



Handwritten Notes
On
Optical Isomerism

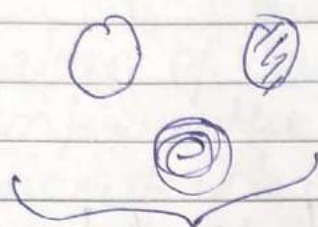


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Optical Isomerism

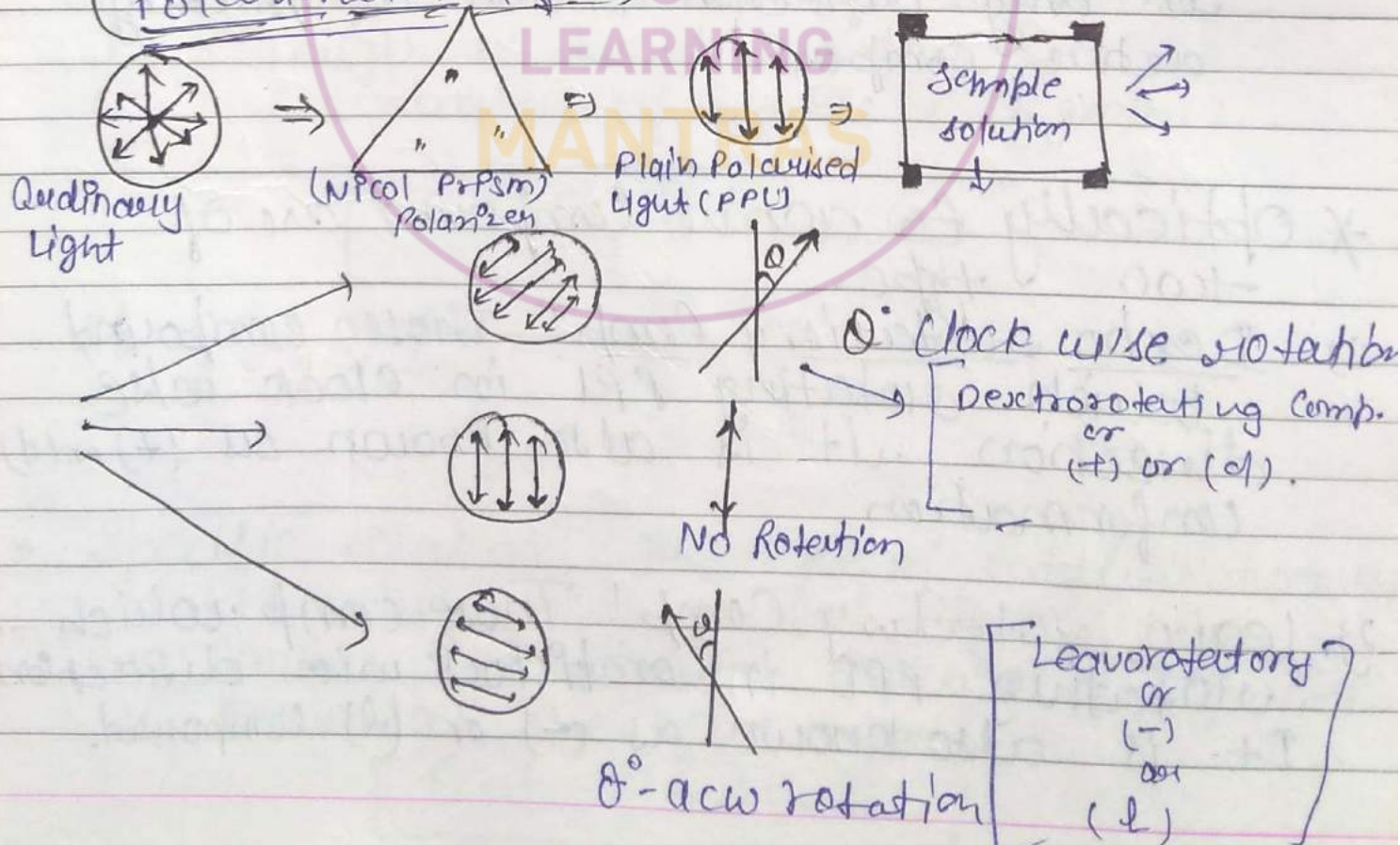


Optical Isomerism

M.f \therefore Same
 Structural formula \therefore Same
 Behaviour toward PPL \therefore Different

The compound having same molecular formula, same structural formula but diff. behaviour toward plain polarised light are known as optical and the phenomenon as optical isomerism

(Polarimeter device)



* ON the basis of behavior towards P.P.L
Compound are of two type.

- 1) optically inactive compound.
- 2) optically active compound.

1) Optically Inactive Compound!

Those comp. which not rotated PPL in any direction known as optically inactive.

2) Optically Active Compound!

Those compound which are rotating PPL in any direction known as optically active compound.

* Optically active compound are of two type.

1) Dextro rotatory Comp! Those compound which rotating PPL in clock wise direction. it is also known as (+) or (d) conformation.

2) Leavo rotatory Comp! Those comp. which rotating PPL in anticlock wise direction. It is also known as (-) or (l) compound.

* Factors affecting angle of Rotation in PPL

1. Nature of Solvent. (should be inactive)
2. Nature of Compound.
3. Temperature
4. Concentration of Sample Solution. ($\alpha \propto \text{Concentration}$)
5. Length of Sample Tube. ($\alpha \propto \text{length of Sample tube}$)

$$\alpha \propto c \quad \text{--- (i)}$$

$$\alpha \propto l \quad \text{--- (ii)}$$

$$\alpha \propto c.l$$

$$\boxed{\alpha = [\alpha] \cdot c.l} \quad \text{--- (iii)}$$

$[\alpha]$ \Rightarrow A constant k/a specific rotation.

l \Rightarrow length of Sample (in dm).

c \Rightarrow Concentration of Sample (in gm/ml)

$$\boxed{\alpha = \frac{\alpha}{c.l}} \quad \text{--- (iv)}$$

If $c = 1 \text{ gm/ml}$
 $l = 1 \text{ dm}$

then $\alpha = \alpha$

* Specific Rotation: The angle of rotation observed in PPL when it is passed through 1dm long sample tube having solution of 1 gm/ml is known as specific rotation.

$\boxed{\alpha \text{ at fixed value}}$

A/b \rightarrow $\alpha = 2$
 $\text{Path} = 6$

$\left(\frac{c - g/m}{1 - d/m} \right)$

$\frac{200 \text{ mm}}{200 \text{ mm}} \rightarrow 20 \text{ dm}$

* Specific rotation for a compound is always constant.

Ques: 6gm of 2-Butanol is dissolved in 40ml to make 40ml solution. and when it is placed in a sample tube of 200mm long. then angle of rotation observed is -4.05° . then what will the specific rotation of 2-Butanol.

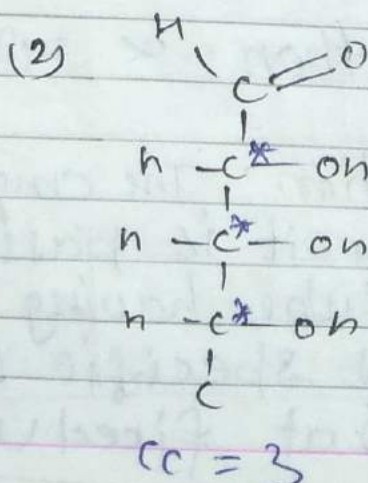
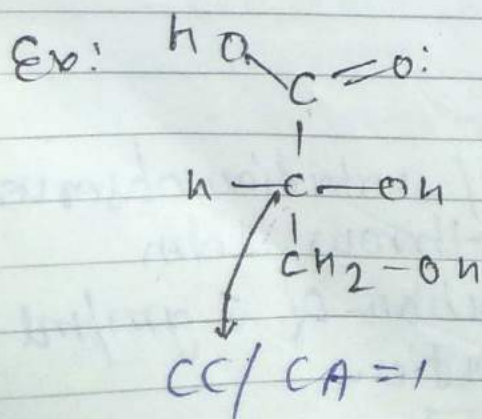
Ans $\alpha = \frac{-4.05}{\frac{6}{40} \times 2} = -13.5^\circ$

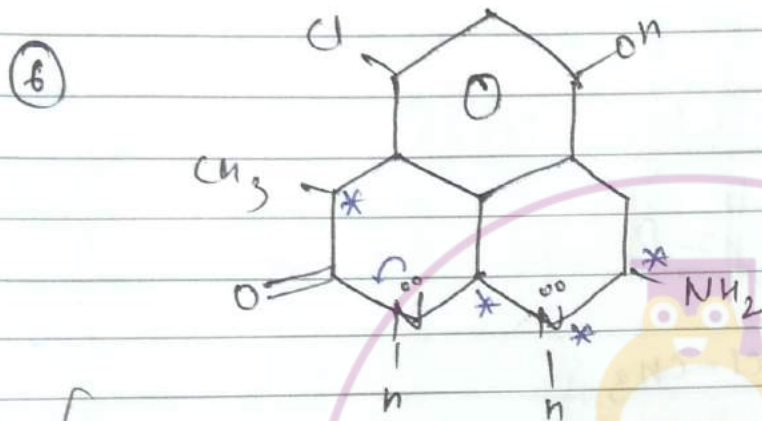
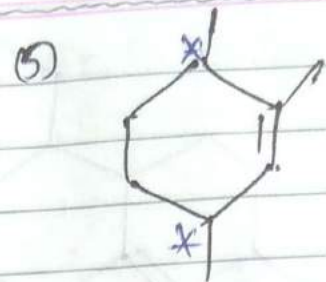
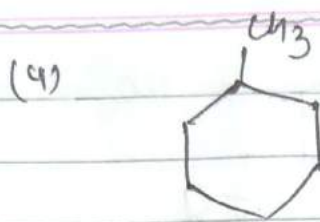
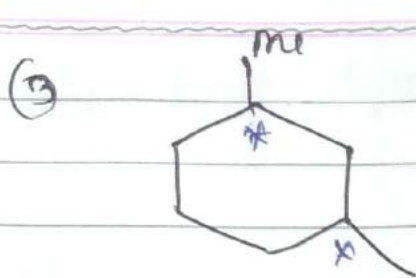
* Basic Information for Optical Isomerism

1. Chiral atom / Chiral centre :

atom with four different gp. is known as Chiral Atom / Chiral centre.
 Condition:

1. four different gp.
2. sp^3 hybridisation.
3. I.P also count in atom

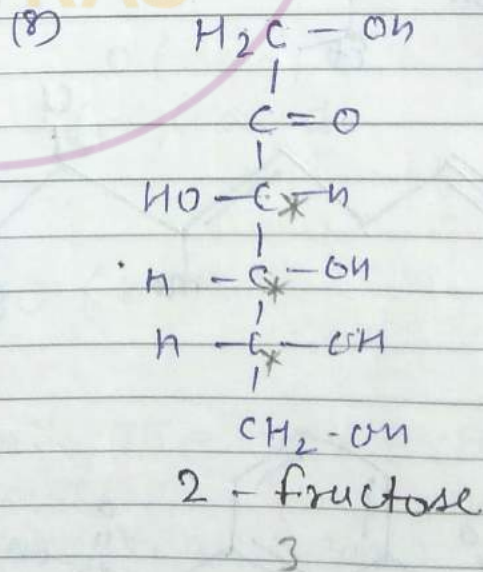
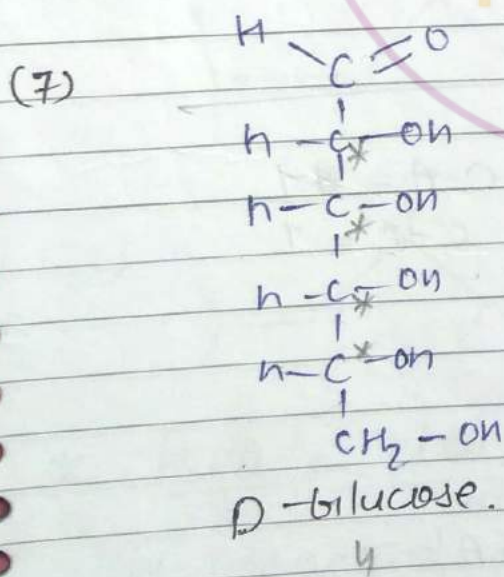




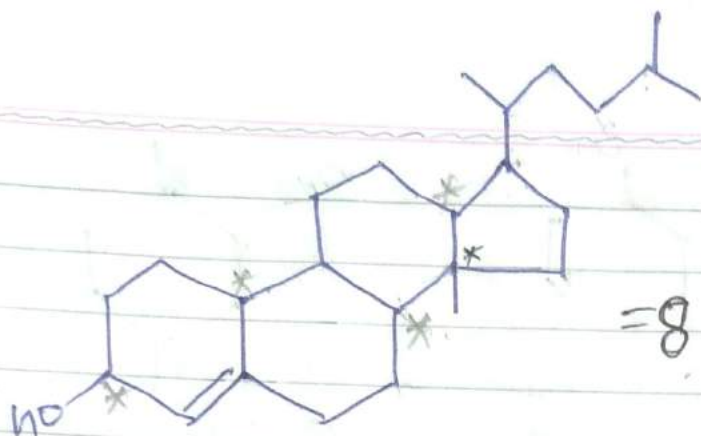
Chiral atom = 4
Chiral Carbon = 3
= Chiral Centre = 4

Note: The lone pair participate in Resonance does not Count in chiral centre.

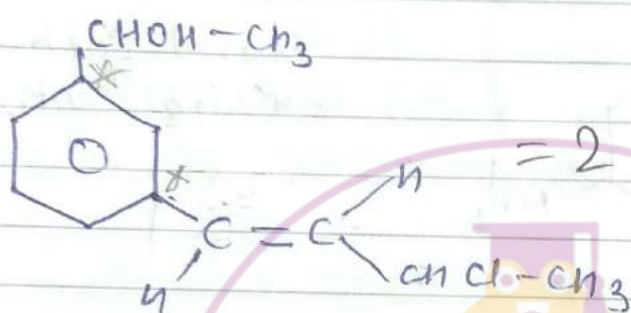
Que: decide chiral atom / chiral Carbon in following structure.



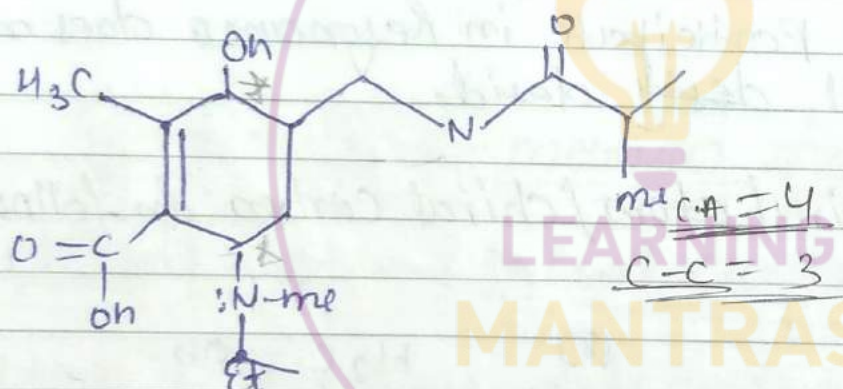
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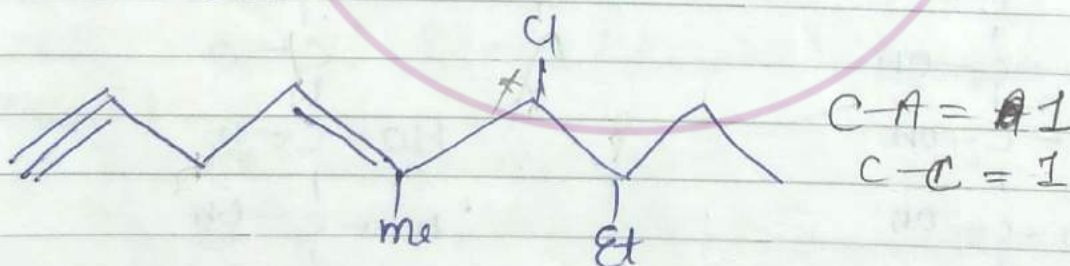
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(11)

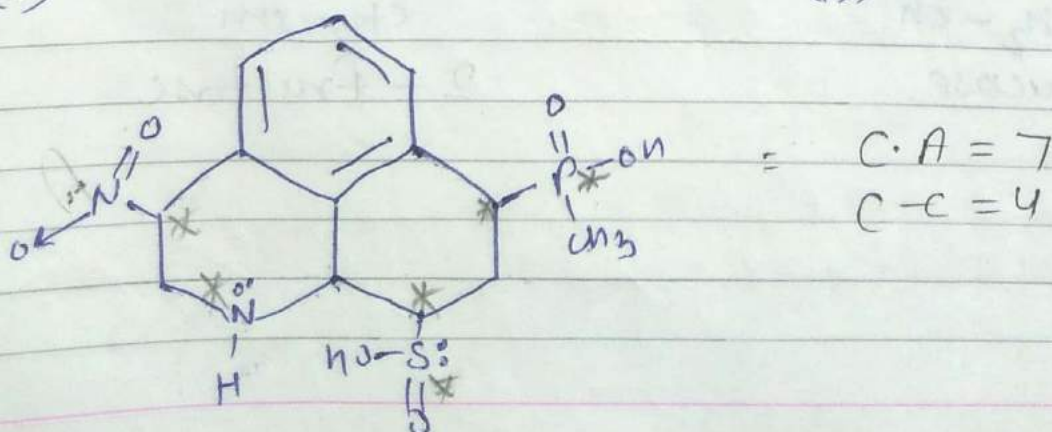


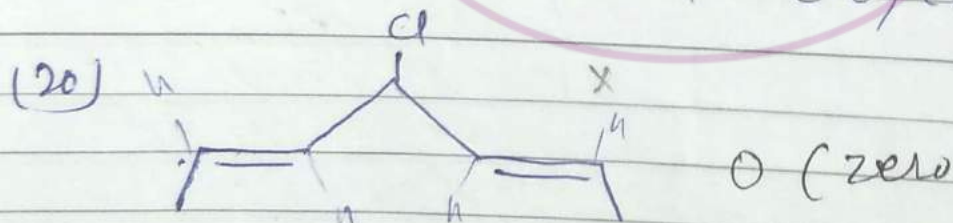
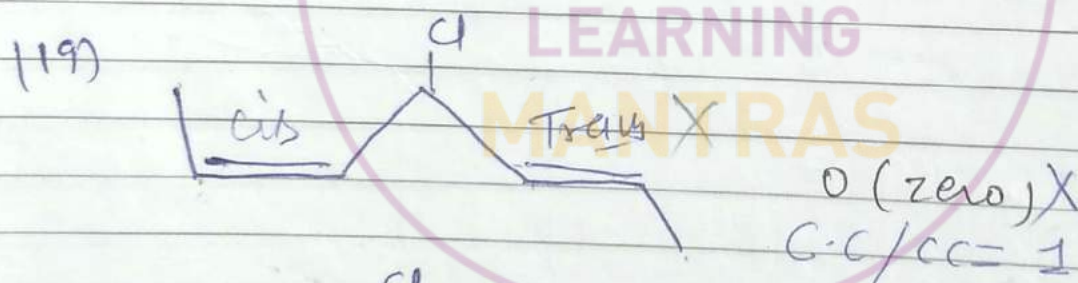
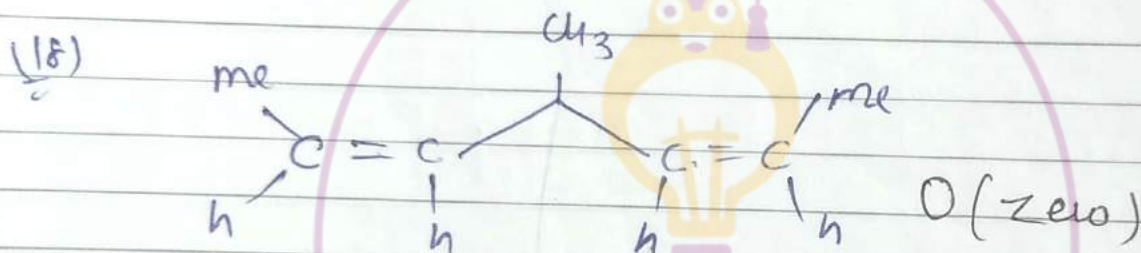
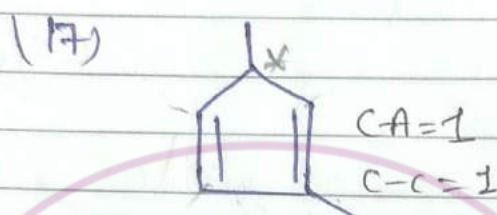
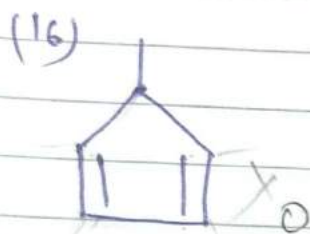
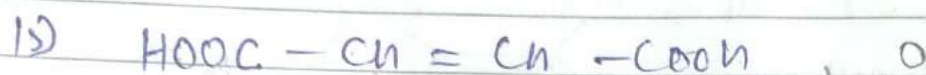
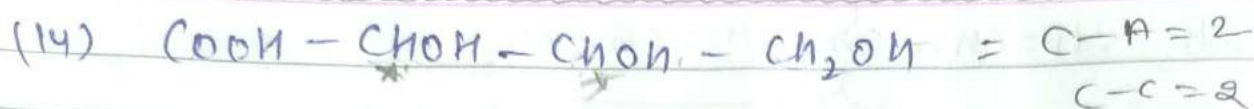
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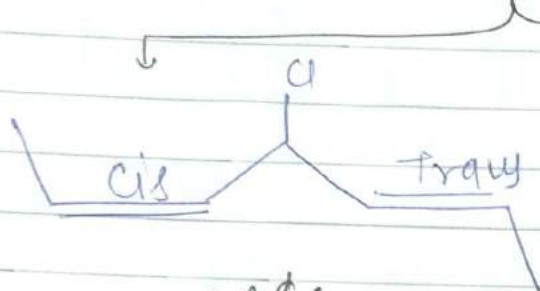
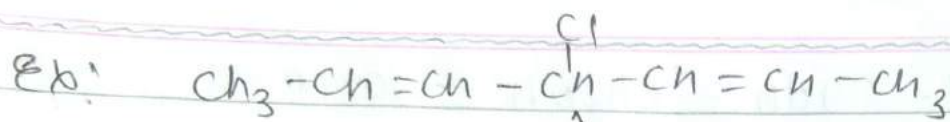
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(14)

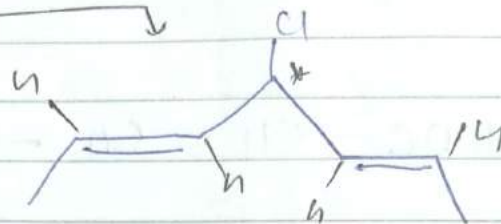




* Pseudo chiral atom :- The carbon of which chirality depends on stereochemistry of attached gp. is known as pseudo chiral atom.



