

AULE GENEVIEVE MSURSTIMA

19/MT501/108

MBBS

CHEM 102

ASSIGNMENT ON STEREOCHEMISTRY AND FUNCTIONAL GROUP.

i) Name the functional groups present in each of the following molecules.

i) $\text{CH}_2=\text{C}(\text{OH})\text{CHO}$ - aldehyde, alkene

hydroxyl functional group, carbonyl functional group ($\text{C}=\text{O}$) but for aldehyde ($\begin{matrix} \text{R} \\ \diagdown \\ \text{C}=\text{O} \\ \diagup \\ \text{H} \end{matrix}$) and carbon-carbon ^{double} bonds (for alkenes).

ii) $\text{C}_6\text{H}_5\text{CH}(\text{NH}_2)\text{COCH}_3$

amine functional group, carbonyl functional group ($\text{C}=\text{O}$) but for ketone

($\begin{matrix} \text{R} \\ \diagdown \\ \text{C}=\text{O} \\ \diagup \\ \text{R}' \end{matrix}$), carbon-carbon single bond (for alkanes)

iii) $\text{CH}_3\text{C}=\text{CHCH}(\text{OH})\text{CHO}$

hydroxyl functional group, carbonyl functional group ($\text{C}=\text{O}$) but the one for aldehyde ($\begin{matrix} \text{R} \\ \diagdown \\ \text{C}=\text{O} \\ \diagup \\ \text{H} \end{matrix}$) and carbon-carbon double bond (for alkenes)

2) A 0.856g sample of pure (2R, 3R)-tartaric acid was diluted to 10cm^3 with water and placed in a 1.0dm polarimeter tube, the observed rotation at 20°C was $+1.0^\circ$. Calculate the specific rotation of (2R, 3R)-tartaric acid

SOLUTION

$$\text{Specific rotation} = \frac{\text{observed rotation (degrees)}}{\text{concentration } \text{g cm}^{-3} \times \text{path length of sample cell in dm}}$$

$$\text{observed rotation} = +1.0^\circ$$

$$\text{conc. in } \text{g cm}^{-3} = \frac{\text{mass (g)}}{\text{volume } \text{cm}^3} = \frac{0.856\text{g}}{10\text{cm}^3} = 0.0856\text{g cm}^{-3}$$

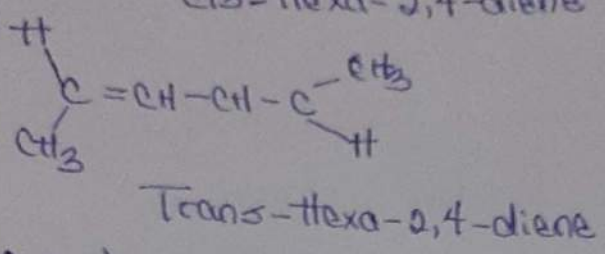
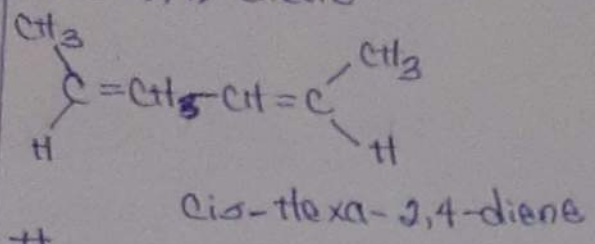
$$\text{path length of sample cell} = 1.0\text{dm}$$

$$\text{specific rotation} = \left(\frac{+1.0^\circ}{(0.0856 \times 1.0)} \right) \text{g}^{-1}\text{cm}^2\text{dm}^{-1}$$

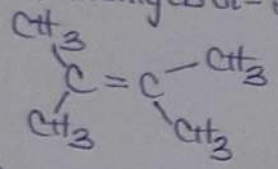
$$= 11.68^\circ\text{g}^{-1}\text{cm}^2\text{dm}^{-1}$$

3) Draw the possible geometric isomers (where possible) for each of the following compounds.

i) Hexa-2,4-diene



ii) 2,3-dimethylbut-2-ene



Geometric isomers is not possible for 2,3-dimethylbut-2-ene