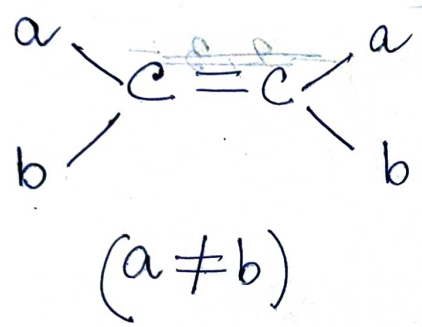


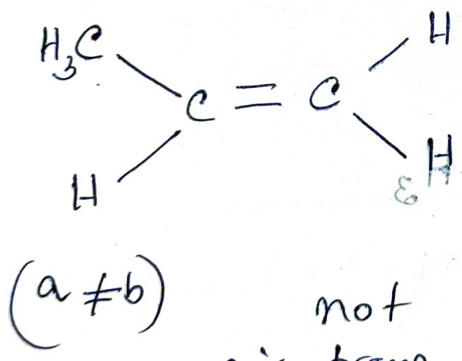
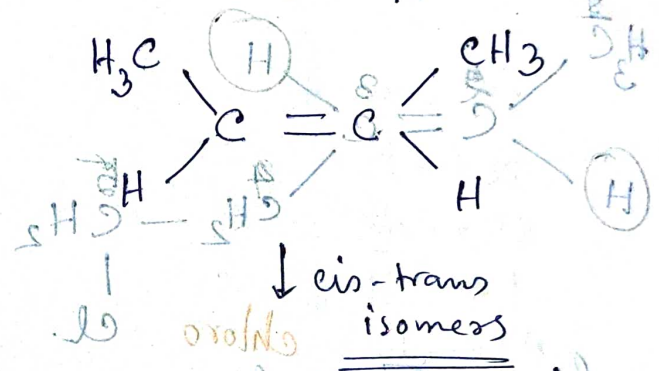
cis-trans isomerism (geometrical isomerism):

identical formula (not shape)
Isomers

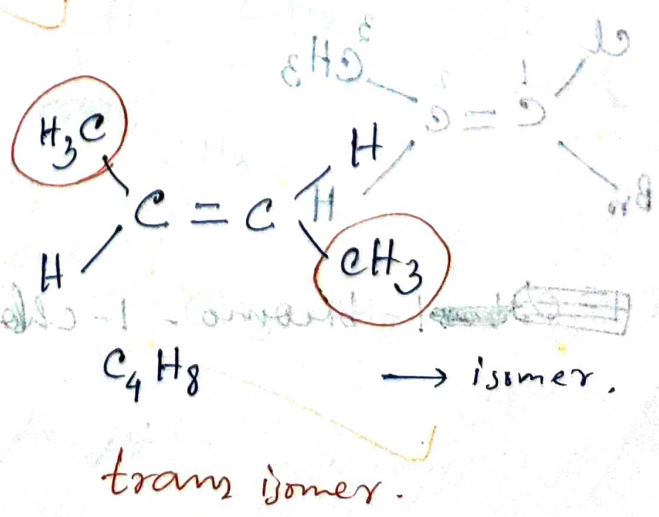
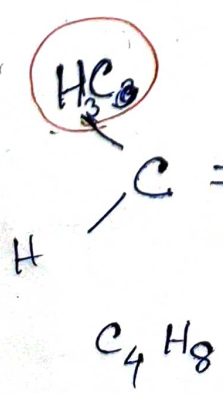
= (double bond)