$\text{CA/CC} = 1$



$\text{CC/CA} = 0$

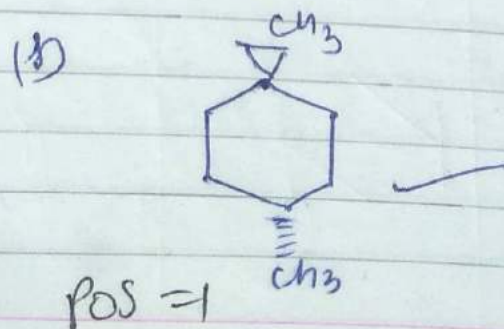
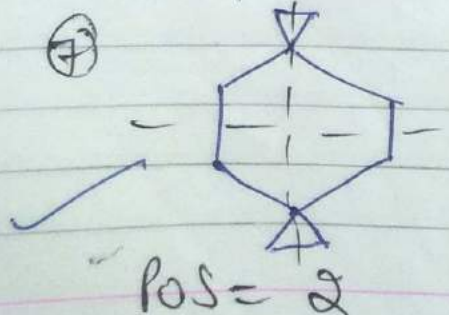
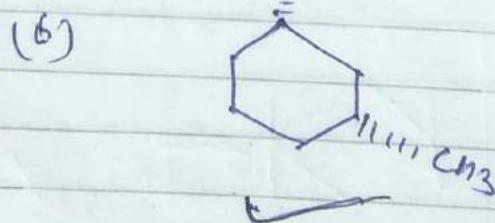
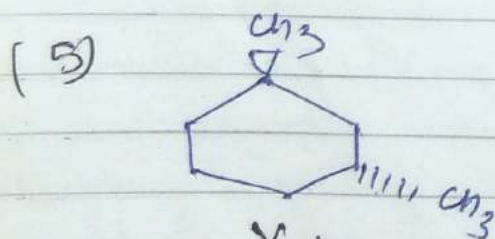
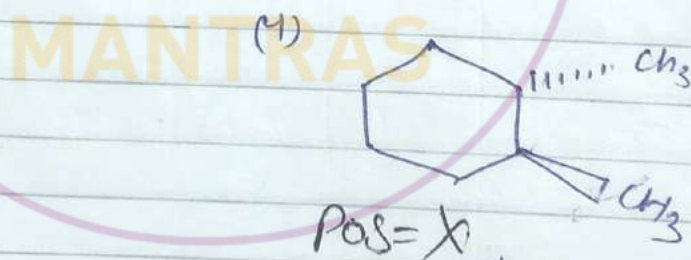
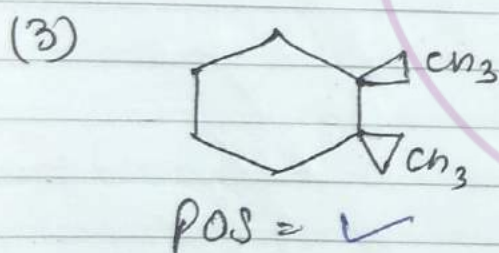
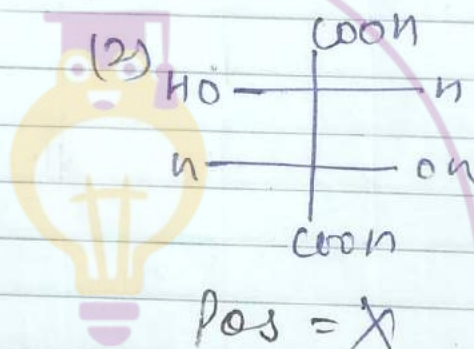
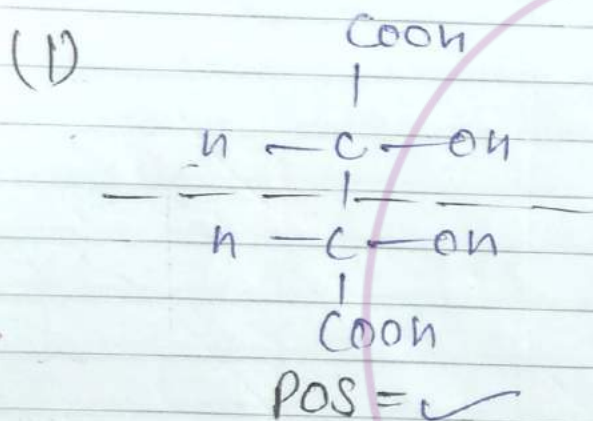


LEARNING
MANTRAS

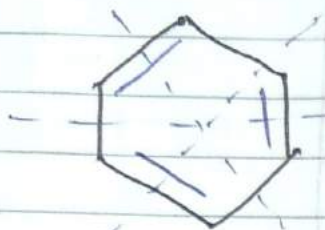
* Elements of Symmetry =

	Symbol	Sh-
Plane of Symmetry	(σ)	→ POS
Centre of Symmetry	(CP)	→ COS
Alternating Axis of Symmetry	(S_n)	→ AAOS
Axis of Symmetry	(C_n)	→ AOS

* Plane of Symmetry: (POS) → (σ) :



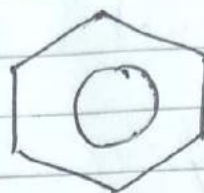
(9)



R.S of Benzene

POS $\Rightarrow 4$

(10)

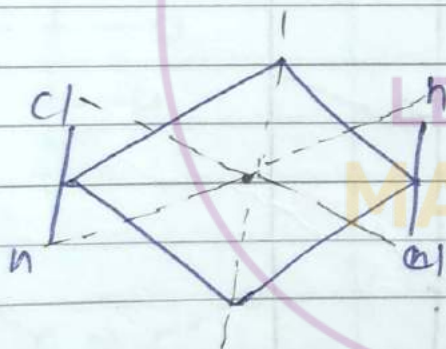


(7)

R.H of Benzene

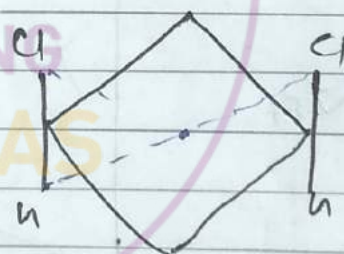
* Planar system always have slice Plane of Symmetric

* Centre of Symmetry \rightarrow (centre of Inversion)
 $(\text{C.O.S}) \rightarrow (\text{C.I})$



C.O.S =

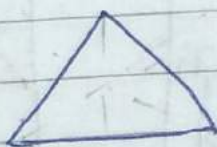
Trans-1,4-DCCB



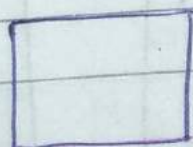
C.O.S = X

Que! Decide compounds
 compounds

1)



2)



POS

C.O.S

chiral
or


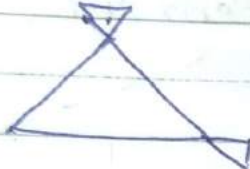
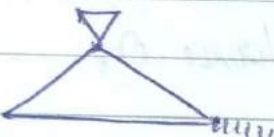
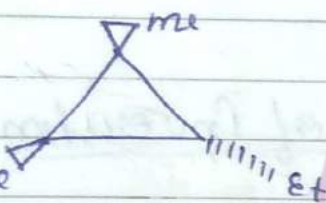
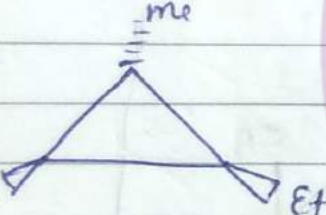
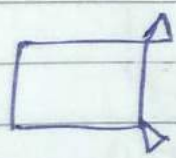
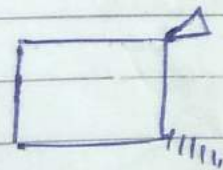
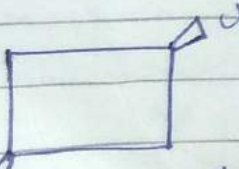
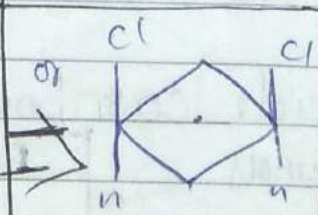
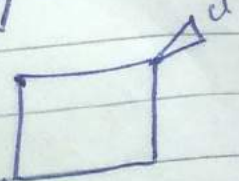
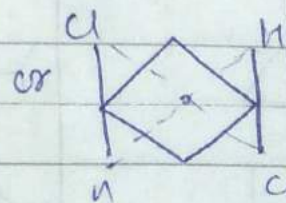
4

X

X

5

X

Compound	Pos	CoS	Chiral op
3) 	✓ 5	X	X not g
4) 	✓	X	X
5) 	X	X	✓
6) 	✓ 1	X	X
7) 	X	X	✓
8) 	✓	X	X
9) 	X	X	✓
10) 	✓ 2	X	
11) 	✓ 1		X

Chiral
o.a

	Pos	cos	
12)	✓	X	X
13)	X	X	
14)	X	X	
15)	X	X	✓
16)	✓=2	X	X
17)	✓=1	✓	X
18)	✓=2	X	X
19)	✓=6	X	X

Open chain non planar system has no COS Generally. Cis system has no cos.
sp³ hybridisation no cos. (non planar)

Chiral

	Pos	CoS	OA
20) <chem>ClC(Cl)(Cl)Cl</chem>	$\checkmark = 3$	X	X
21) <chem>ClC(Cl)(Cl)Cl</chem>	$\checkmark = 2$	X	X
22) <chem>F[C@H](Br)I</chem>	X	X	\checkmark
23) <chem>Cc1ccccc1C</chem>	$\checkmark = 2$	X	X
24) <chem>Cc1ccccc1C</chem>	$\checkmark = 2$	X	X
25) <chem>Cc1ccccc1C</chem>	$\checkmark = 3$	\checkmark	X
26) <chem>C1=CC2CCC1C2</chem>	X	X	\checkmark
27) <chem>C1=CC2CCC1C2</chem>	$\checkmark = 1$	X	X
28) <chem>C1=CC2CCC1C2</chem>	X	X	\checkmark

Cis \Rightarrow Neither COS

Trans - Both

Chiral
o.a.

Pos

Cos

29)



✓₂

X \Rightarrow because cis

X

30)

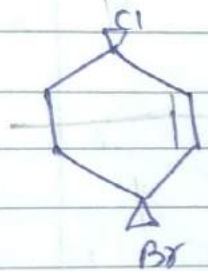


✓

✓
X

X

31)



X

X \Rightarrow

✓

32)



X

X = MTR

X

33)



✓₂

X

X

34)

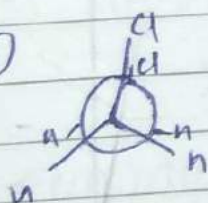


X

X

✓

35)

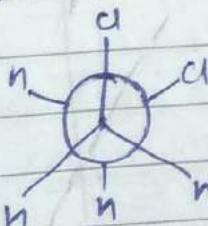


✓

X

X

36)



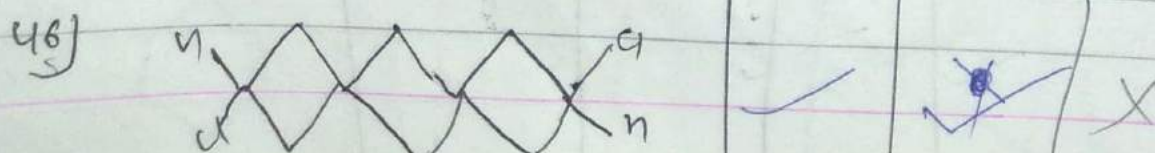
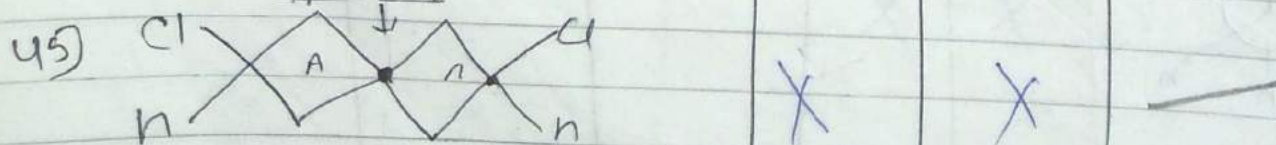
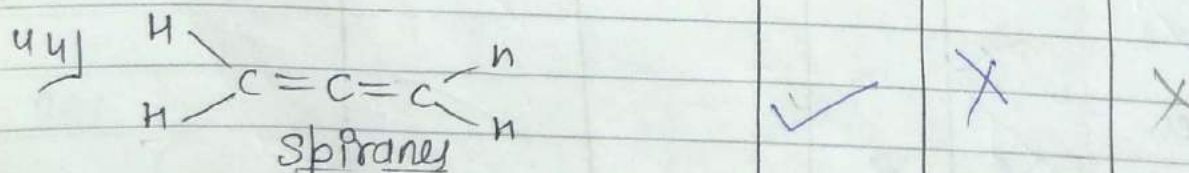
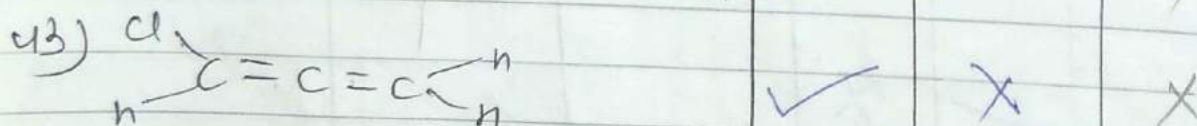
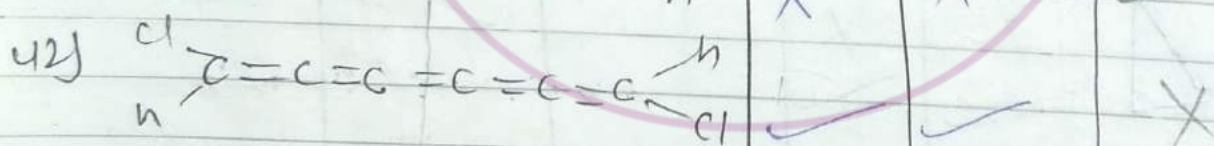
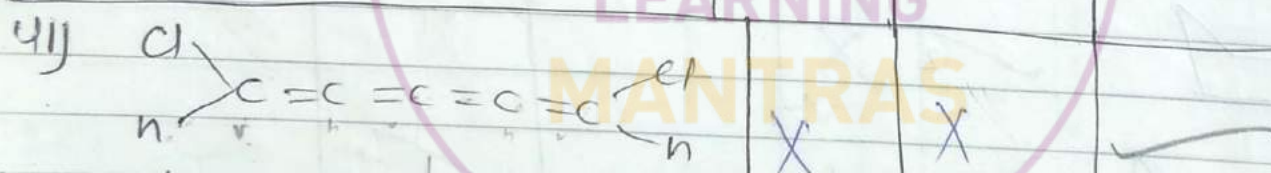
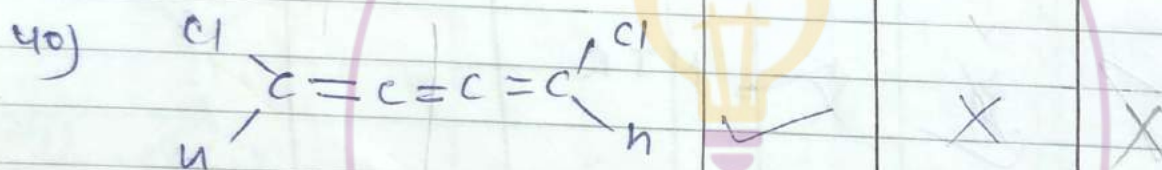
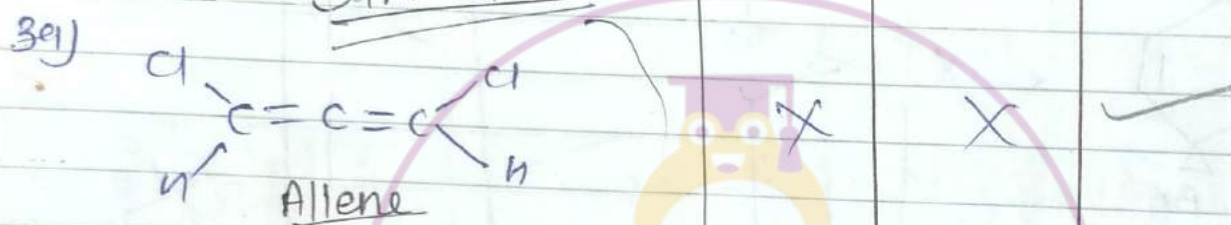
X

X

✓

	Pos	Cos	Chiral aa
37)	X	X	✓
38)	✓	✓	X

Cummulenes



		Pos	Cos.	Chiral C
47)		✓	X	X
	<u>Allylic</u>			
48)		X	X	✓
49)		✓	X	X
50)		=4 ✓	X	X
51)		=2 ✓	X	X
52)	Screw	X	X	✓
53)	Kid's tricycle	X	X	✓
54)	ceiling fan	X	X	✓
55)	Hockey stick	X	X	✓

* cummulenes / spirane / Alkylidene:

When total no. of db/cycle/
/cycle + db.
↓
odd No.
↓
Planar system
Pos (v)

When total no. of db/
cycle/cycle + db
↓
even no.

Pos x
When valency of terminal-c
are diff.

Pos = v

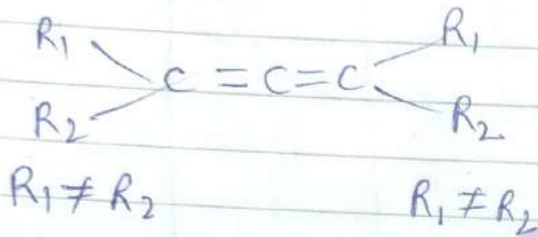
even no.

POS X

POS ✓

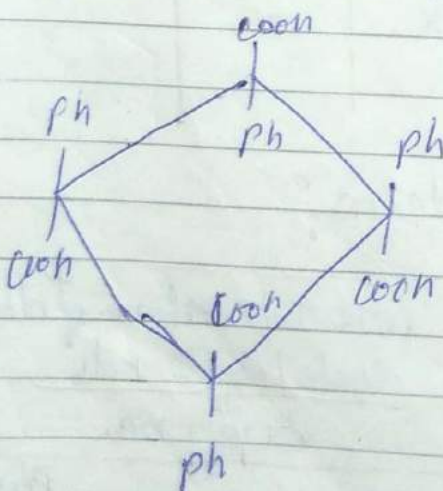
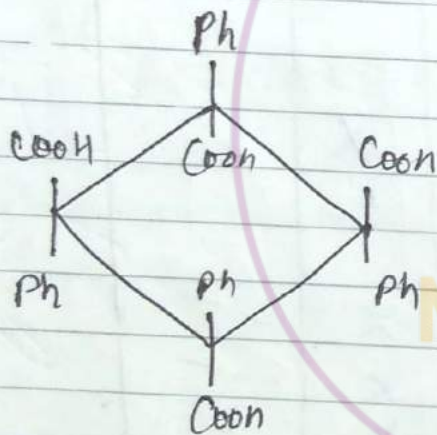
When valency of terminal-C are different

When R_1, R_2 of $\overset{\text{ant}}{\text{C}}$ are ~~both~~ same

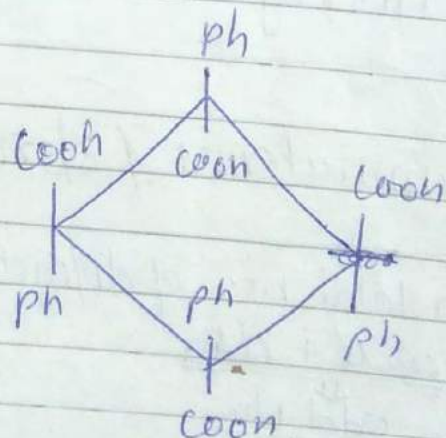


$$R_1 = R_2$$

Alternating axis of symmetry (AAOS) $\rightarrow S_n$



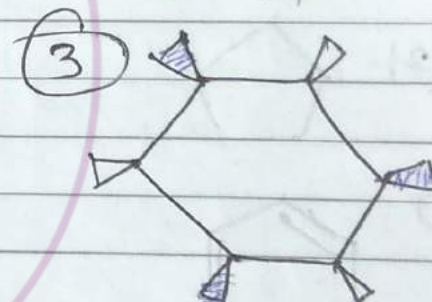
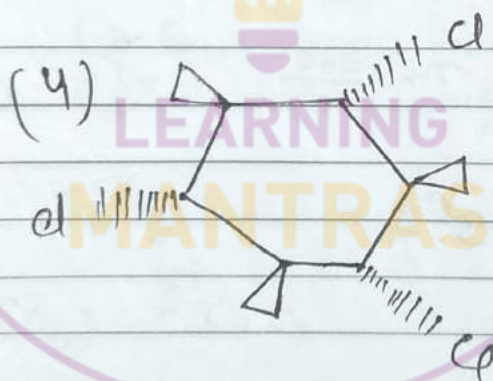
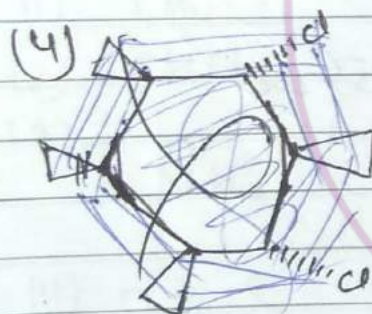
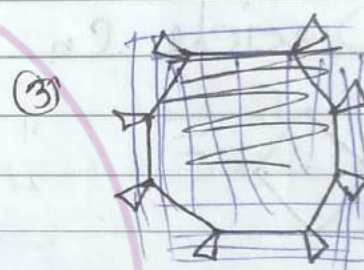
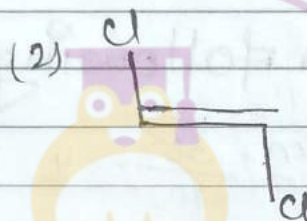
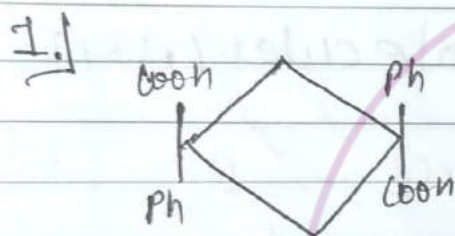
90°



- AAOS Present
- S_n Present = ($n = 4$)
- $n = \frac{360}{90} = 4$
- \rightarrow four fold AAOS (S_4)

{ Those Comp. which have either POs and/or have AAOS }

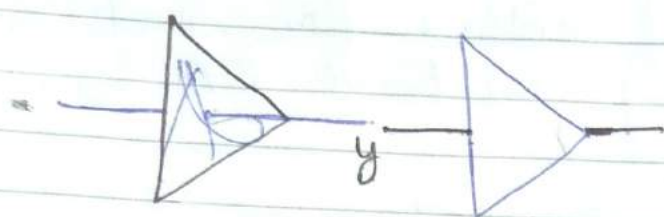
Ques: decide four fold AAOS
 \rightarrow Sn in following molecules



* Axis of Symmetry : (AOS) $\rightarrow C_n$:



- Along n Axis
- AOS = $\sqrt{\quad}$
- $C_n = ?$
- $n = \frac{360}{120} = \frac{360}{120} = 3$
- Three fold AOS - C_3



Along y -axis

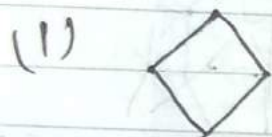
AOS = \checkmark

$$n = \frac{360}{180} = 2$$

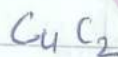
$$C_n = C_2$$

$$\text{Total } C_n = C_1, C_2, C_3$$

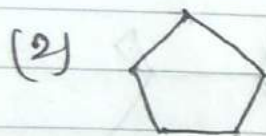
* Decide C_n in following molecules (r.m.)



