{\parallel}{\text{C}}-\text{NH}-\text{CH}_2-\text{CH}_2-\text{CH}_3$
 C = N propyl Acetamide
 I = N propyl Ethanamide

(b) $\text{CH}_3-\text{CH}_2-\overset{\text{O}}{\parallel}{\text{C}}-\text{NH}-\text{CH}_2-\text{CH}_2-\text{CH}_3$
 I = N propyl propanamide



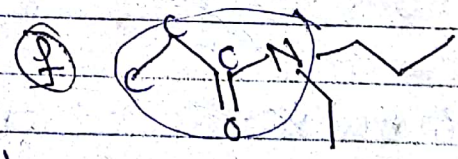
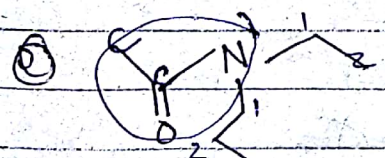
I = N,N dimethyl butanamide

I = N ethyl N methyl butanamide



I = N ethyl butanamide

I = N ethyl N propyl pentanamide



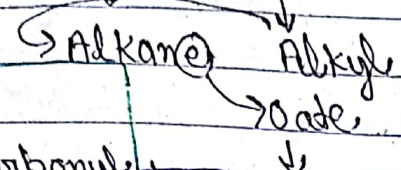
I = N ethyl N propyl ethanamide

I = N ethyl N propyl propanamide

EAA → Ethyl Aceto Acetate
 AAE → Ethyl f. Aceto Acetate Ester

XIII

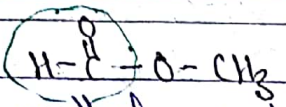
Ester :- $(R-\overset{\overset{O}{\parallel}}{C}-O-R)$ / $R-COOR$ / RCO_2R



Suffix = oate
 Prefix = Alkoxy Carbonyl

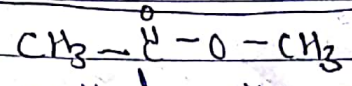
I = Alkyl Alkanoate

Ex: (i)



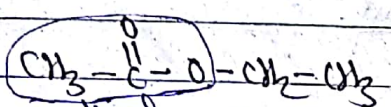
I = methyl methanoate
 C = methyl Formate

(ii)



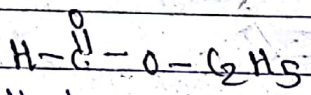
I = methyl ethanoate
 C = Methyl Acetate

(iii)



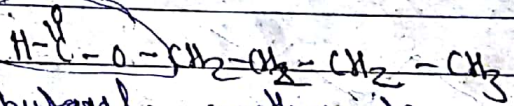
I = ethyl ethanoate
 C = ethyl Acetate

Q.1 (a)



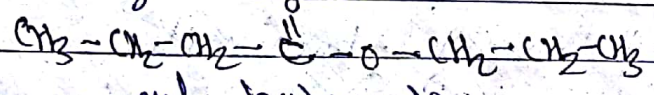
I = ethyl methanoate
 C = ethyl Formate

(b)



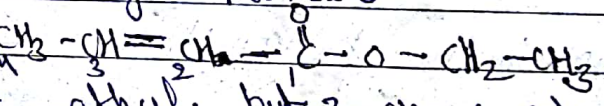
I = butyl methanoate
 C = butyl Formate

(c)



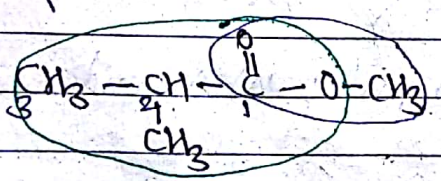
I = propyl butanoate

(d)



I = ethyl but-2-en-1-oate

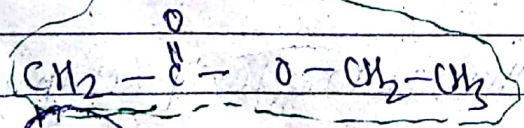
(e)



I = Methyl (2-methyl propanoate)

Q.2 (a)

Ethyl Aceto Acetate



EAA