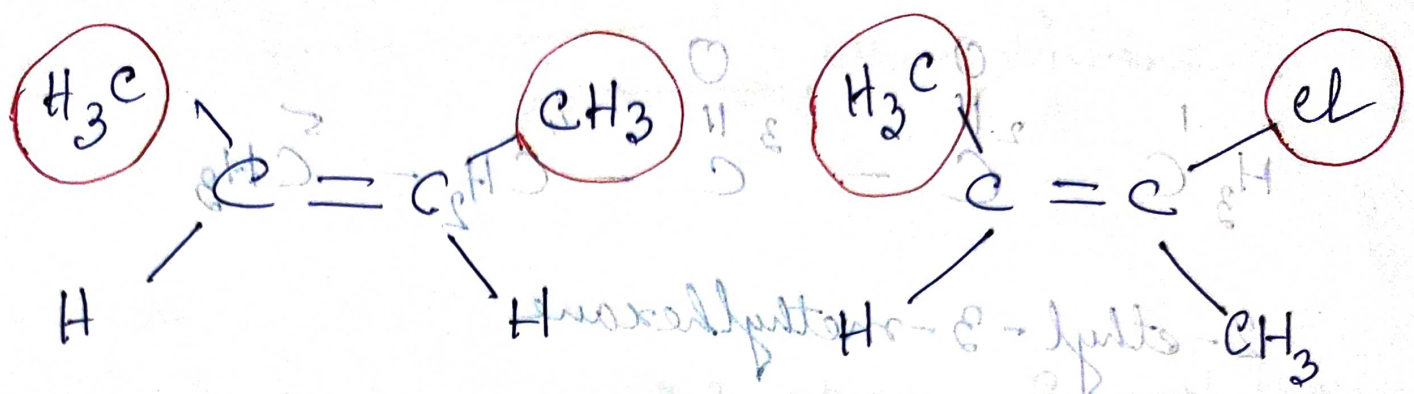
Let, a = CH₃
b = H



not cis-trans isomers.



E/Z Nomenclature

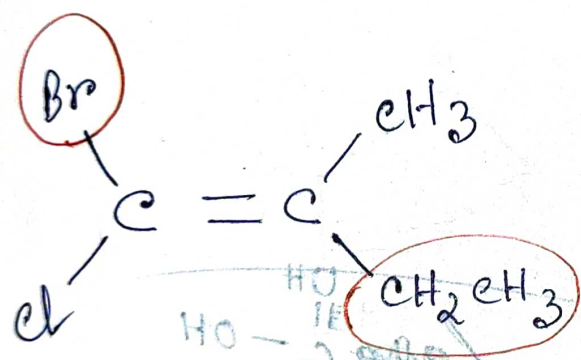


cis, Z

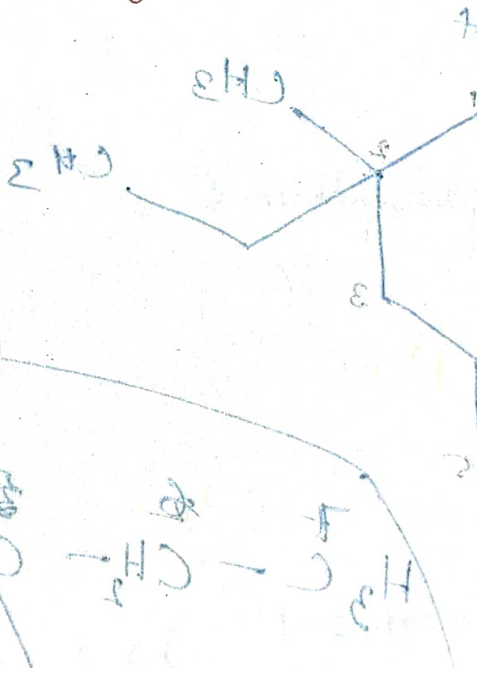
trans, E

Z

same group same side



E



- C → 12
- Cl → 35.5
- Na → 23
- H → 1
- N → 14
- O → 16
- Ca → 40

R-S Nomenclature:-

Rules:

(0) Nearer end of an axis on a plane precedes the farther end (proximity rule)