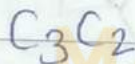
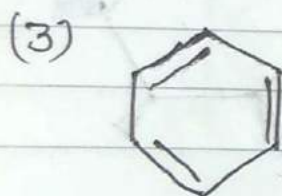
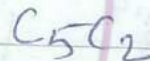
$$\frac{4}{2} \quad n = \frac{360}{90} = 4$$



$$\rightarrow n = \frac{360}{180} = 2$$

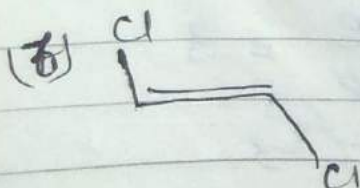
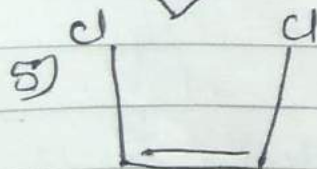
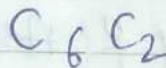
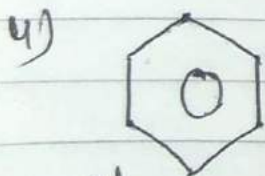


$$\rightarrow n = \frac{360}{72} = 5$$

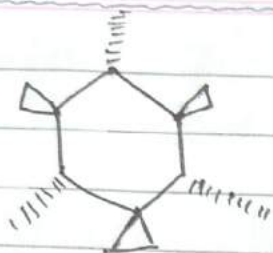


$$\frac{6}{2}$$

Res. of Benzene

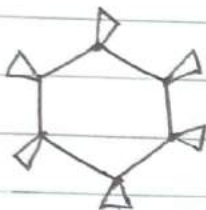


7)

 $C_3 C_2$

Linear structure
 C_∞, C_2

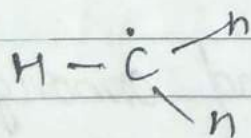
8)

 ~~$C_3 C_2$~~ C_6 \rightarrow Trigonal P. $\rightarrow C_3, C_2$ \rightarrow Tetrahedral $\div C_3, C_2$ \rightarrow Pyramidal $\div C_3$ \rightarrow V-shaped $\div C_2$ 9) $H-H$ ~~C_3~~ C_∞

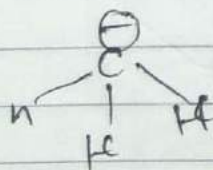
10)

 C_2 11) CH_4 $\rightarrow C_3, C_2$ 12) $CHCl_3$ $\rightarrow C_3$ 13) $\rightarrow \ddot{N}H_3$ C_3 14) $O=S(=O)_2$ (SO_3) $\rightarrow C_3 C_2$

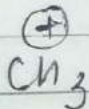
15)

 (in_3) $C_3 C_2$

16)

 (CH_3^+) C_3

17)

 $C_3 C_2$

Essential Condition of optically active

H.W! Race : 6.

(18)



C_2 trans

(19)



C_2



$C_3 C_2$

* Note: Essential condition for optically active compound which have neither POS nor AOS nor COS are optically active.

* All optically compound are known as chiral, dissymmetric compound.



Pos = X

Cos = X

AOS = X

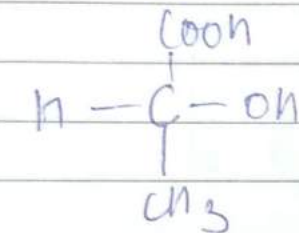
→ chiral compound /
Dissymmetric comp

Ques! Identify chiral/optically active compound among following.

Ques 1-55 Ques

Detail study for 1 chiral or 2 chiral atom compound.

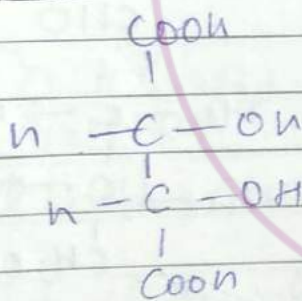
* Single Chiral Compound ^{atom.}



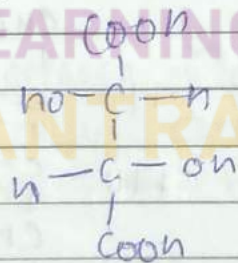
Pos = X
Neg = X
OA / chiral.

Single chiral atom compound are always optically active.

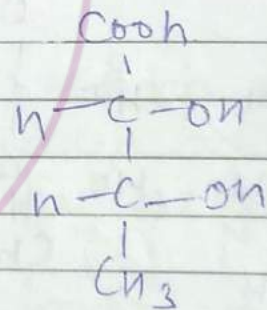
2) Two chiral compound



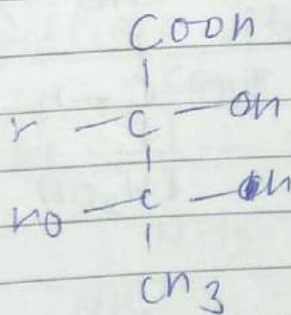
Pos = ✓
OA / chiral



Pos = X
Neg = X
OA
Chiral



⇒ 11



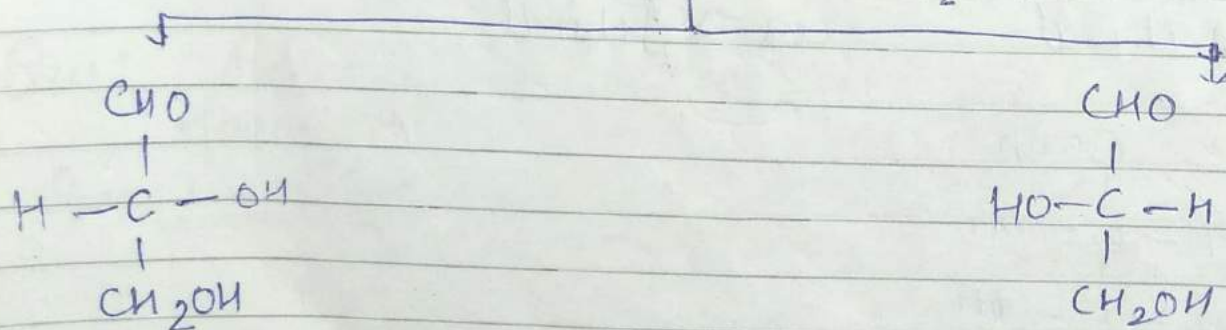
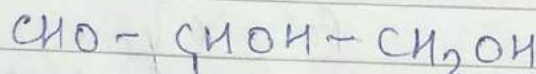
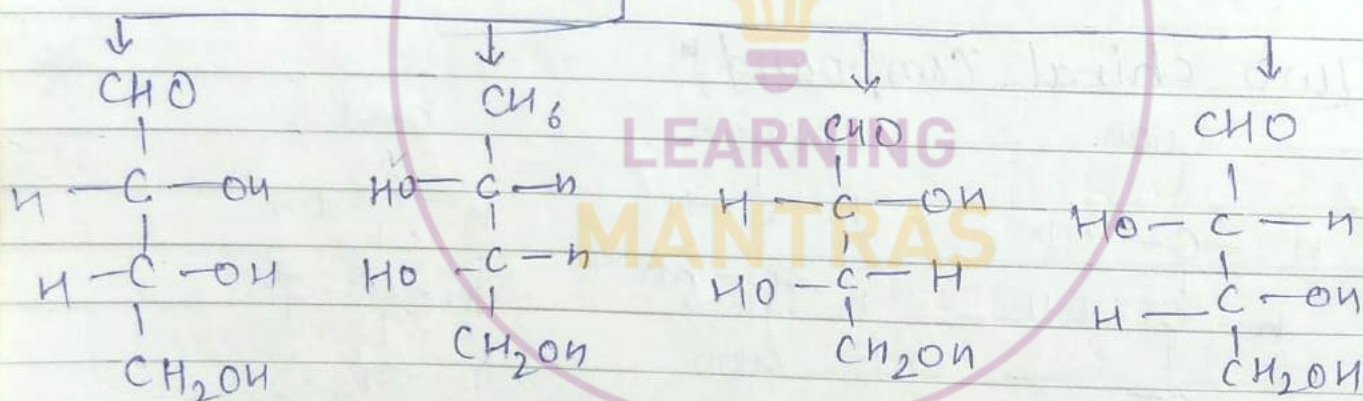
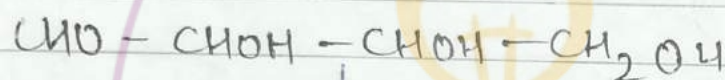
N. Very Important.

HW: 10:00:00

* Method of representation of molecules:

1. Ball stick.
2. wedge dash
- 3.
- 4.

5] Fischer projection formula:



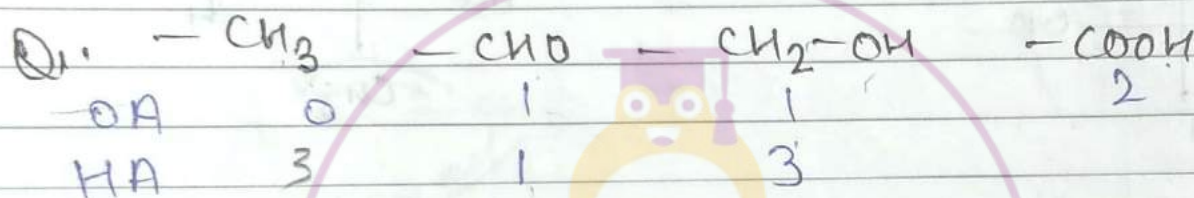
- COOH

Priority State: No. of O.A = O.A \propto O.S

→ No. of H.A = H.A \propto $\frac{1}{O.S}$

→ (+)ve charge = (+)ve charge \uparrow , O.S \uparrow

→ (-) " " = (-)ve " \uparrow O.S \downarrow

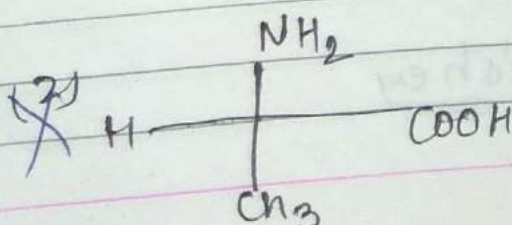
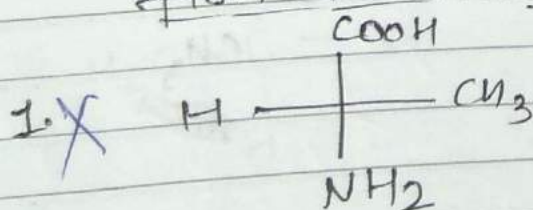


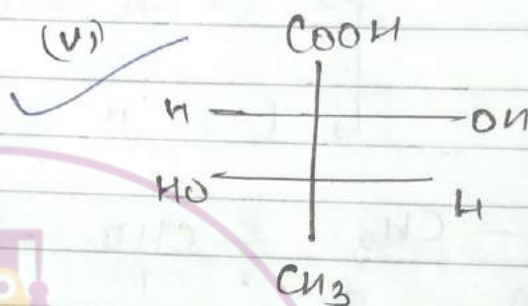
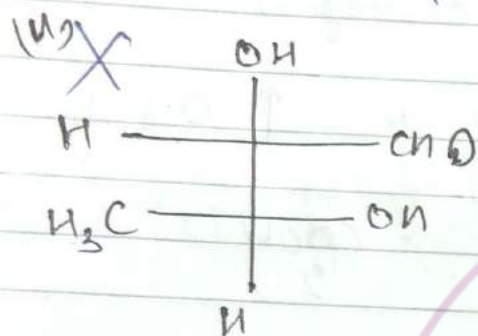
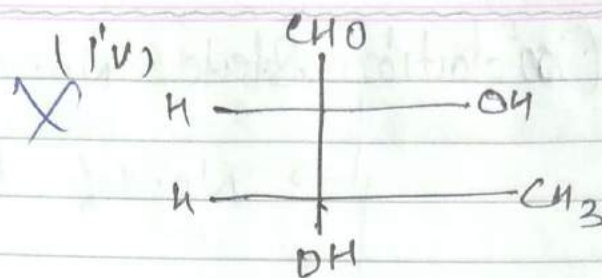
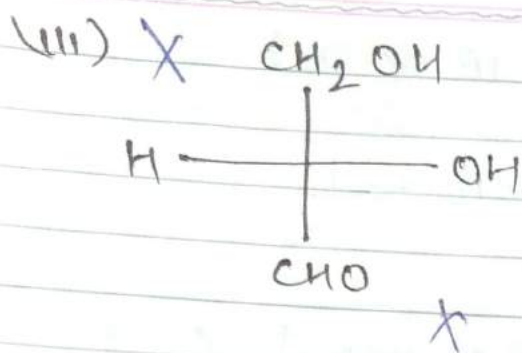
Ans: $\text{—COOH} > \text{—CHO} > \text{—CH}_2\text{OH} > \text{—CH}_3$

* In fisher structure carbon chain are written vertically and group on chiral carbon either left or right.

* In fisher projection formula more oxidised end must be at top.

Q. Which of the following structure(s) is/are fisher structure:-

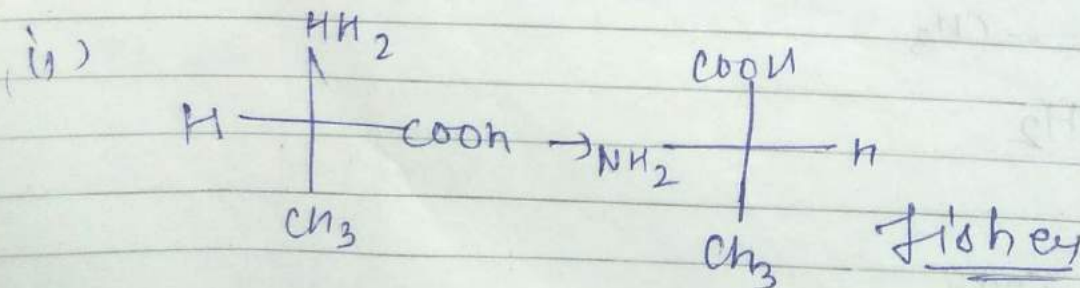
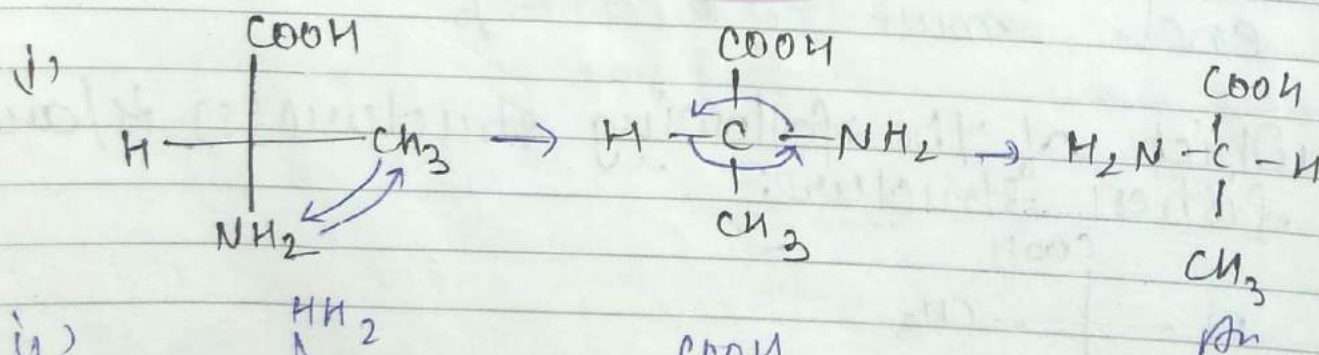


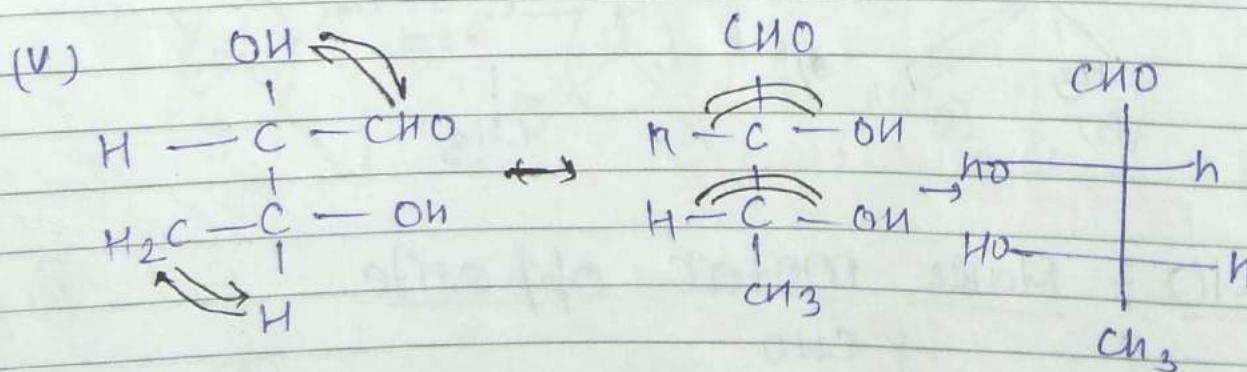
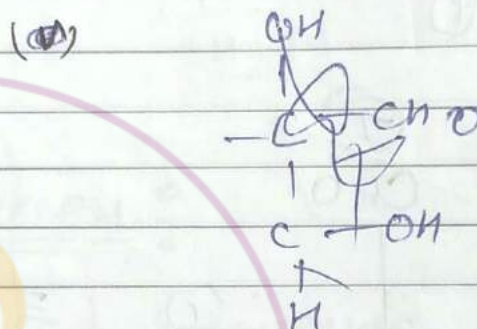
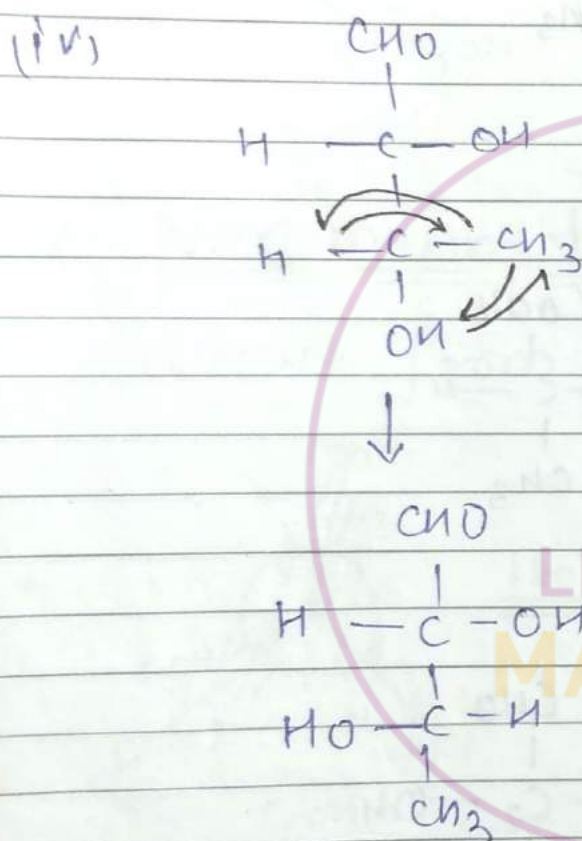
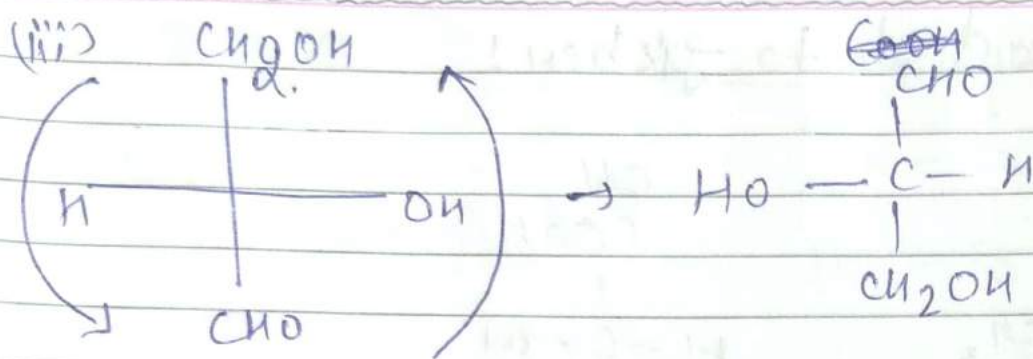


* Method to convert different structure into Fisher structure :

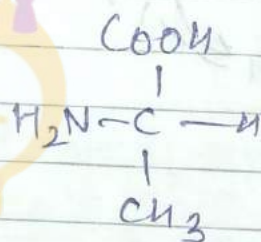
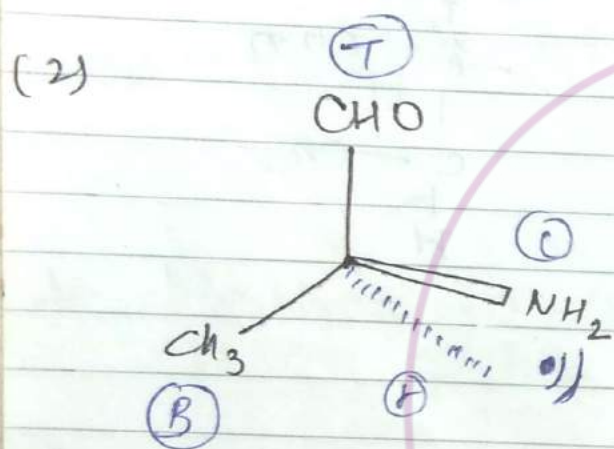
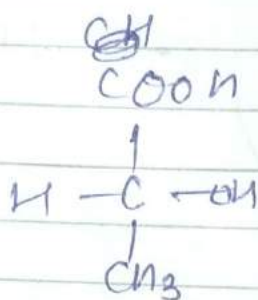
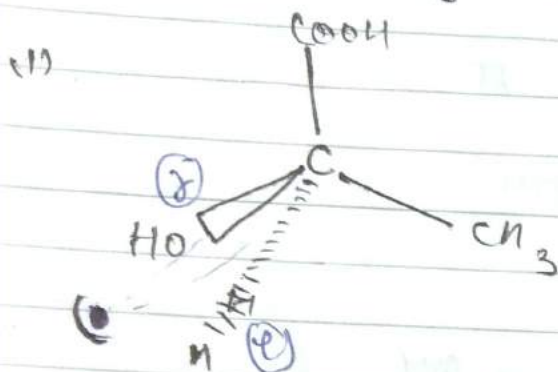
1. from 2D non fisher to fisher :

Freedom → even interchange on each c.
 → 180° Rotation

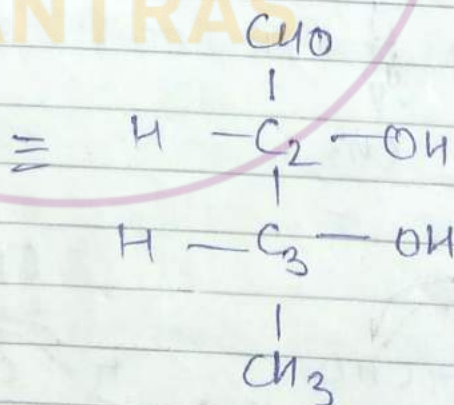
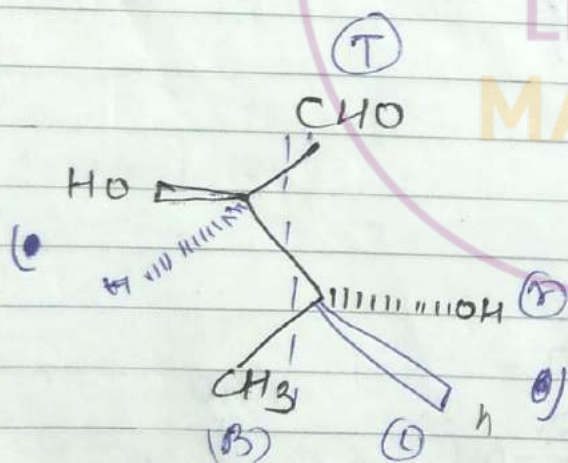




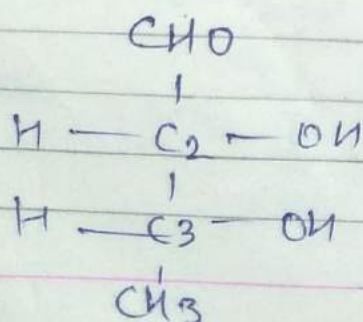
* For wedge dash to fisher!