(1) Higher atomic number precedes lower. ~~eg.~~

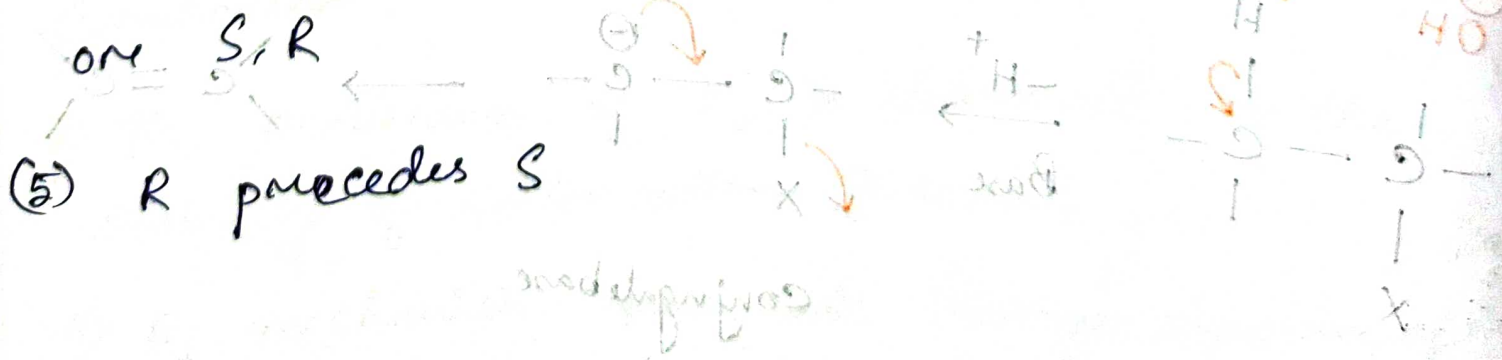
eg: $S > F > O > N > C > H$

(2) Higher atomic mass number precedes lower.

eg. $T > D > H$

(3) Cis precedes trans; and Z precedes E

(4) Like pairs RR or S, S. precedes unlike R, S



(5)

Atoms and groups with increasing priority:-

0. Lone pair

1. H

2. D

3. CH₃

4. CH₂CH₃

5. CH₂(CH₂)_nCH₃

6. CH₂-CH=CH₂

7. CH₂-C≡CH

8. CH₂C₆H₅

9. CH(CH₃)₂

10. CH=CH₂

11. C(CH₃)₃

12. C≡CH

13. C₆H₅

14. CH₂OH

15. CH=O

16. COR

17. CONH₂

18. CO₂H

19. CO₂R

20. NH₂

21. NHCH₃

22. N(CH₃)₂

23. NO

24. NO₂

25. OH

26. OCH₃

27. OC₆H₅

28. OCOR

29. F

30. SH

31. SR

32. SOR

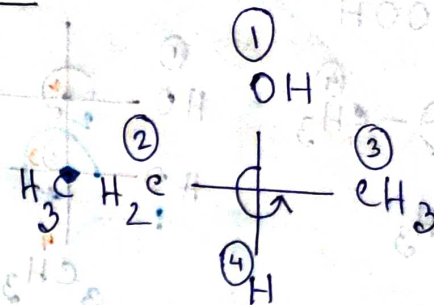
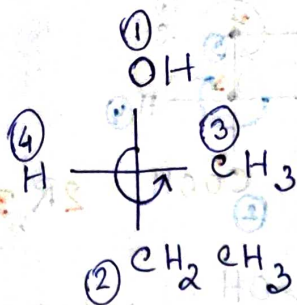
33. SO₂R

34. Cl

35. Br

36. I

R-S Nomenclature Example:



anticlockwise \rightarrow S

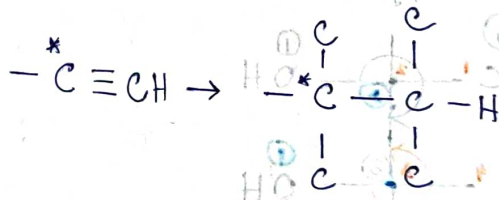
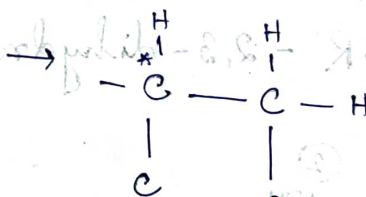
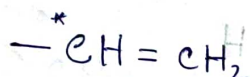
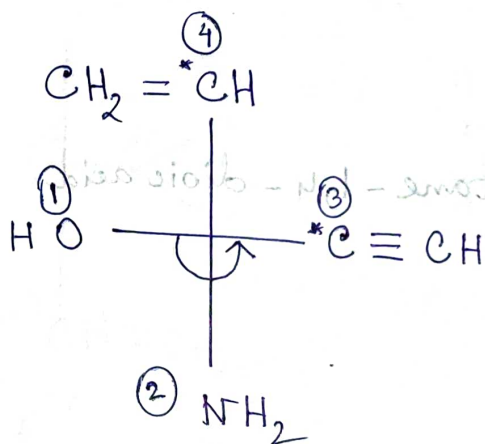
anticlockwise \rightarrow S

④ no. gr. \rightarrow horizontally present.

④ no. gr. \rightarrow vertical line.

R

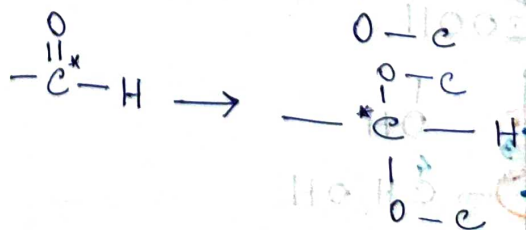
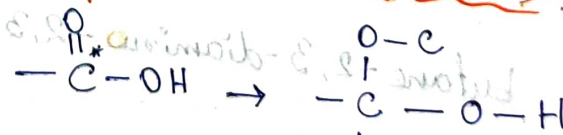
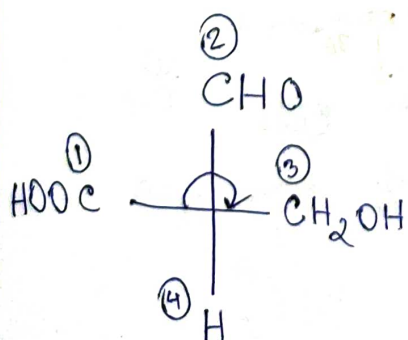
S



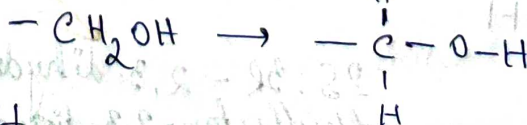
S

$(-\text{OH}) > -\text{NH}_2 > -\text{C}\equiv\text{CH} > -\text{CH}=\text{CH}_2$ Fantom atom.

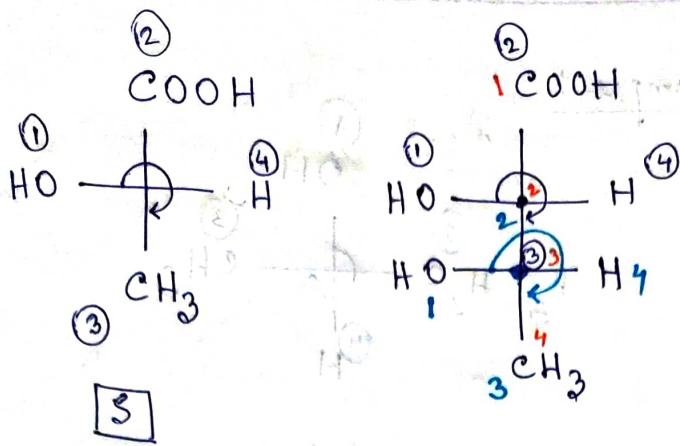
CIP Rule (Cahn-Ingold-Prelog Rule)