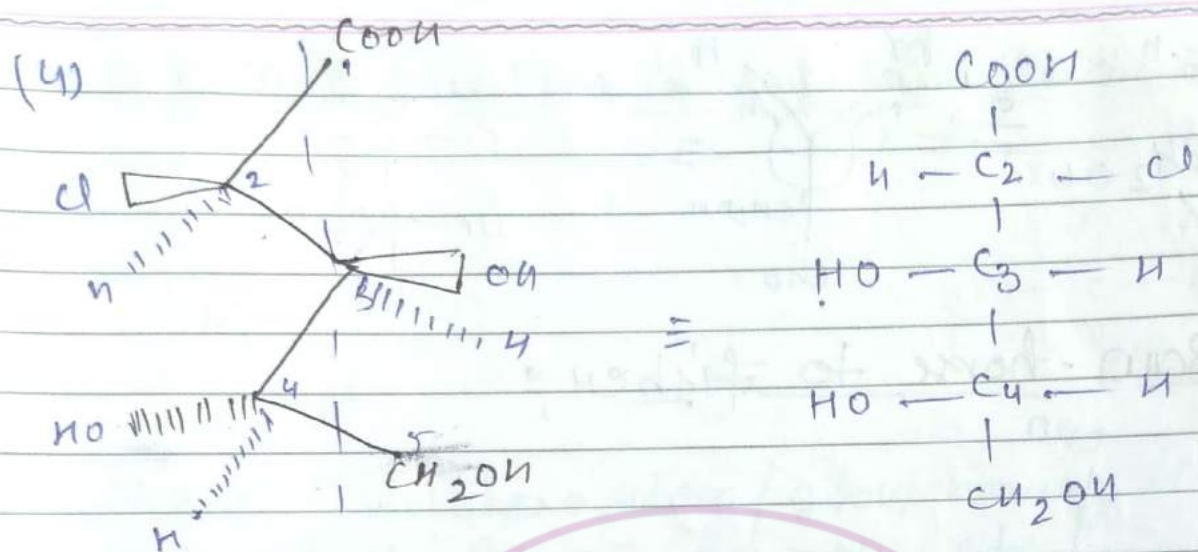


LEARNING
MANTRAS



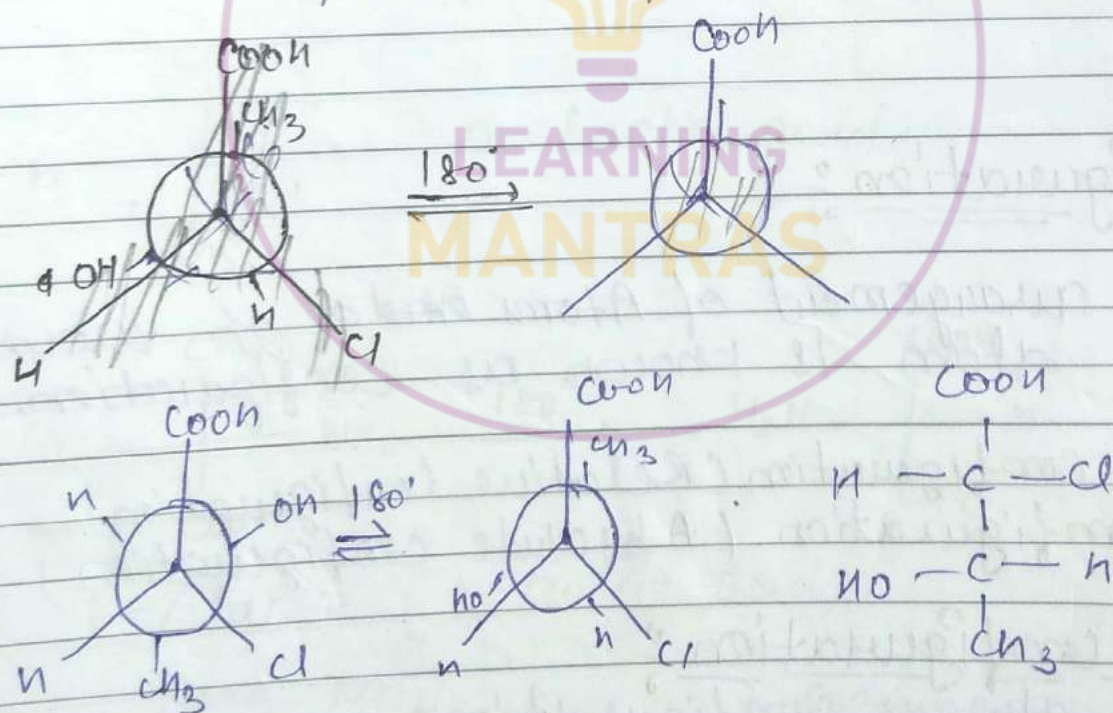
* MMO : Make motak opposite.



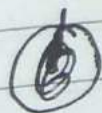


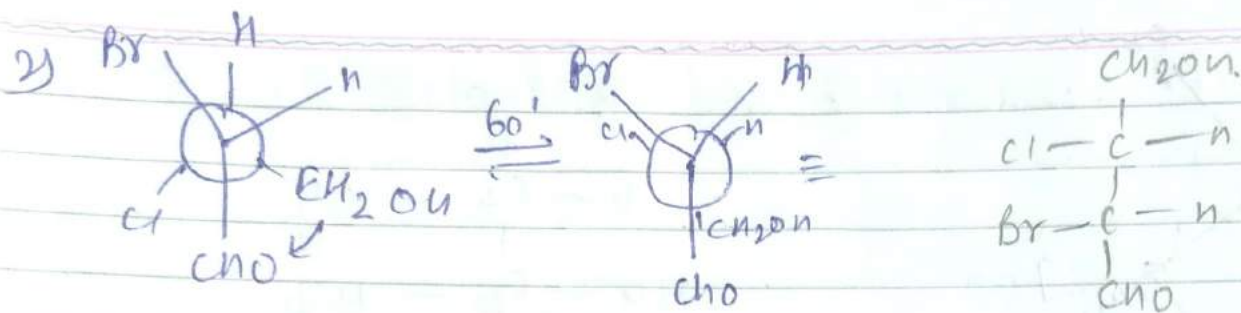
* Newmann to fisher :

Start from Eclipsed structure

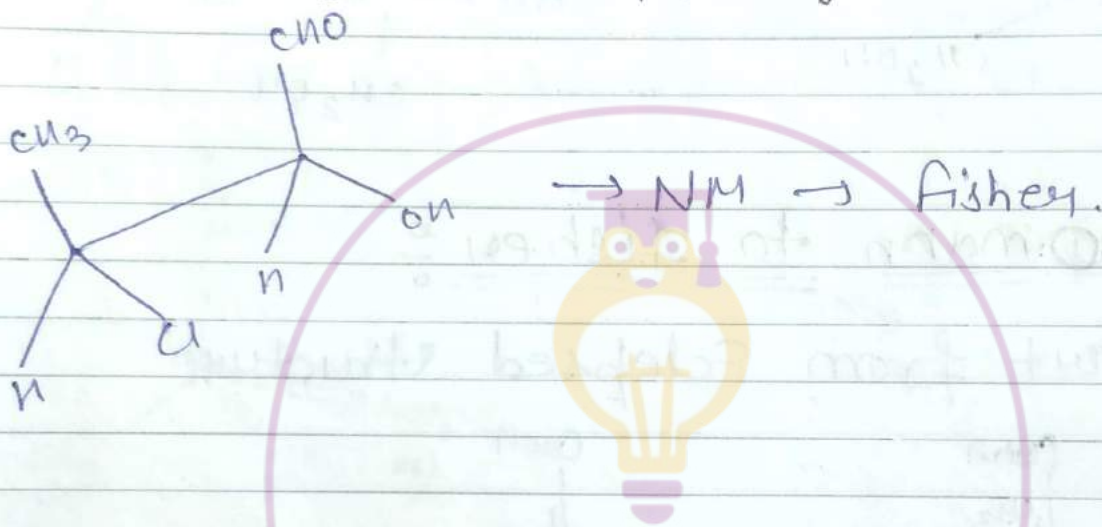


(2)





* From Saw-horse to Fischer:



* Configuration:

The arrangement of atoms and of gp around chiral atom is known as configuration.

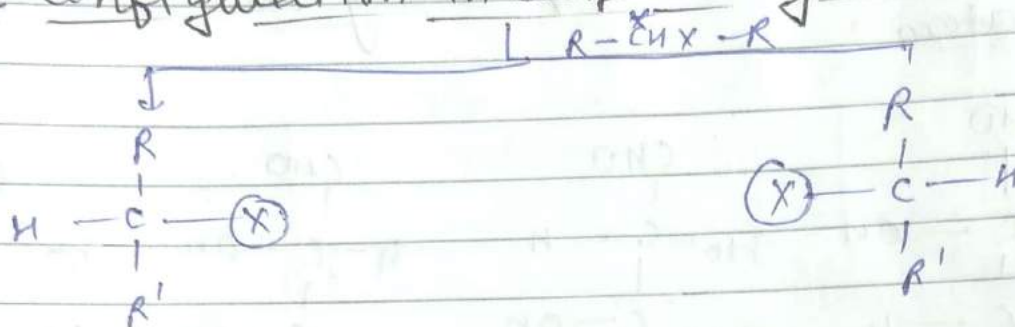
- (1) D/L configuration. (Relative configuration).
- (2) R/S configuration. (Absolute configuration)

* D/L configuration:

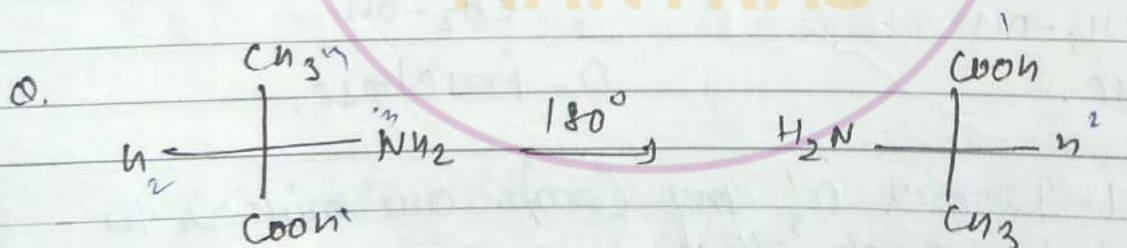
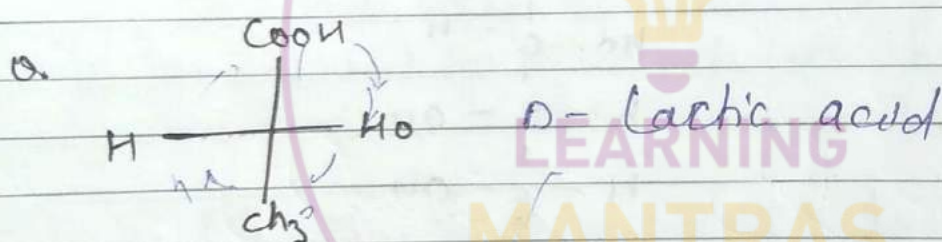
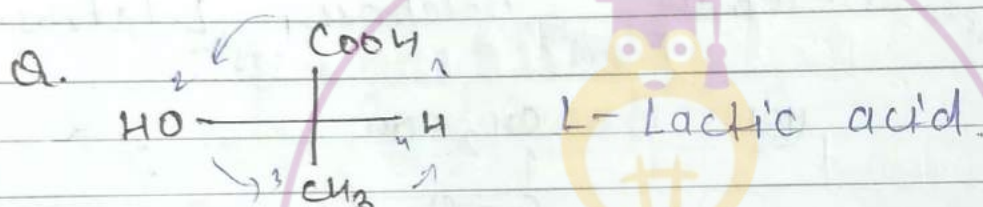
D/L always configuration see only fisher structure

O₂

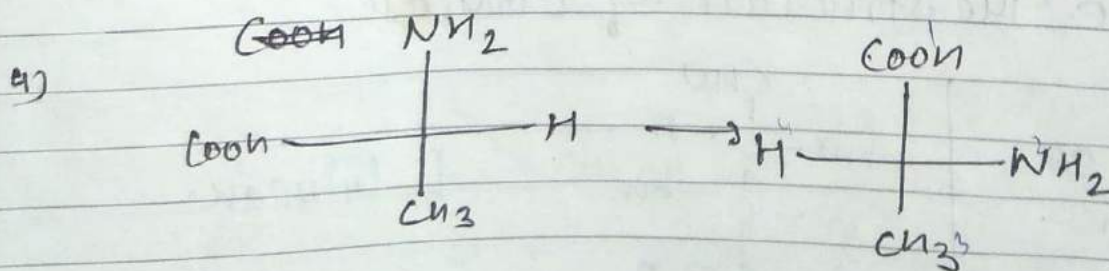
* D/L Configuration in comp. having 1 chiral atom



where X = hetero atom / group having Hetero atom.
Ex: $-F, -Cl, -Br, -I, -OH, -NH_2$ etc.



L-Alanine (L-Amino acid).



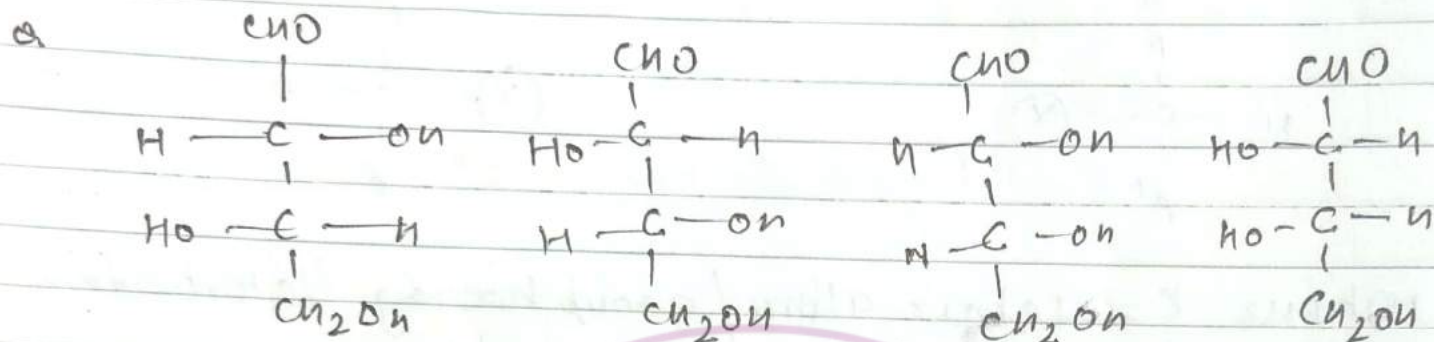
D-Amino acid (D-Alanine)

O₃

only main

only advance

* D/L Configuration in Comp. having more than one chiral atom !



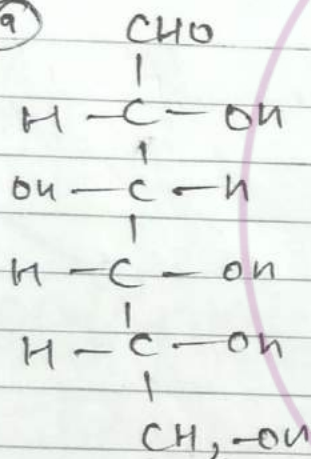
Threo L-Tetrose.

Threo D-Tetrose

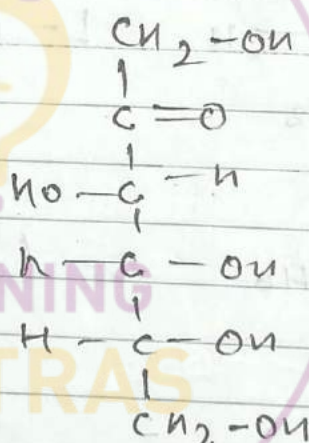
Erythro D-Tetrose

Erythro L-Tetrose

Que: (a)



(10)

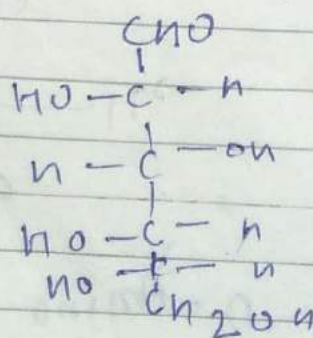
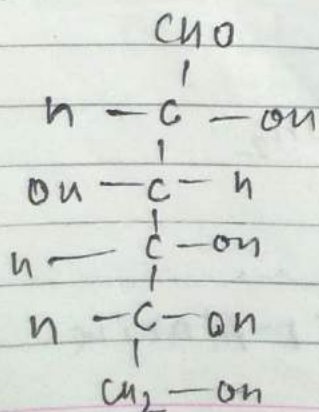


D-Glucose.

D-Fructose.

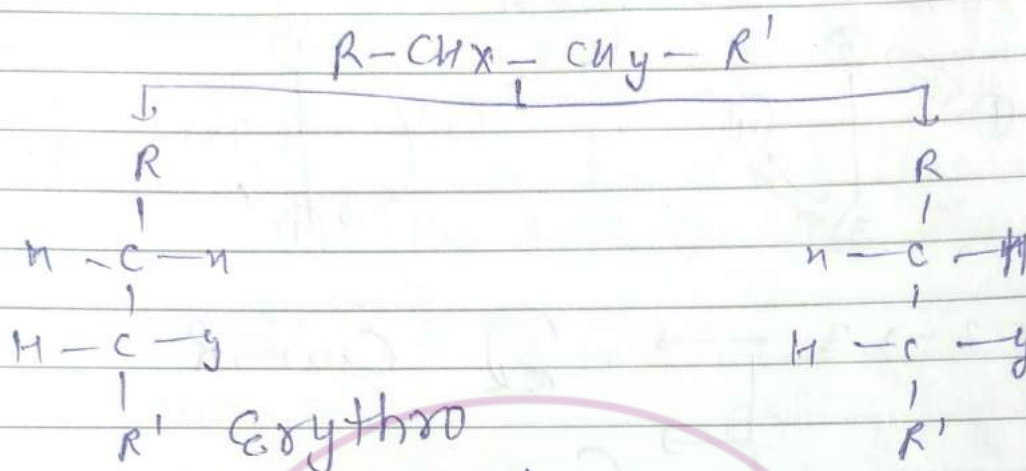
Note: D/L Isomers of any Comp. are mirror to image to each other.

Que: make the structure of L-Glucose.



L-Glucose

* Erythro/Threo Configuration :



* R/S Configuration :

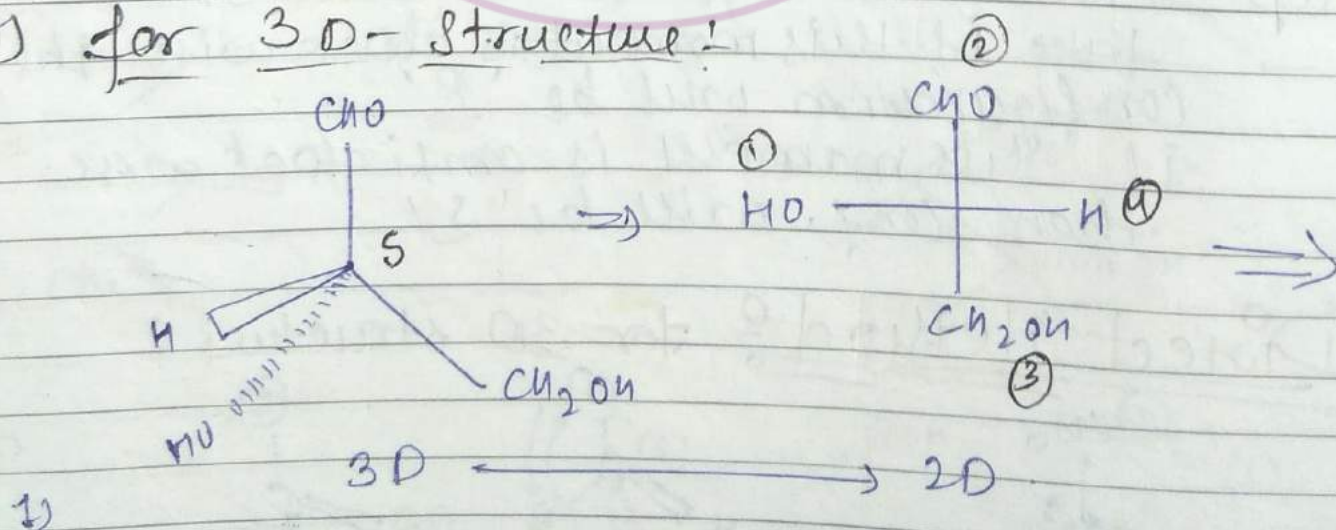
Not in Advanced

Threo

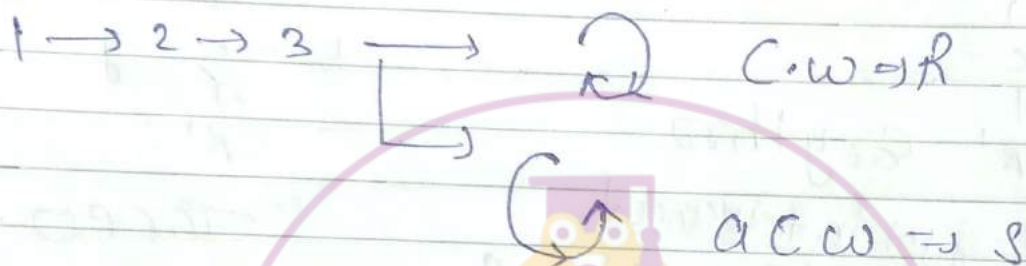
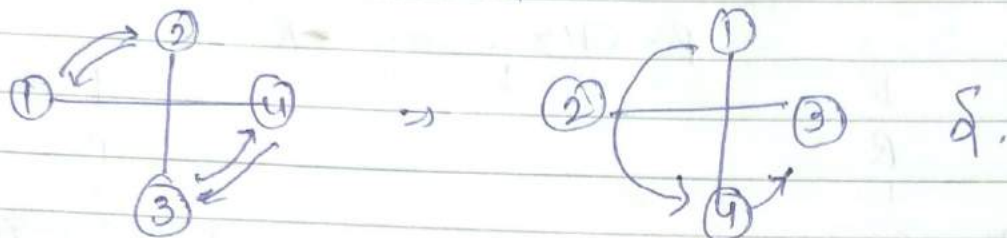
With the help of R/S configuration the configuration of each chiral center can be decided

'R' \Rightarrow Rectus \Rightarrow CW / RHS
'S' \Rightarrow Sinister \Rightarrow CCW / LHS.

(A) for 3D-Structure :



- 2) Decide Priority of atom or gp. \rightarrow 4 are vertices.
 (3) move from $1 \rightarrow 2 \rightarrow 3$. arrow



Step \Rightarrow 1. Convert 3D to 2D

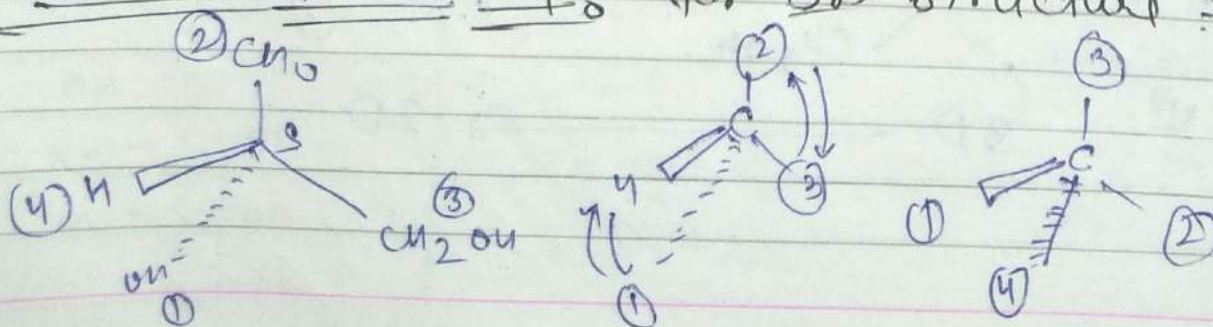
Step - 2. Decide Priority of gps attached with chiral atom.

Step - 3: Now Put 4th Priority gp at bottom line or Vertical line.

Step \rightarrow 4 Now move from 1 \rightarrow 2 to 3.

~~for~~ if this moment is clock wise then configuration will be 'R'
 If this moment is anticlock wise then conf. will be 'S'

Direct Method for 3D structure?

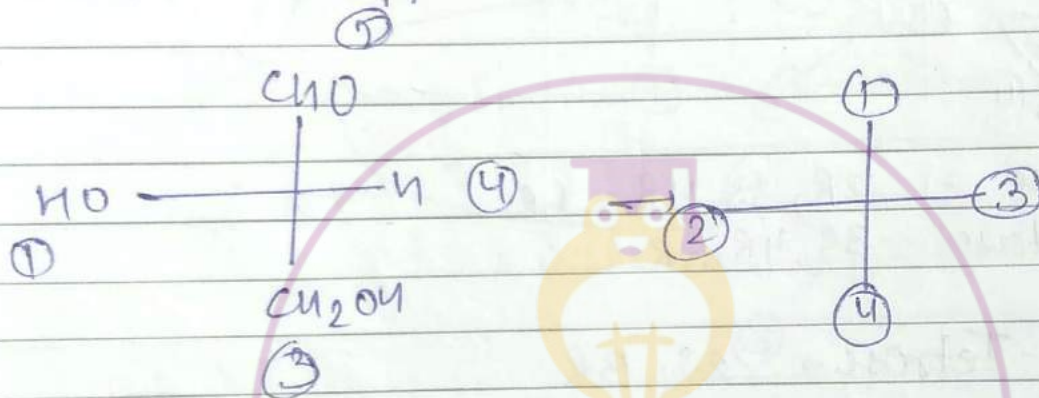


- * If 4th is on Dash then Same
 * If 4th is not on dash the result will be S and S will be R.

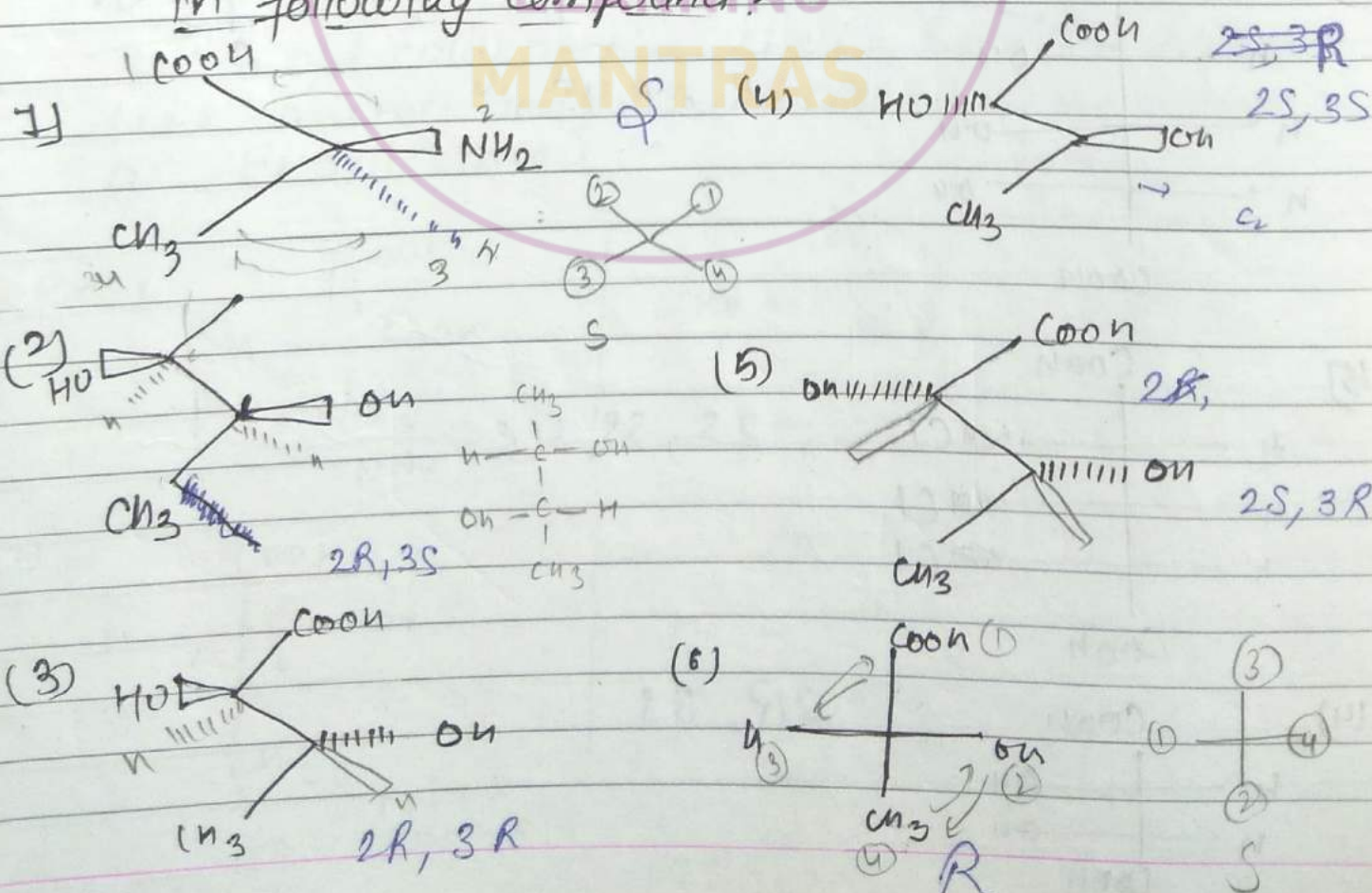
Condition: 4th must be at dash line.

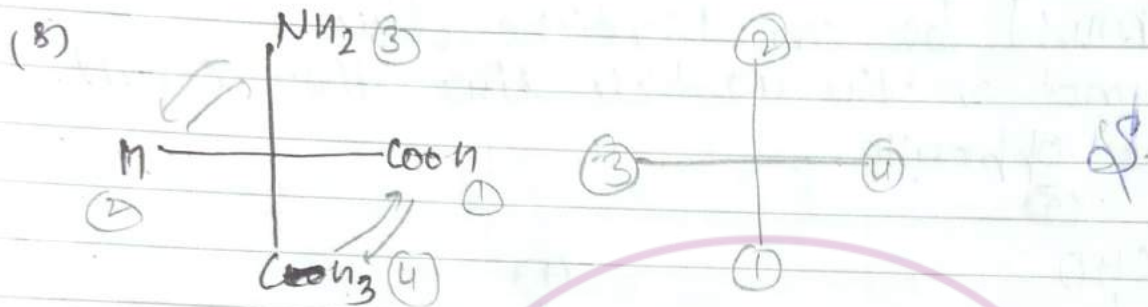
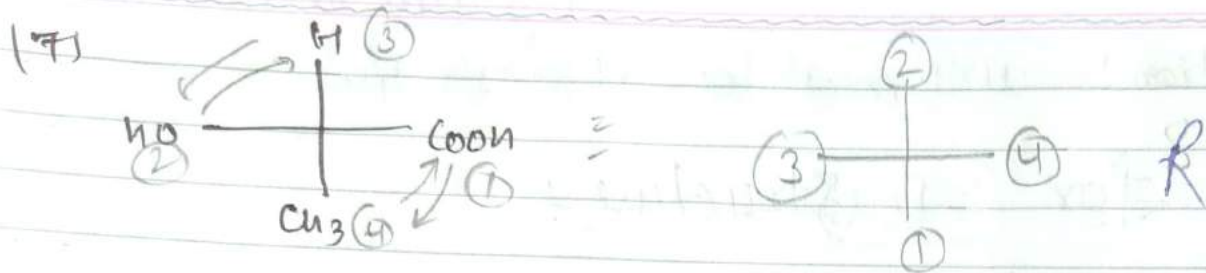
* ~~For~~ for 2D structure:

4th must be on vertical line.
 If 4th is not on the vertical line then result will be opposite.



Q. Decide R/S configuration each chiral carbon in following compound!

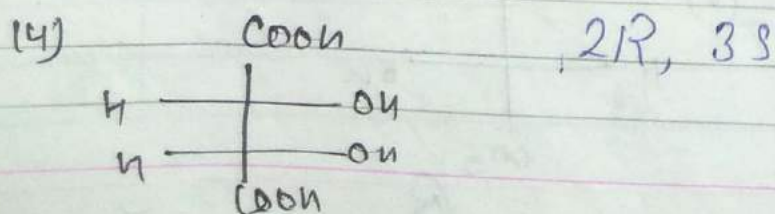
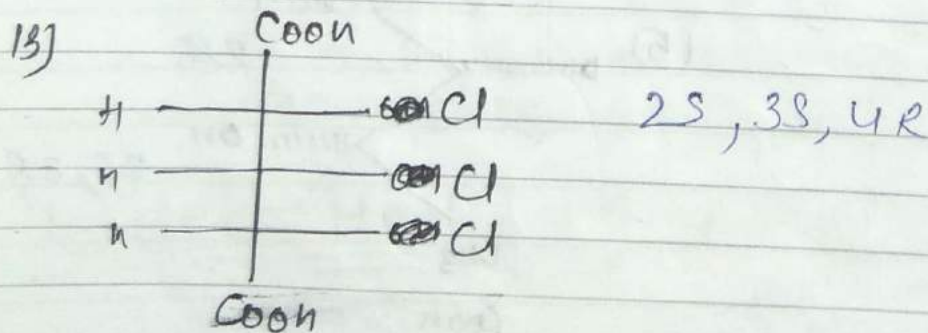
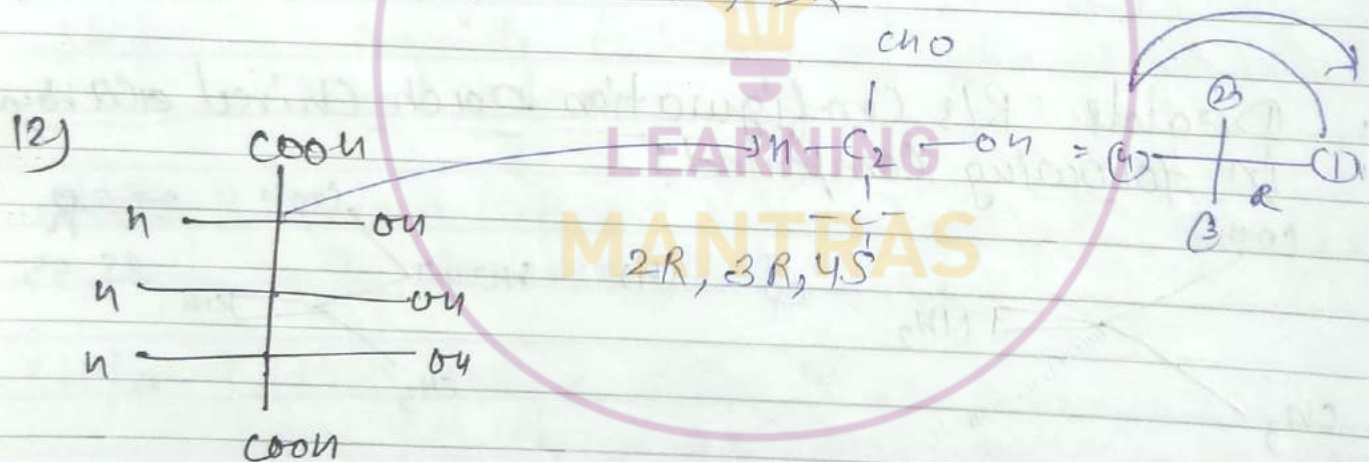


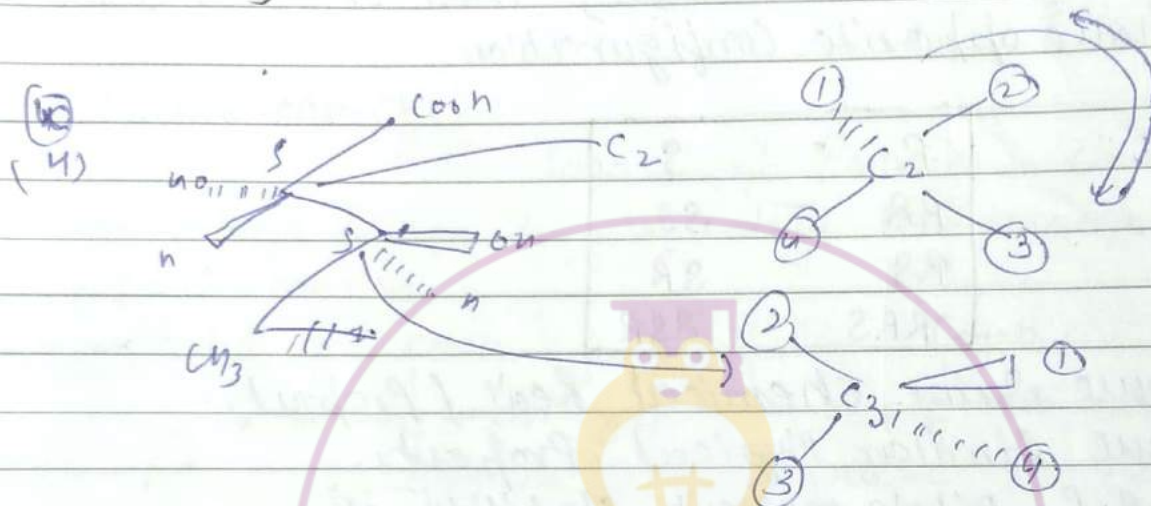
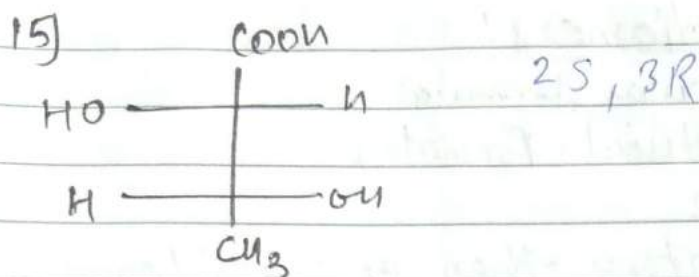


(9) D-Glucose \rightarrow 2R, 3S, 4R, 5R

(10) D-Fructose 3S, 4R, 5R

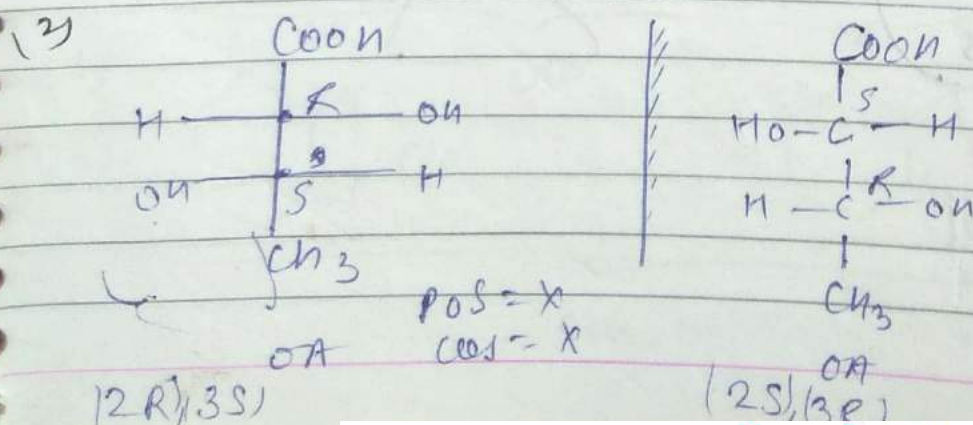
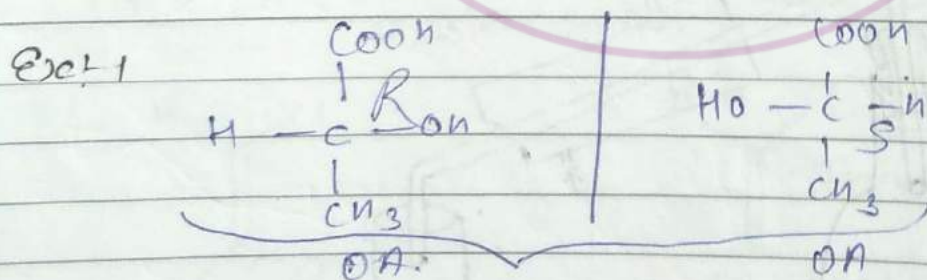
(11) Threo-D-Tetose \rightarrow 2S, 3R





* Enantiomers / Anti-mers / Enantiomorph / Antipods.

Two optically active stereo isomers which are mirror image to each other are known as Enantiomers.



Optical active mirror image

* Characteristic of enantiomers:

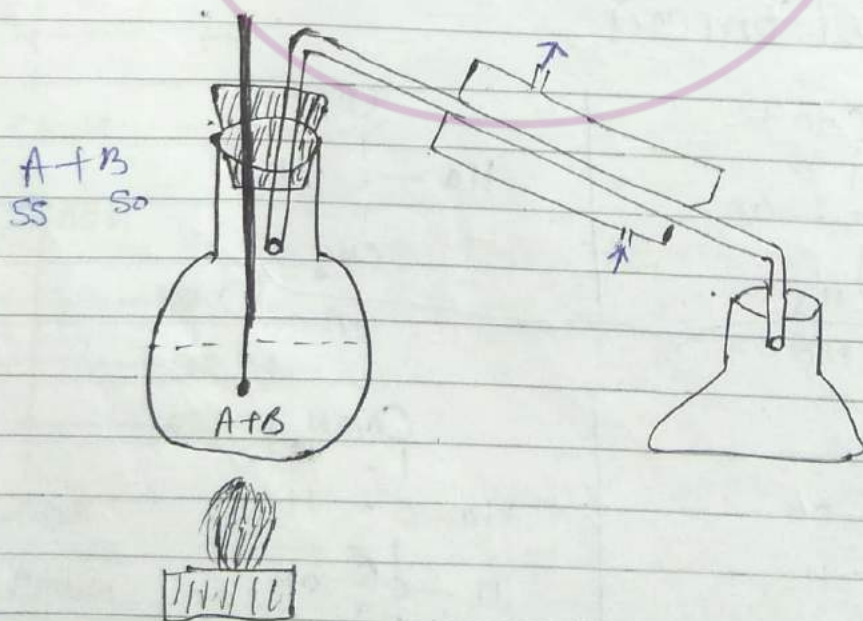
- 1) They have same M.f (Molecular formula)
- 2) They have same S.f (structural formula)
- 3) They are optically active.
- 4) If one is dextro rotatory then other is levo.
- 5) They have opposite configuration.

R	-	S
RR	-	SS
RS	-	SR
RRS	-	SSR

- 6) They have same chemical reactivity / property.
- 7) They have similar physical properties.
(M.P, B.P, Dipole moment, stability etc.)