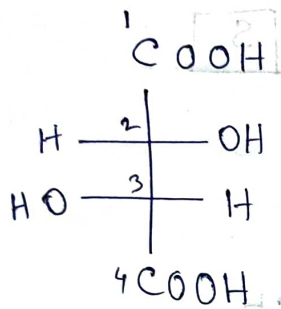
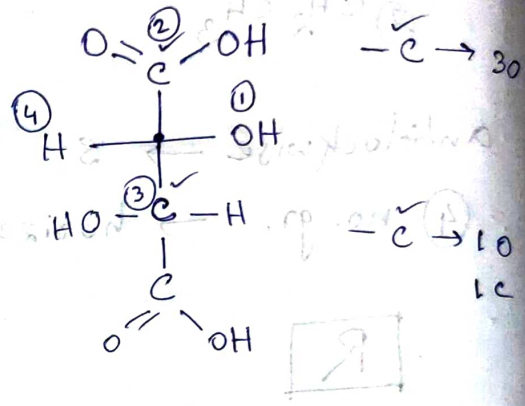
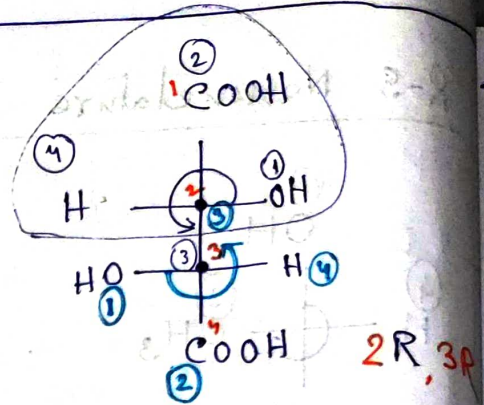
R



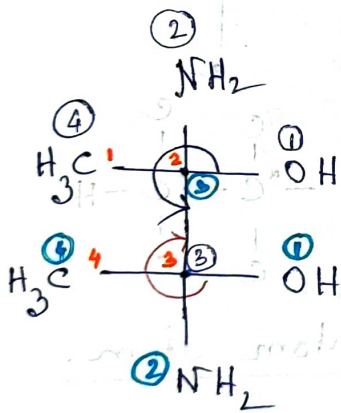
$(-\text{COOH}) > -\text{CHO} > -\text{CH}_2\text{OH} > \text{H}$



2S, 3S

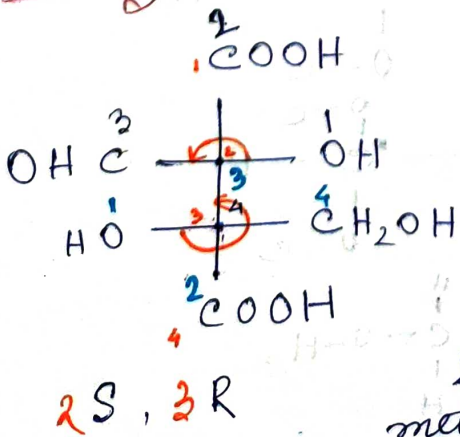


2R, 3R - 2,3-dihydroxybutane-1,4-dioic acid

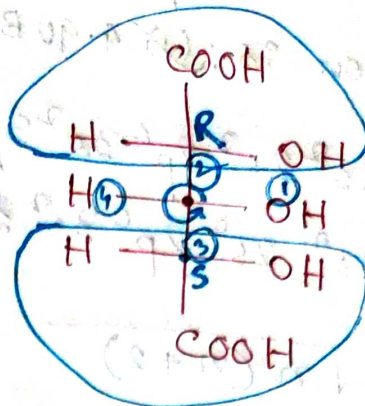
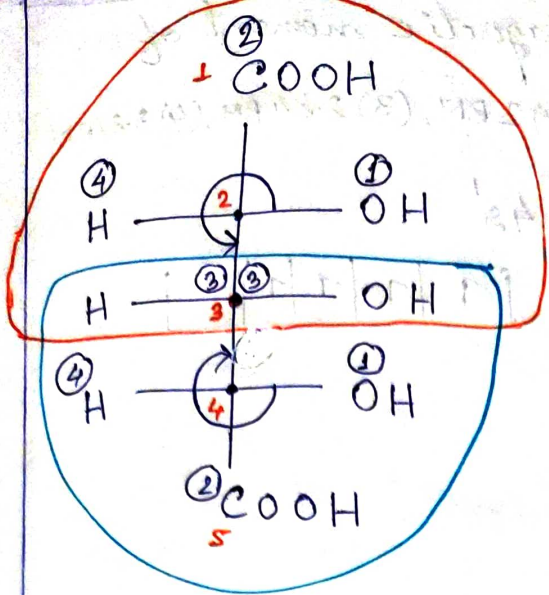


2R, 3R

2R, 3R - butane-2,3-diamine-2,3-diol

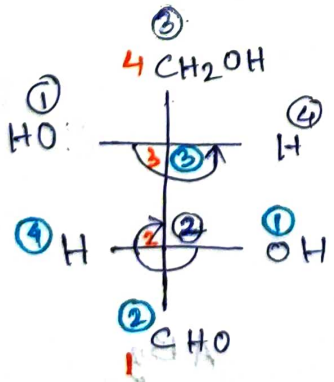
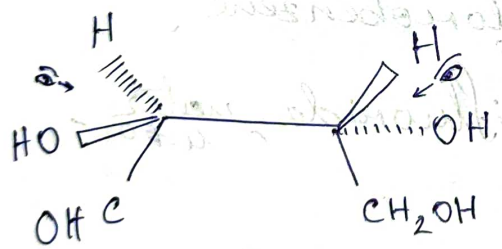
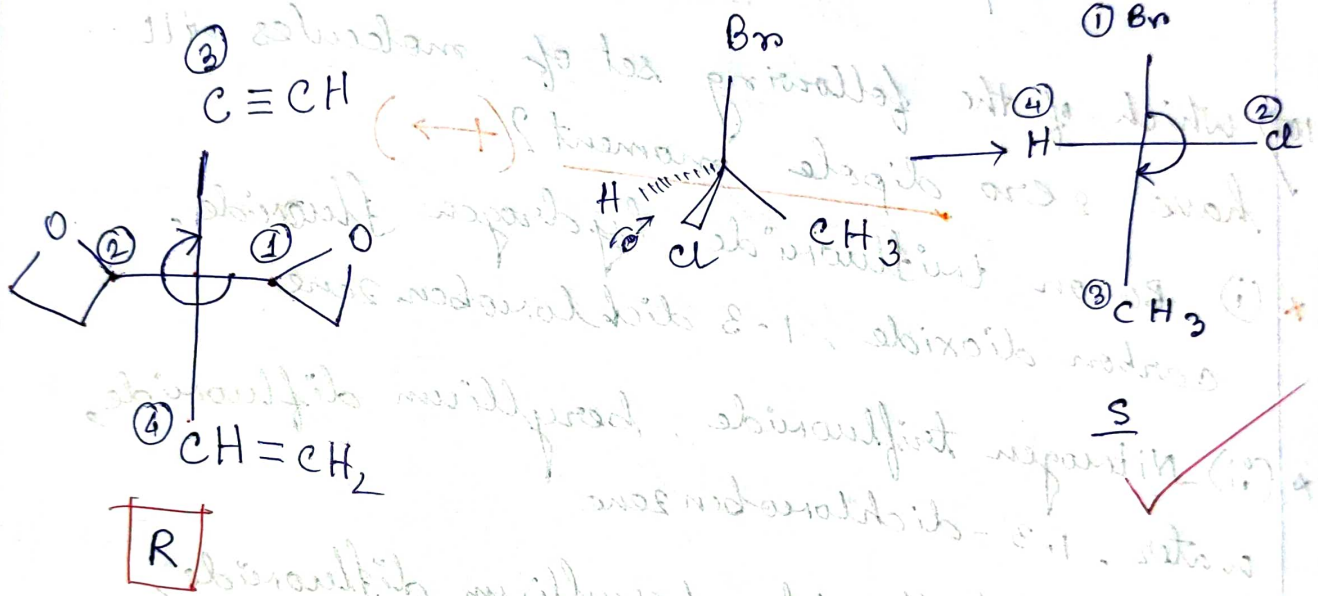