Ques! The melting point of (R)-Butan-2-ol is 55°C . then what will m.p. of (S)-Butan-2-ol

- ☒ (a) $+55^{\circ}\text{C}$
☐ (b) -55°C
☐ (c) 0°C
☐ (d) Can not decided



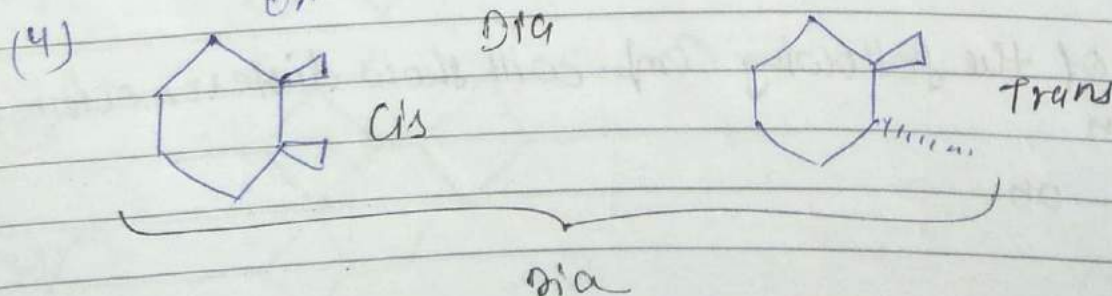
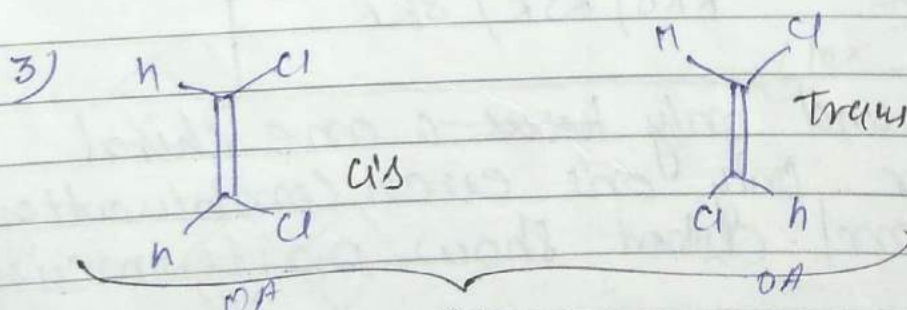
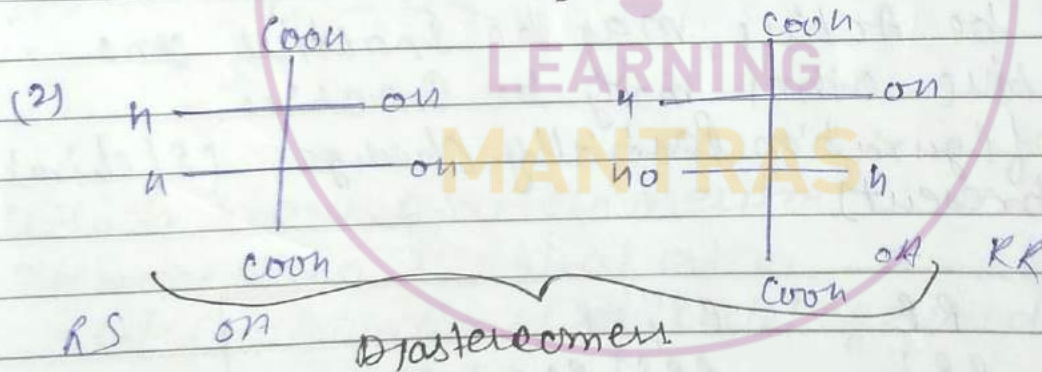
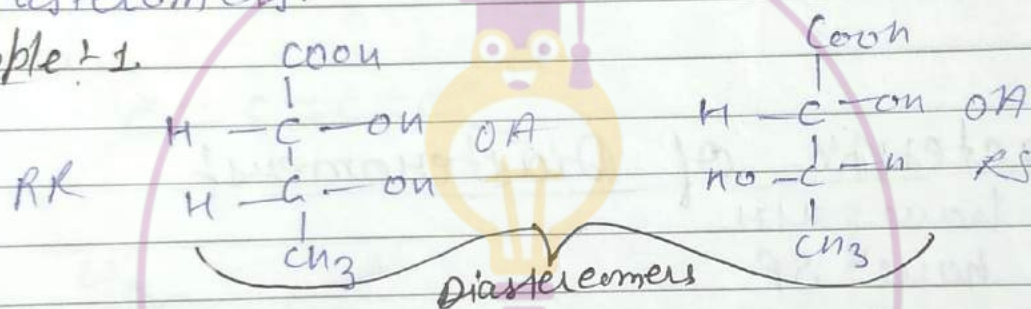
8) Enantiomers cannot separate by fractional distillation.

9) Enantiomers are non super imposable mirror image of each other.

* Diastereomers:

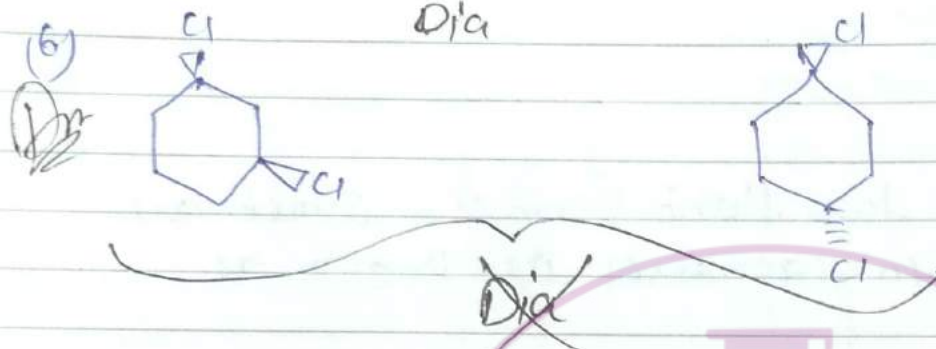
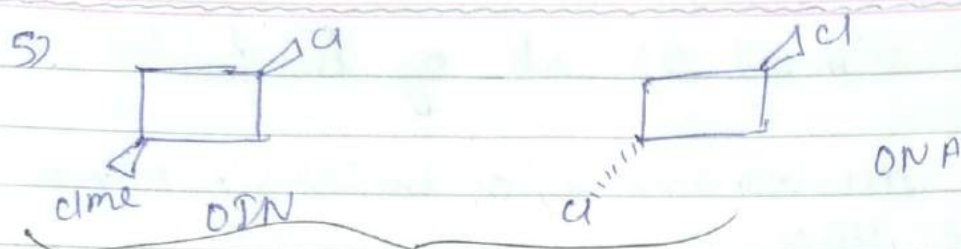
The two stereoisomers which are not mirror image to each other are known as diastereomers.

Example: 1.



Di - 2 or more

Stereo
Mirror Image



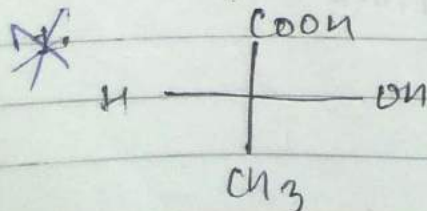
* Characteristic of Diastereomers!

- 1) They have different NMR.
- 2) They have different SF.
- 3) They may be Active, may be Inactive, one may be active other may be Inactive.
- 4) Their configuration partially change (if chiral atom present).

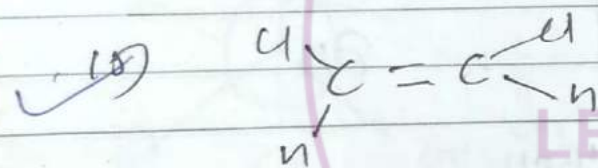
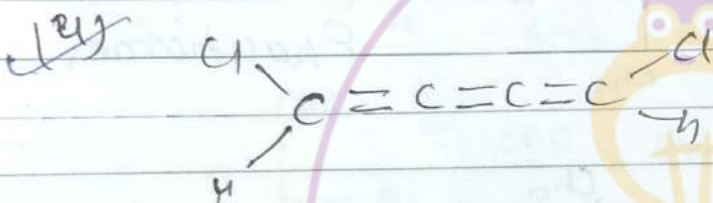
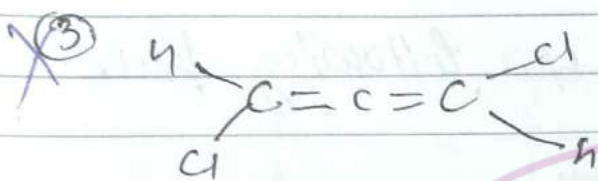
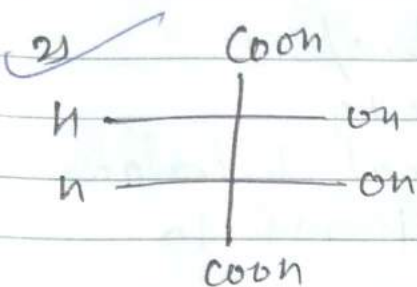
RR	RS/SR
RRR	RRS/RSR/SRR

5) Those compound only have a one chiral centre or only one even combination of double bond. do not show diastereomerism.

Ques! which of the following Comp. will show diastereomerism!



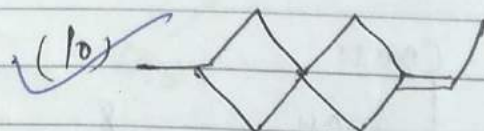
(A's Traces) ✓



1) Which show Diastereomerism.

1) 2 or more than 2 chiral atom.

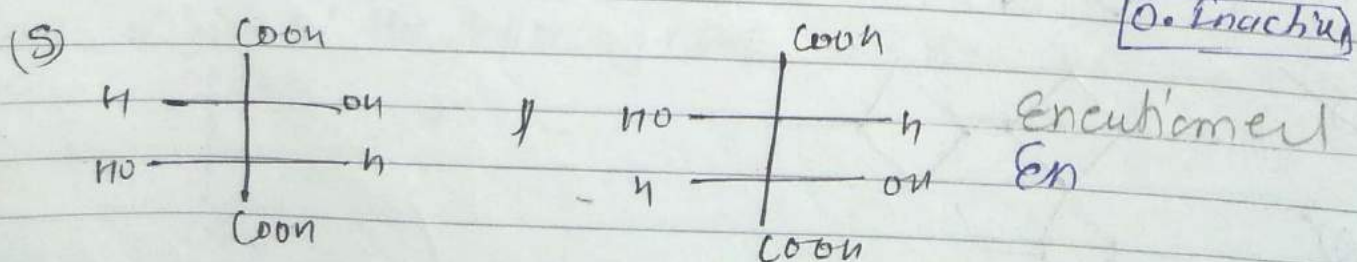
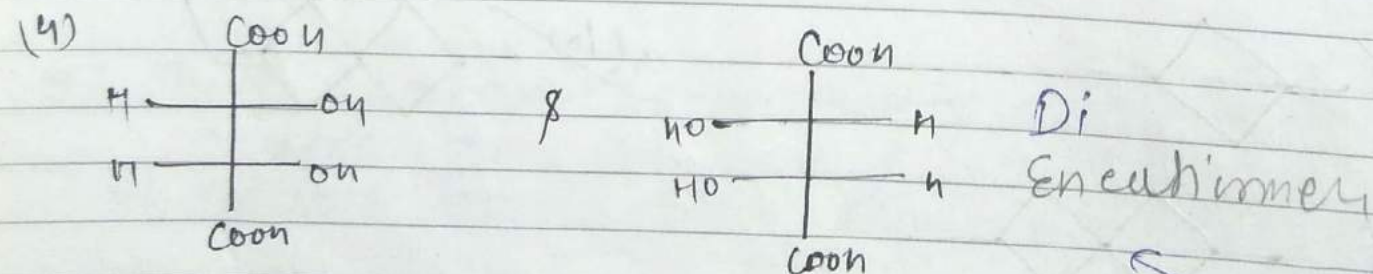
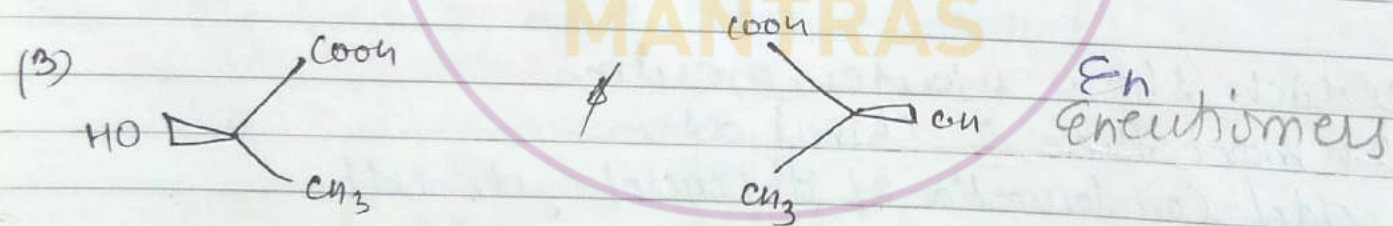
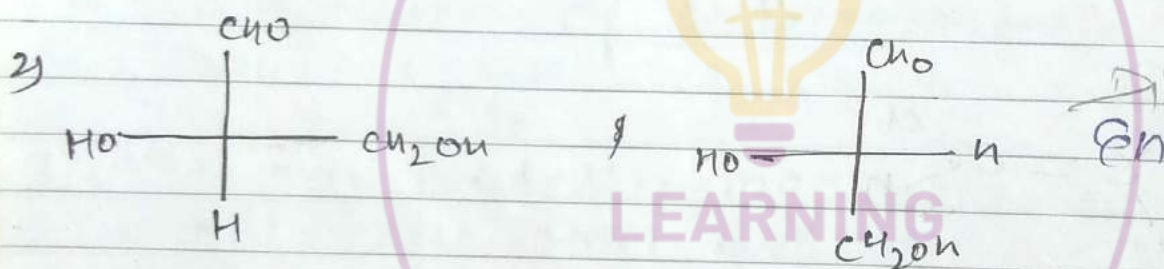
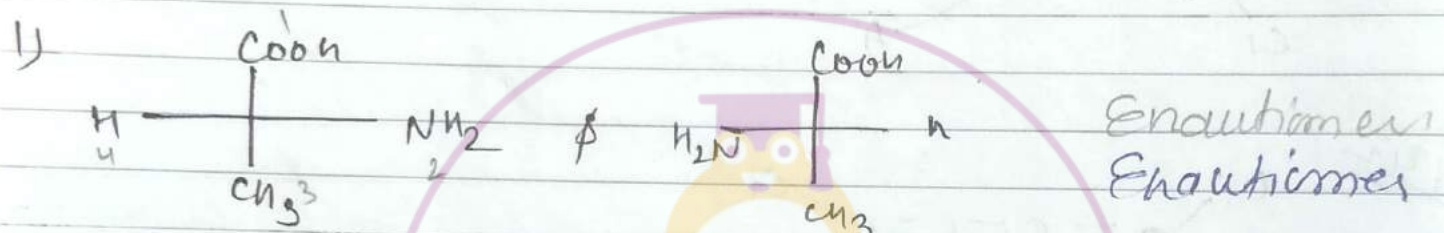
2) Odd combination of db/cycle/cycle + db.



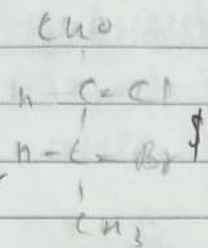
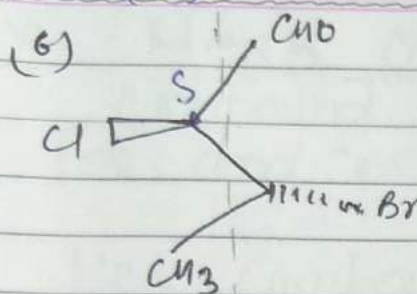
See cos \rightarrow cycle planar syst
POS

- 3) They have different chemical property
- 4) They have different physical property
- 5) They can be separated by fractional distillation
- 6) Mesoisomers are also diastereomers to each other.

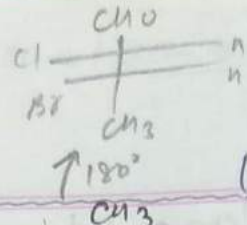
Que 5 Decide Relationship b/w following pairs



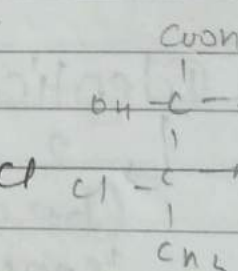
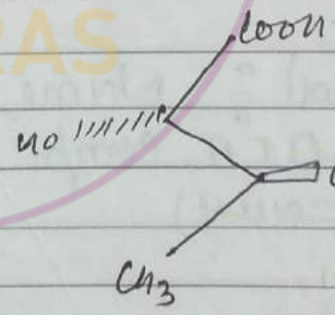
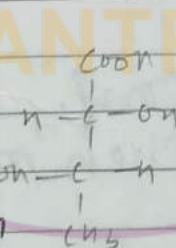
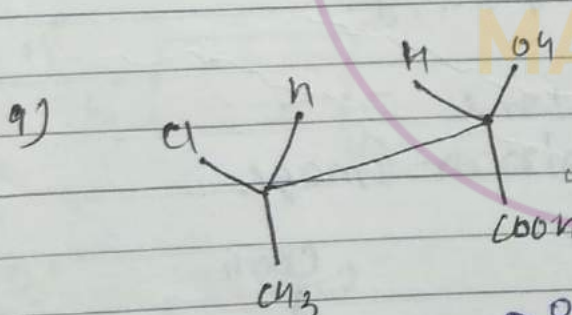
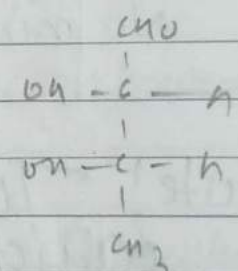
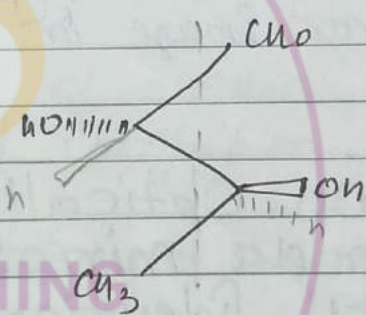
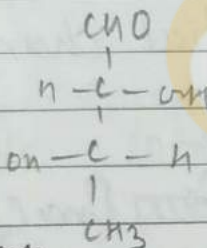
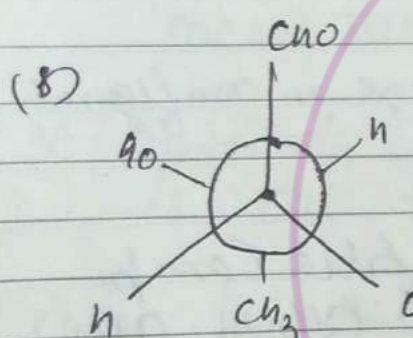
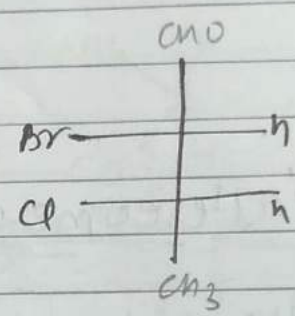
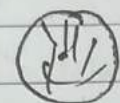
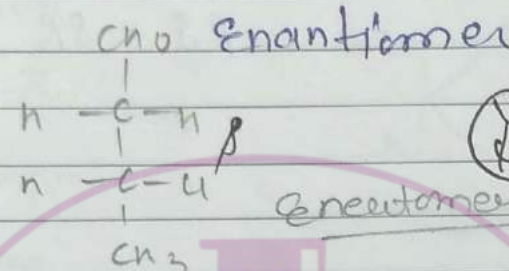
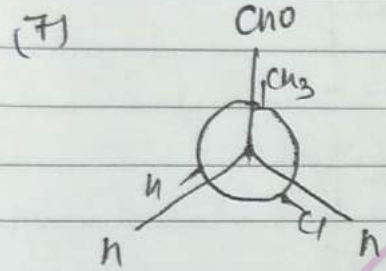
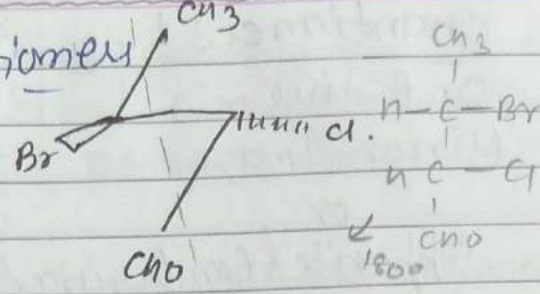
Conf. opt. 30, mirror image.



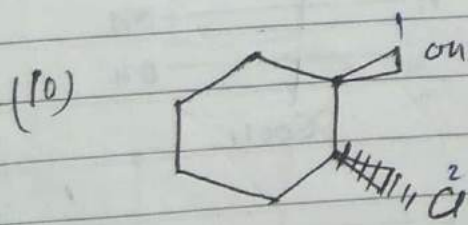
Enantiomers



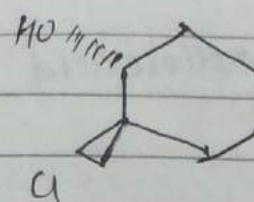
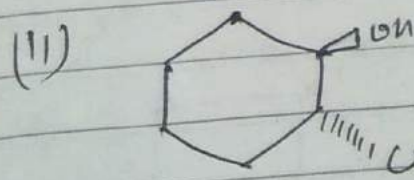
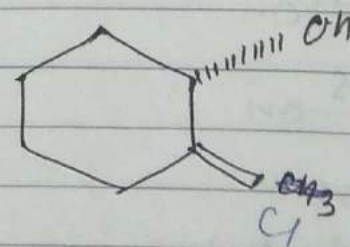
(2R, 3S)



D⁺ D⁻



Enen



Identical

* Enantiomers! 1) Stereo Isomers

1) O. Active

2) Mirror Image

or
Opposite Configuration

R-S

RR-SS

RS-SR

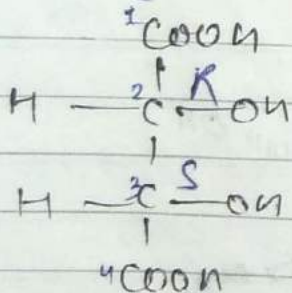
* Diastereomers!

1) Stereoisomers

2) No mirror Image or Partial change in configuration

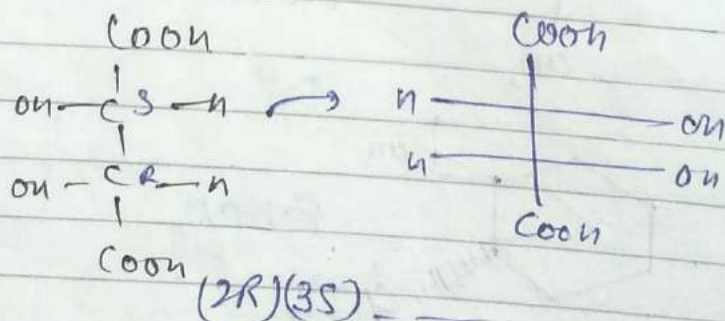
Note! Two Optically ~~act~~ Inactive Comp. which are mirror to each other are known as Identical Compound.

* Identical: Name, Conf., same.
↳ 2 OIA Comp when mirror Image (because)



2R, 3S, -2-3-

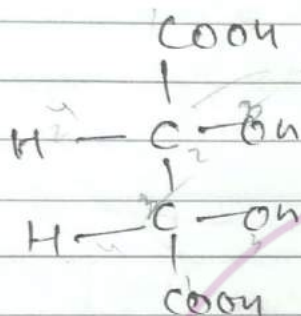
hydroxybutan-2,3-diol acid



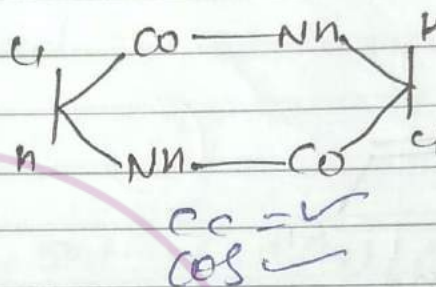
* Meso Compounds:

Compound which have chiral Carbon with pos and cos are known as meso compound.

Meso compounds are optically inactive compound.



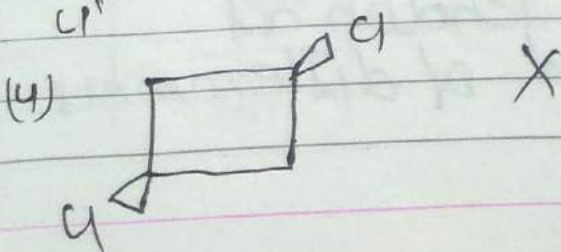
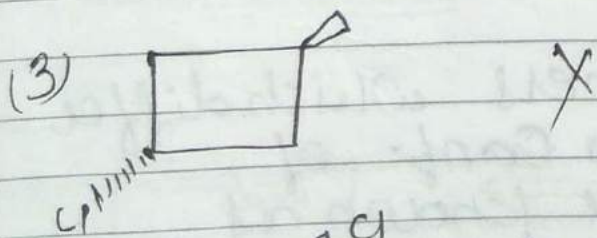
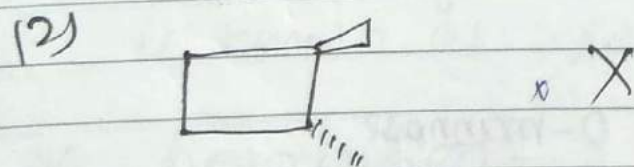
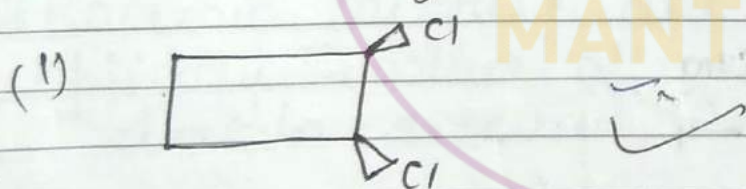
meso-Tartaric acid

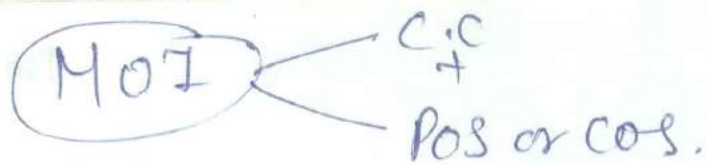


meso compound.

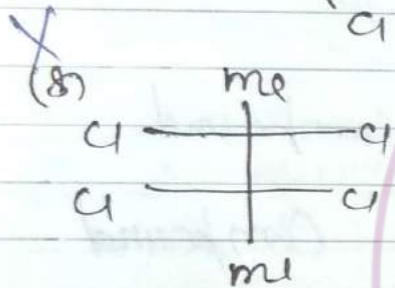
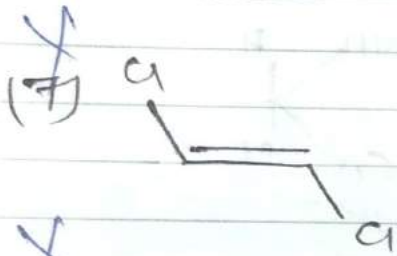
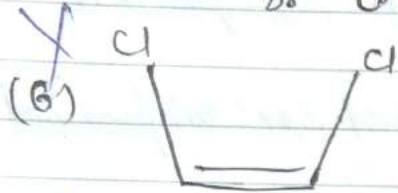
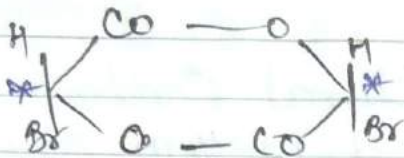
* Optically inactivity of meso compound is due to internal compensation.

Ques: Identify meso-compound among following!

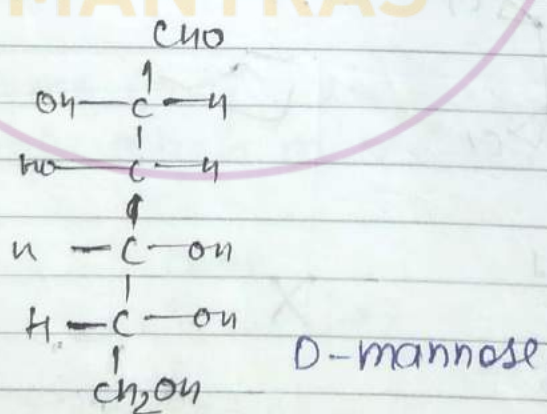
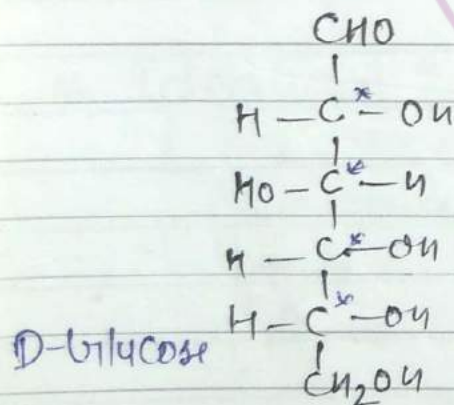




(5) X



* Epimers \Rightarrow



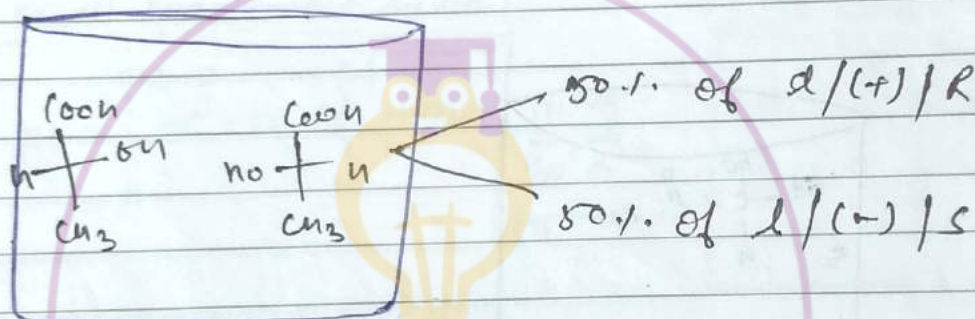
\Rightarrow The two Stereo Isomers which differ to each other due to Conf. of one chiral atom are known as Epimers. They are a kind of diastereomers.

* Racemic Mixture:

The equimolar mixture of D and L Isomers of any compound is known as Racemic Mixture

OR

The mixture having 50% D and 50% L Isomer of any compound.

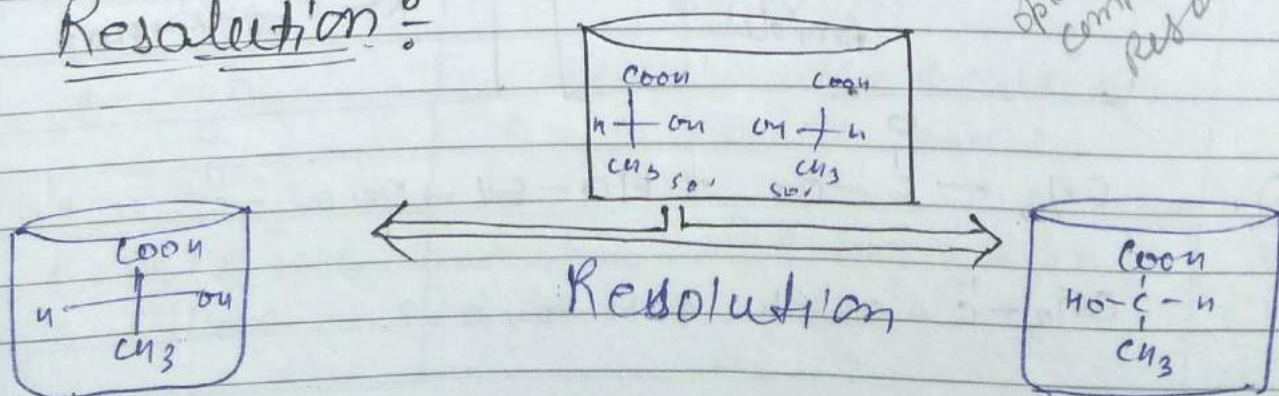


* It is also known as (\pm) or (dl) or (RS) mixture

* Racemic mixture are optically inactive mixture
Optically Inactive of racemic mixture is due to external compensation.

The process by which racemic mixture is prepared is known as Resolution

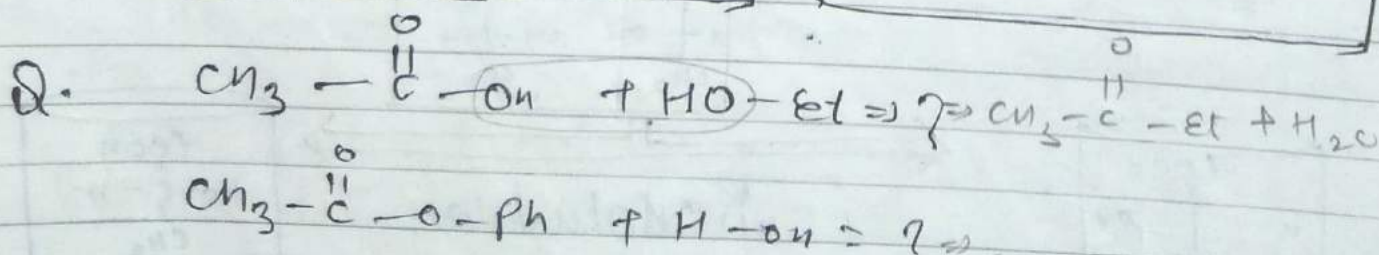
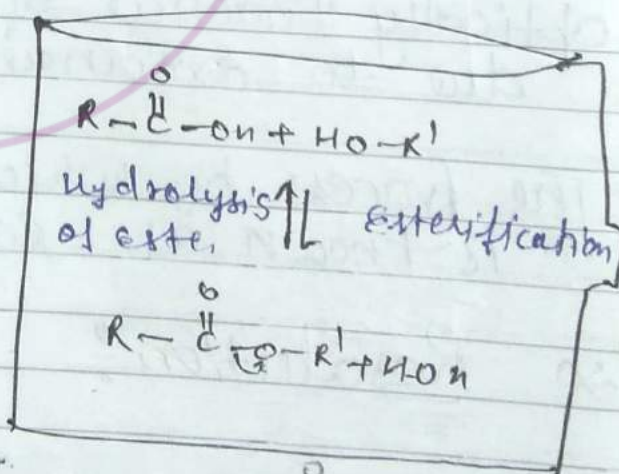
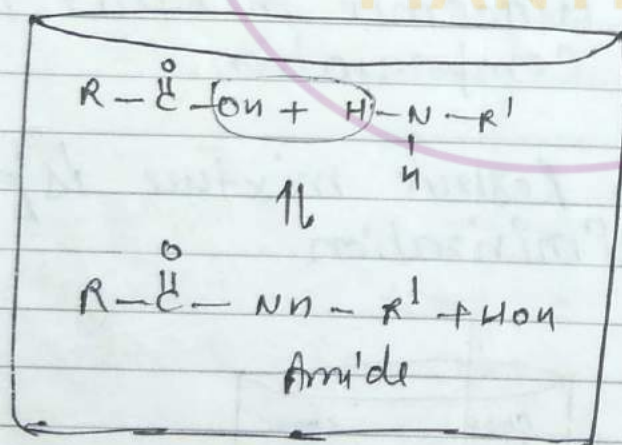
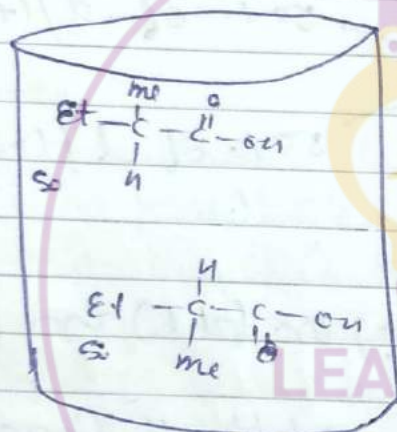
* Resolution:



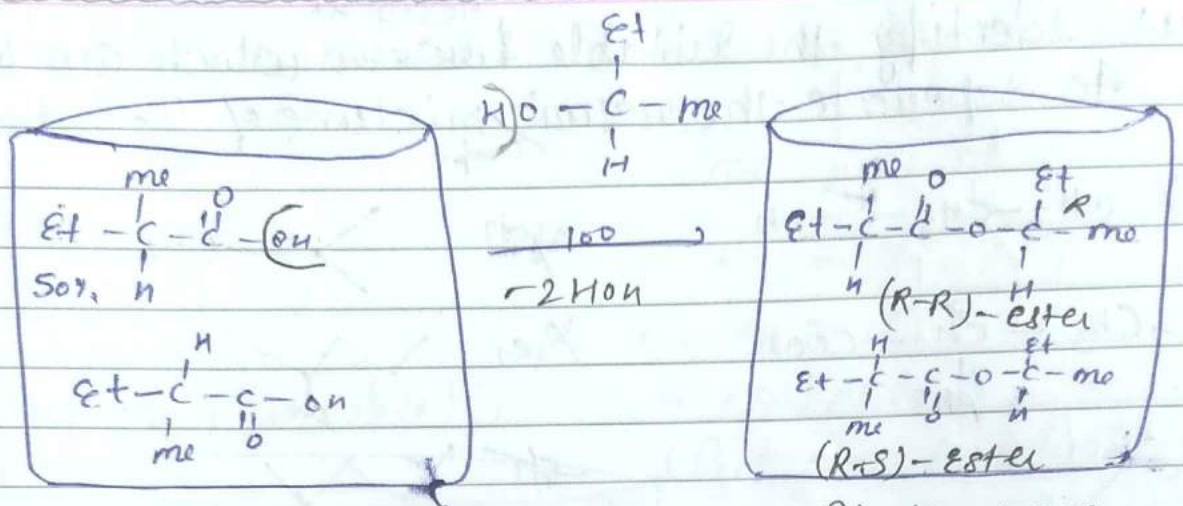
Resolution: \rightarrow By mechanical method \rightarrow
 \rightarrow By Biological " \rightarrow
 \rightarrow By chemical method/
 or

By Diastereomeric formation.

* Resolution of Racemic mixture by diastereomeric formation:

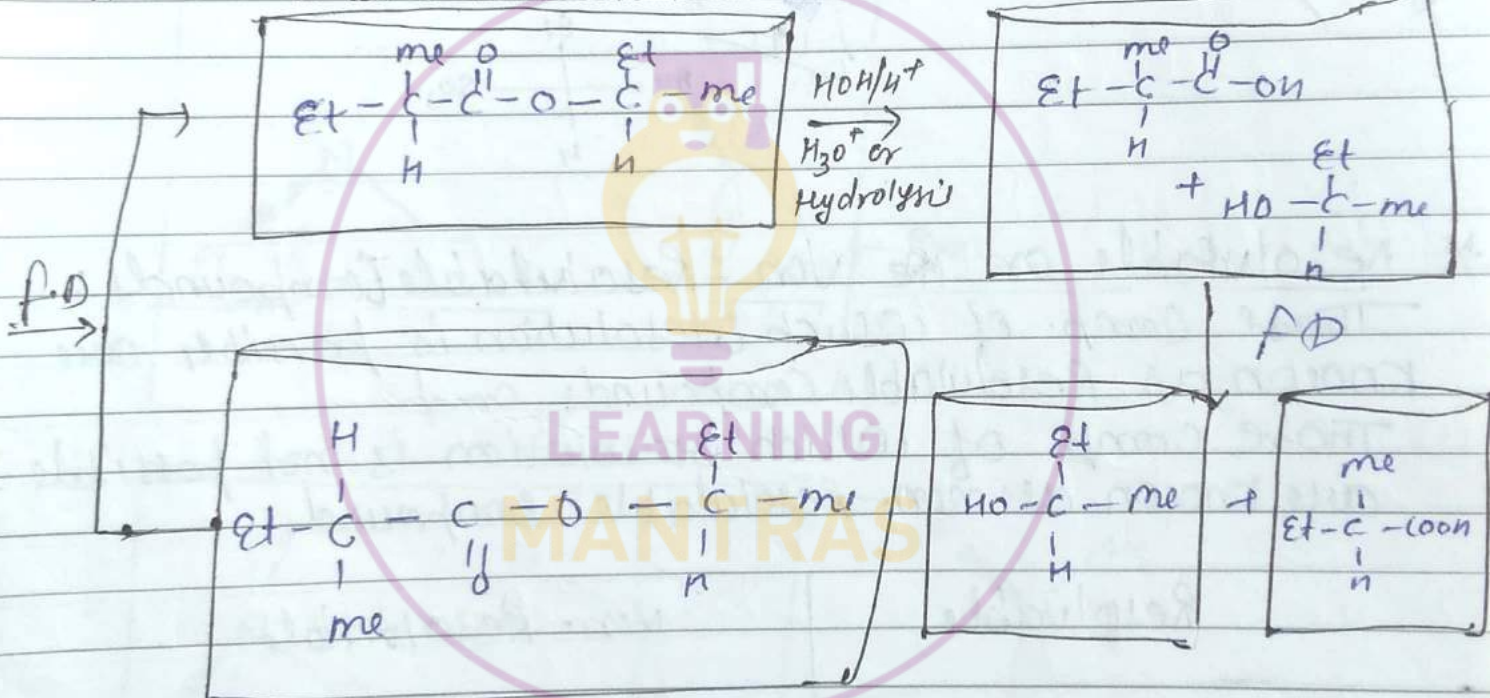


H.W = Race: 9 and 10.



Rm or ± or all Enantiomeric mixture

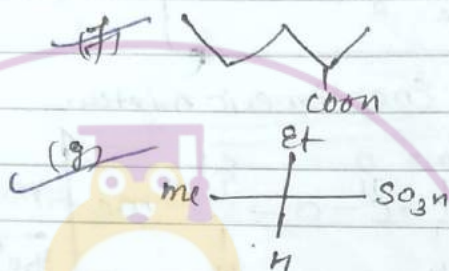
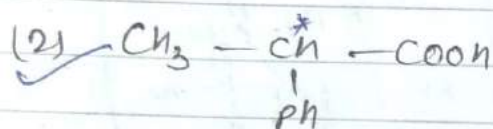
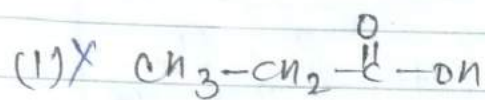
Diastereomers



X

RM	Separator
R-COOH	optically active R-OH / R-NH ₂ (" " " Ball).
R-OH / R-NH ₂ Alcohol	(" " " R-COOH / R-SO ₃ H) (" " " Acid)

Ques: Identify the suitable ^{reagent} ~~reagent~~ which can be used to separate the racemic mixture of 2-Butanol. Base



* Resolvable or Non-Resolvable Compounds:
 Those comp. of which resolution is possible are known as Resolvable Compounds and
 Those comp. of which resolution is not possible are known as non-Resolvable compound.

Resolvable

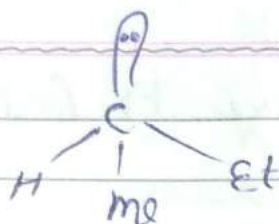
Non-Resolvable

All O.A and chiral compound

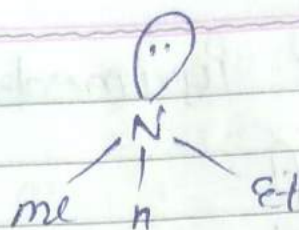
All O.A / Achiral compound.

* Exception: All Pnmedal Compound with elements are 2nd period at Centre are non Resolvable.

Ex:

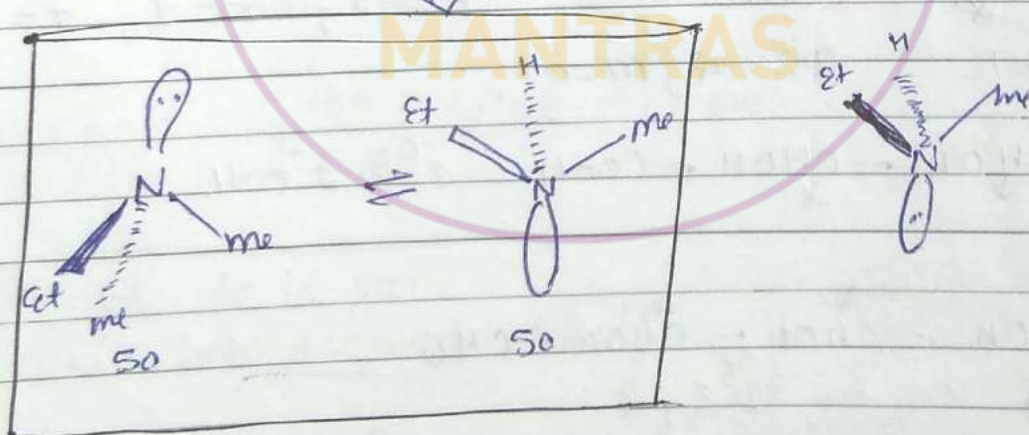
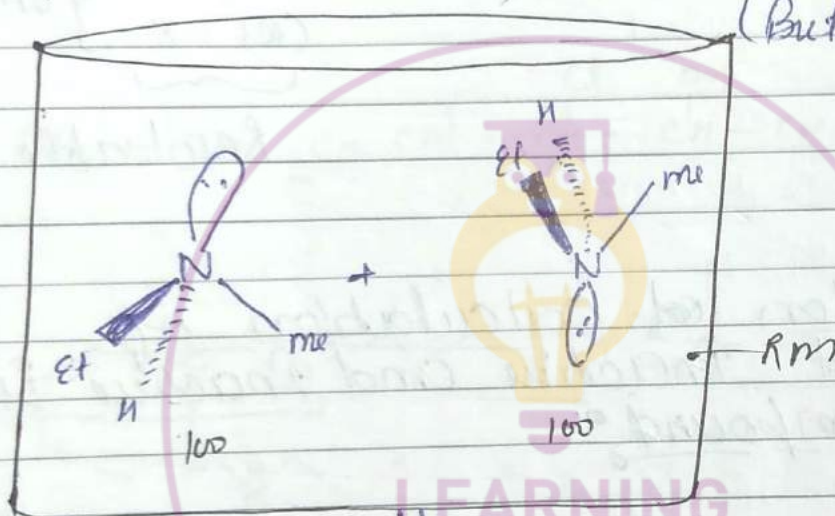


(ii)

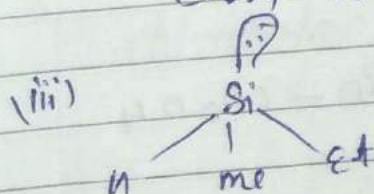


Pos = X
 Cos = X
 OA/chiral
 (But non Resolvable)

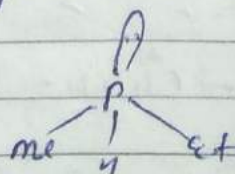
Pos = X
 Cos = X
 OA/chiral
 (But non Resolvable)



* Pyramidal structure having central atom or Bond or higher period are Resolvable



(iv)



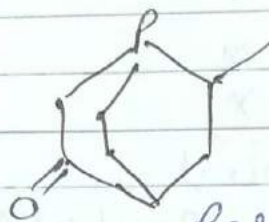
Pos = X
 Cos = X
 OA/chiral
 Resolvable

Pos = X
 Cos = X
 OA/chiral
 Resolvable

* Cyclic Pyrimidal structure are Resolvable.