2S, 3R - 2,3-dihydroxy-2-formyl-3-hydroxy methylbutan-2,3-dioic acid



Syllabus
Full Organic,
Redox

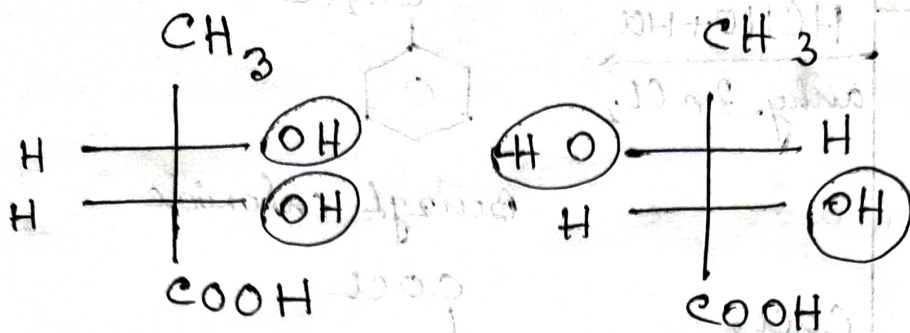
2R, 3R, 4S



~~3R~~, 2S, 3R

2S, 3R - 2, 3, 4-bis(hydroxy)butanal

Threo - Erythro nomenclature :



Erythro

(Same gr. \rightarrow \rightarrow
निचे \rightarrow \rightarrow)

Threo

(Same gr. opposite
 \leftarrow \rightarrow)

D/L nomenclature :

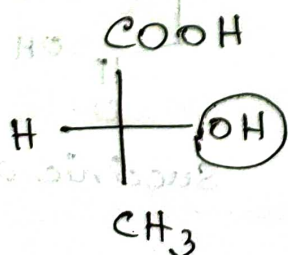


D-glyceraldehyde

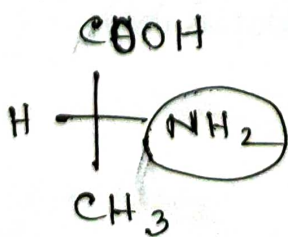
(D \rightarrow dextrorotatory)

L-glyceraldehyde

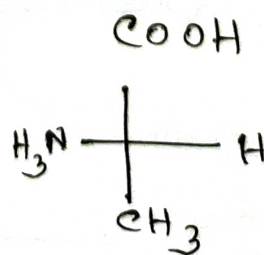
(L \rightarrow Levorotatory)



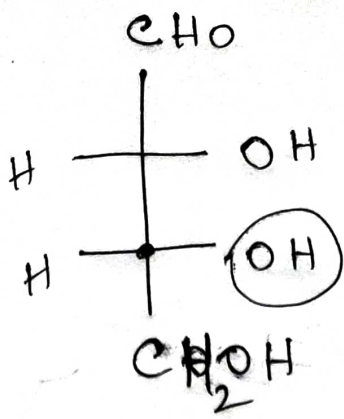
D-lactic acid.