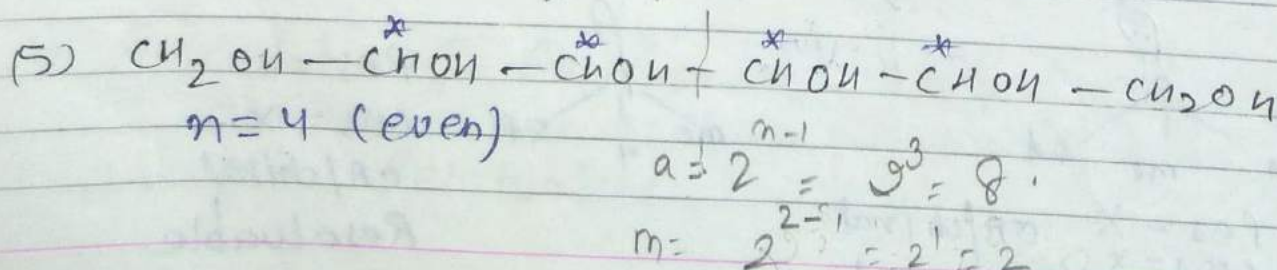
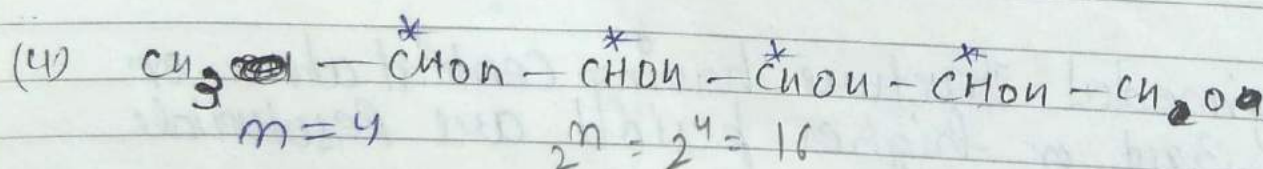
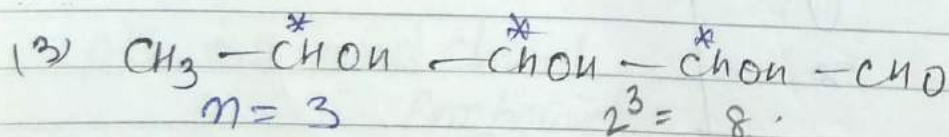
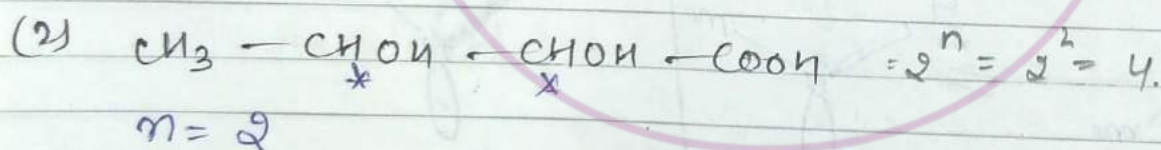
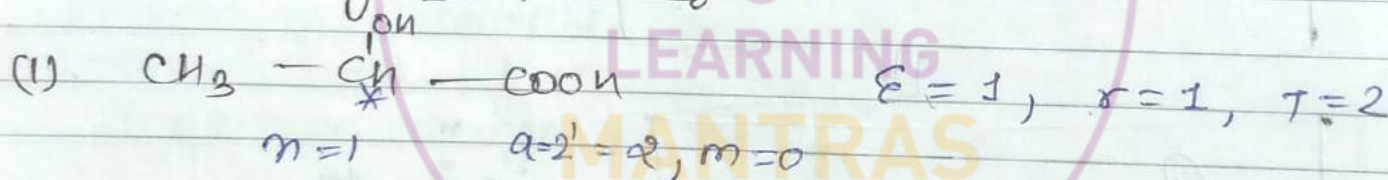


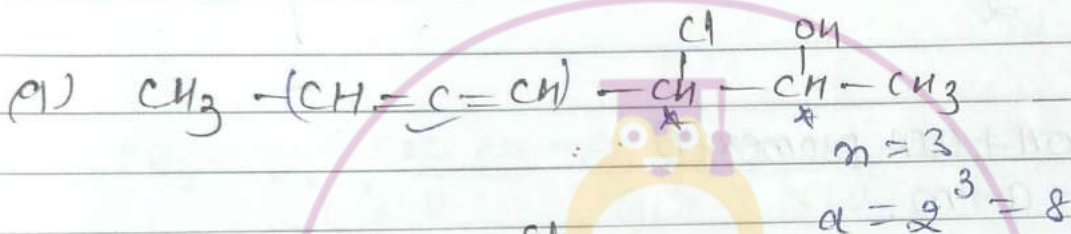
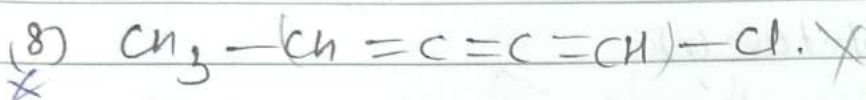
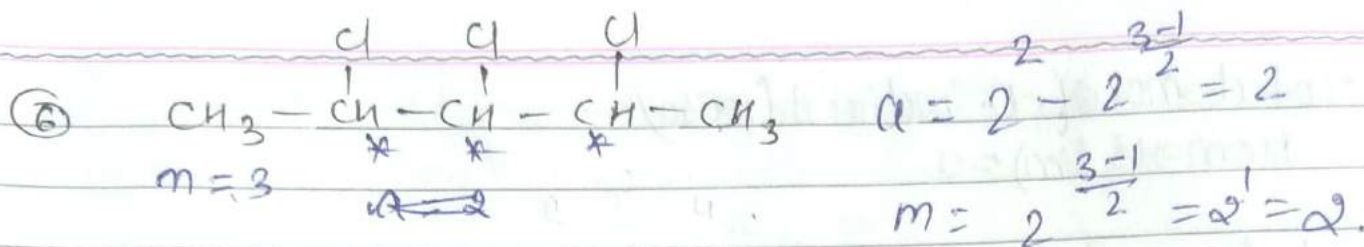
Pos = x
 $\underbrace{\text{Cos} = x}_{\text{OA}}$
 Resolvable



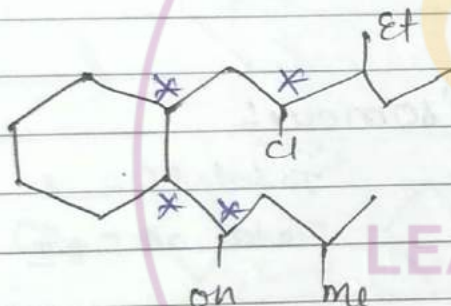
Pos = x
 $\underbrace{\text{Cos} = x}_{\text{OA}}$
 Resolvable

* Determination of calculation of
 * Optical and Inactive and Inactive isomer
 in any compound.



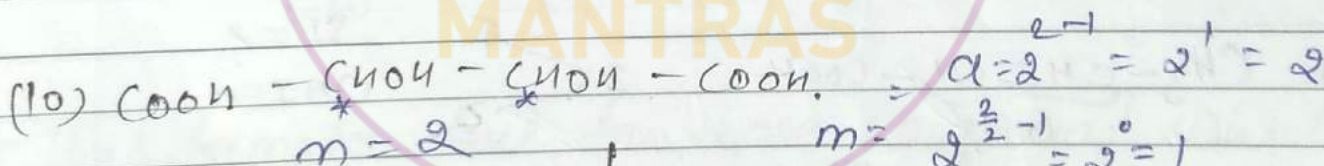
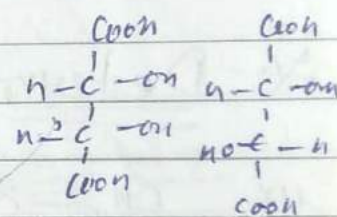


(10)



$n = 4$

$\alpha = 16$



When molecule is not divide into two equal half.

a) Total no. of enantiomers = 2^n
 $n = \text{optical isomer centre}$
 OR
 chiral atom or even combination of db/cycle/c+db

When molecule can be divide into two equal half.

even	odd
$\alpha = 2^{n-1}$	$\alpha = 2^{n-1} - 2^{\frac{n-1}{2}}$
$m = 2^{\frac{n}{2}-1}$	$m = 2^{\frac{n-1}{2}}$

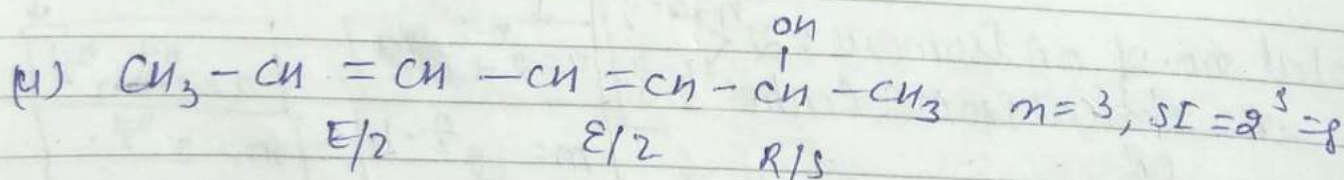
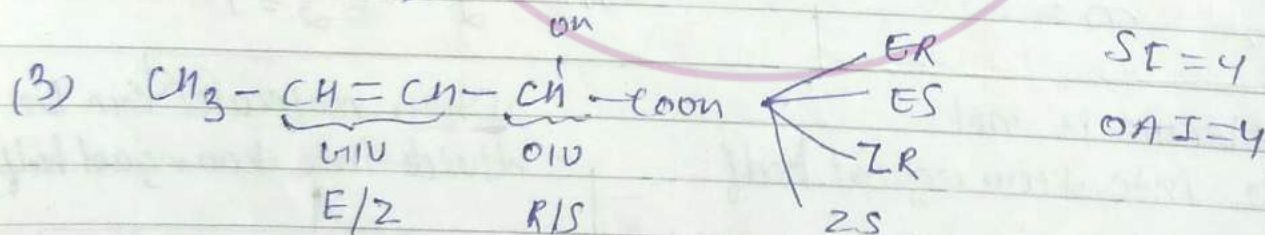
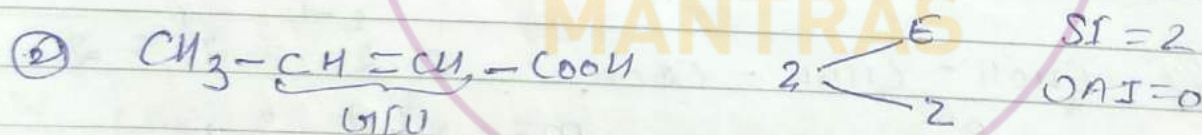
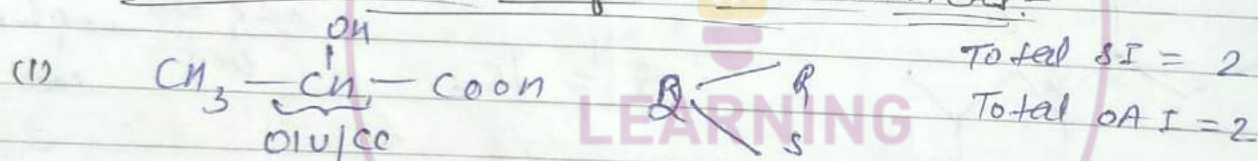
b) Total no. of optically inactive (meso) isomers (m) = 0

c) Total no. of Enantiomeric pair (EP) = $\frac{a}{2}$

d) Total no. of Racemic mixture (r) = $\frac{a}{2}$

e) Total no. of OA + OAI Isomer (T) = a + m

* Determination of Stereoisomers:



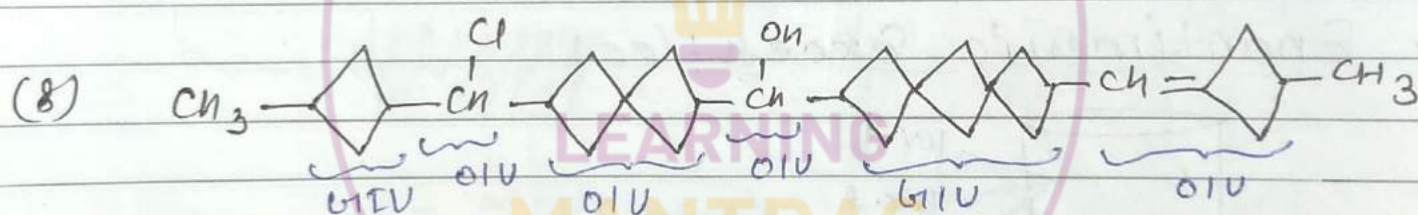
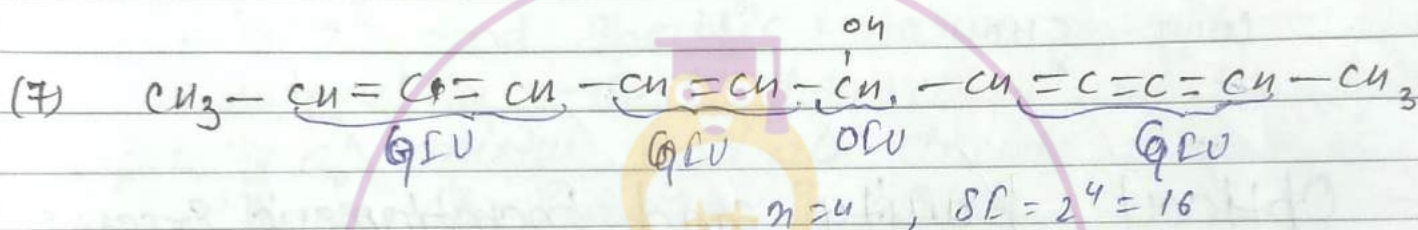
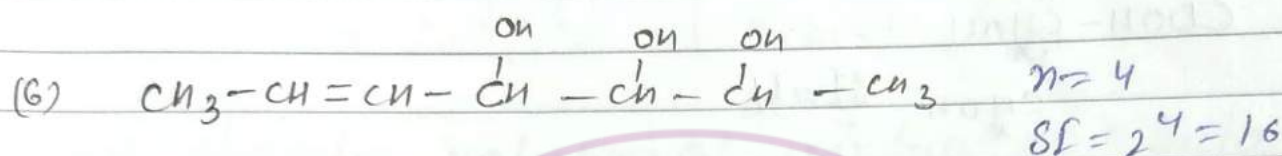
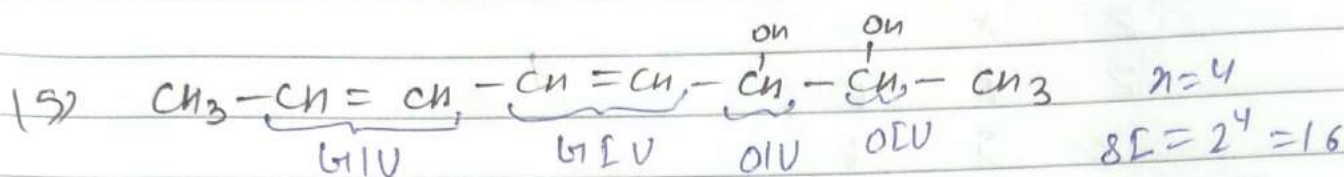
$$\text{Total no of S.I} = 2^n$$

$$n = \text{No. of (H/U + OH)}$$

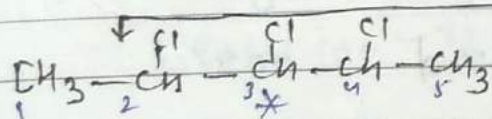
even = 01
odd = 01

LSU = odd Combination of db/cycle/cycle + db

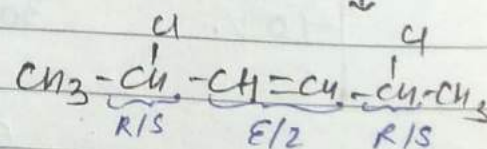
OSU = even " " " " " "



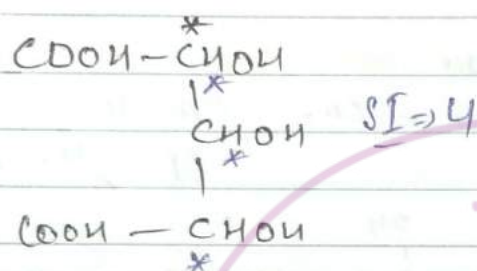
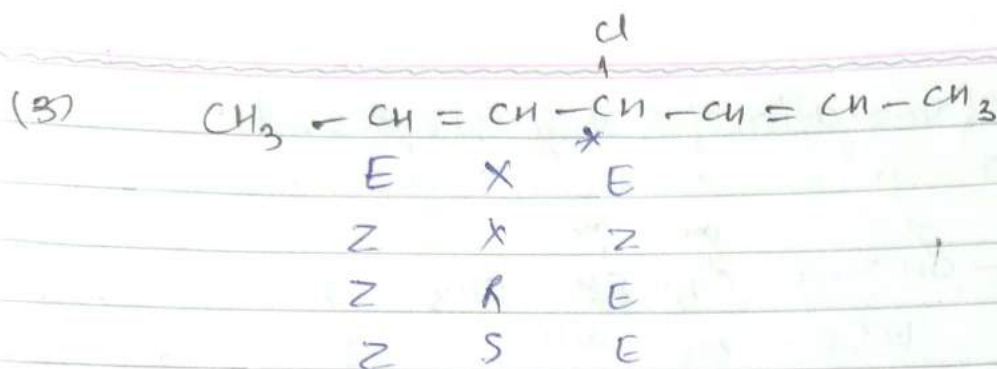
This formula fail for those compound which can be divided into two equal half.



R	X	R (2R) (4R) - 2,3,4-TCP
S	X	S (2S) (4S) (4S) " "
R	R	S (2R) (3R) (4S) " "
R	S	S (2R) (3S) (4S) + " "

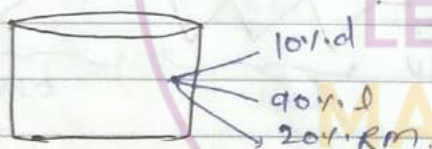


R	E	R	(2R) (3E) (5S) - 2,5-dichloro hex-3-ene
S	E	S	" "
R	E	S	" "
S	E	R	" "
R	Z	R	" "
S	Z	S	" "
R	Z	S	" "
S	Z	R	" "



* Optical purity and Enantiomeric Excess:

1. Enantiomeric Excess (ee)



d %	l %	R.M %	ee	op
10 %	90 %	20 %	80 % d	80 % d
50 %	50 %	100 %	0 %	0
10 %	30 %	60 %	40 % d	40 % d

$$ee = \left| \frac{d - l}{d + l} \right| \times 100 \quad \text{--- (1)}$$

2. Optical purity (op):

$$\boxed{op = \frac{\alpha_{\text{mixture}}}{\alpha_{\text{pure}}} \times 100} \quad \text{--- (2)}$$

op = ee

$\alpha_{\text{mixture}} = \text{S.R. of mixture}$

$\alpha_{\text{pure}} = \text{S.R. of pure isomer}$

We know $op = ee$, then

$$\frac{[d-l]}{[d+l]} = \left(\frac{\alpha_{\text{mixture}}}{\alpha_{\text{pure}}} \right) \quad \text{--- (3)}$$

Ques! ^{D/L} specific rotation of mixture of 2 Butanol is -9.72° and specific rotation of -1° α Butanol is -13.5° then what will the optical purity of mixture and also decide d and l isomer of 2 Butanol in mixture.

M-1

~~$\frac{d-l}{d+l} = \frac{-9.72}{-13.5}$~~

~~$\frac{d-l}{d+l} = \frac{13.5}{9.72}$~~

$$op = \frac{\alpha_{\text{mix}}}{\alpha_{\text{pure}}} \times 100 = \frac{-9.72}{-13.5} \times 100 = \frac{-972}{-13.5}$$

$$= 72\% \text{ l.}$$

$$op = 72\% \text{ l.}$$

$$ee = 72\% \text{ l.}$$

$$RM = 100 - 72 = 28\% \begin{cases} 14\% \text{ d} \\ 14\% \text{ l} \end{cases}$$

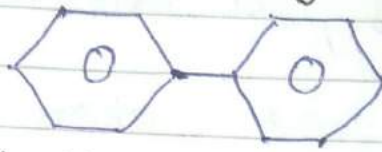
$$\text{Total \% of L} = 72 + 14 = 86\%$$

11

$$d = 14\%$$

Ques

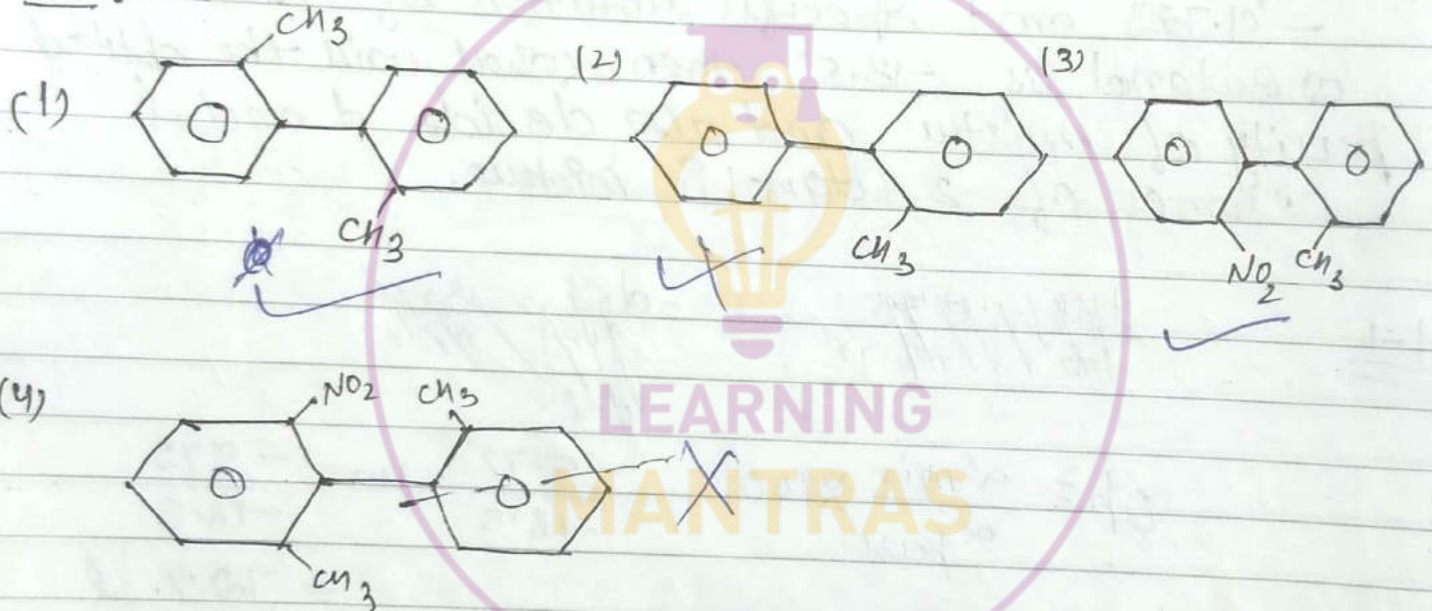
* Optical Activity in Biphenyl!



Condition: to be optically active.

- 1) Large gp at any/both ortho position to both ring.
- 2) No POS along Intermolecular bond in both ring.

Que:



Learning Mantras
Our Guidance, Your Success