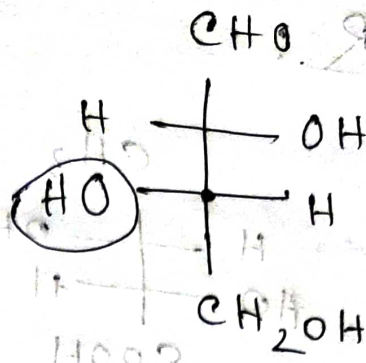
D-alanine



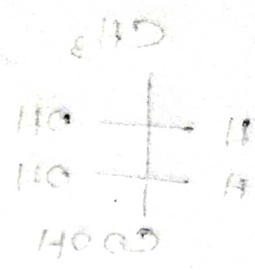
L-alanine



D

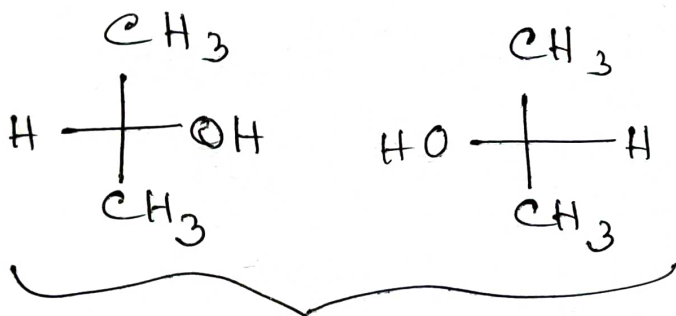


L



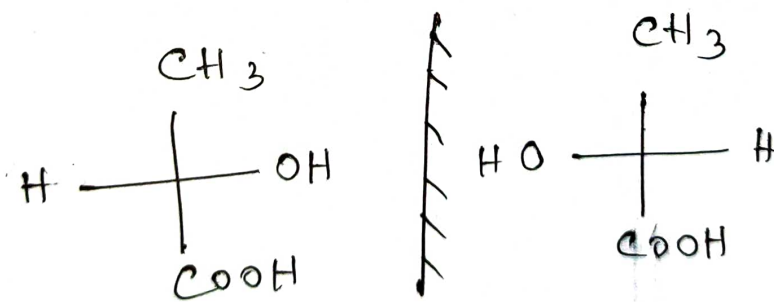
Two different compounds
 are not superimposable
 mirror images of each other

HOMOMER



Same structure.

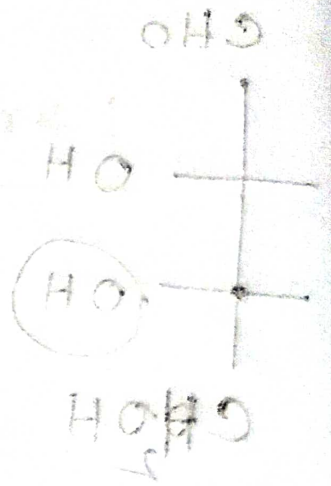
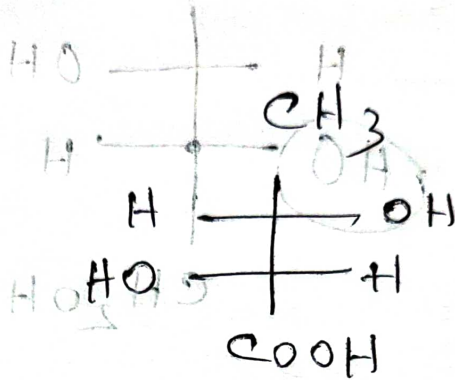
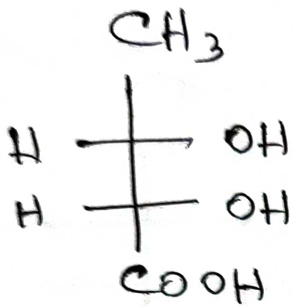
ENANTIOMER



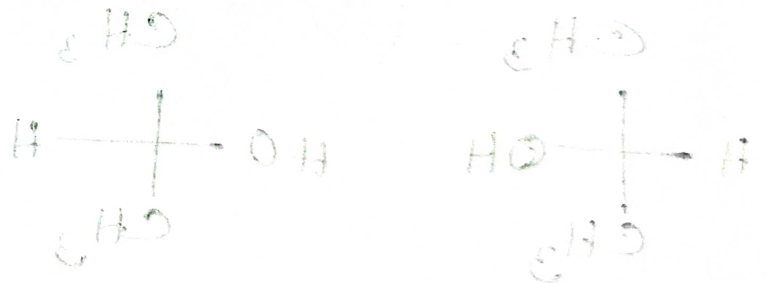
- i) Mirror image relationship
- ii) Superimposable structures.

↓
 Two kind of compound are possible for
 a chiral centre.

DIASTEREOMER



- i) No mirror image relationship.
- ii) Two different compounds
- iii) Non-superimposable structures.



Same structure.

ENANTIOMER

