

NAME: UCHENDU ANITA CHIZARAM

MATRIC NO: 19/MHS01/408

DEPARTMENT: MEDICINE AND SURGERY

COURSE CODE: CHM102

ASSIGNMENT

b NAME: UCHENDU ANITA CHIZARAM
MATIC NO. 1711115011408
DEPARTMENT: MEDICINE AND SURGERY
COURSE CODE: CHM 102

ASSIGNMENT

1. Name the functional groups present in each of the following molecules

- $\text{CH}_2 = \text{C}(\text{OH})\text{HCHO}$
- $\text{C}_6\text{H}_5\text{CH}(\text{NH}_2)\text{COCH}_3$
- $\text{CH}_3\text{C} = \text{CHCH}(\text{OH})\text{CHO}$

ANSWER

MOLECULES	FUNCTIONAL GROUPS
$\text{CH}_2 = \text{C}(\text{OH})\text{HCHO}$	- Aldehyde (-CHO) - Hydroxyl group (-OH) - Alkene
$\text{C}_6\text{H}_5\text{CH}(\text{NH}_2)\text{COCH}_3$	- Carbonyl group (-CO) - Amine (-NH ₂)
$\text{CH}_3\text{C} = \text{CHCH}(\text{OH})\text{CHO}$	- Hydroxyl group (-OH) - Aldehyde (-CHO)

2. A 0.856g sample of pure (2R,3R)-tartaric acid was diluted to 10cm³ with water and placed in a 1.0dm polarimeter tube. This observed rotation at 20°C was +1.0°. Calculate the specific rotation of (2R,3R)-tartaric acid

ANSWER

$$[\alpha] = \frac{\alpha}{cL}$$

Where $[\alpha]$ is specific optical rotation

α observed rotation
 c concentration in g cm⁻³
 L pathlength in (dm)

$$\alpha = +1.0^\circ$$
$$L = 1.0 \text{ dm}$$
$$c = ?$$

b)

$$\text{concentration in } \text{g cm}^{-3} = \frac{0.886 \text{ g}}{10 \text{ cm}^3} = 0.0886 \text{ g cm}^{-3}$$

2)

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

$$[\alpha]_{\lambda}^T = \frac{\alpha}{c \cdot l}$$

$$[\alpha]_{\lambda}^T = \frac{+1.0^{\circ}}{0.0886 \text{ g cm}^{-3} \times 1.0 \text{ dm}} = +11.682^{\circ} \text{ g}^{-1} \text{ cm}^3 \text{ dm}^{-1}$$

or $11.7^{\circ} \text{ g}^{-1} \text{ cm}^3 \text{ dm}^{-1}$

T

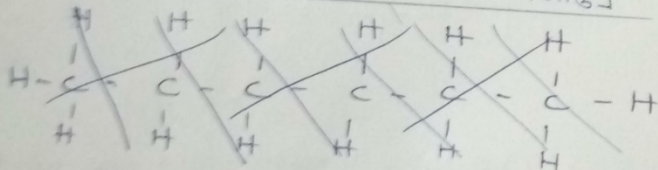
Therefore the specific rotation of (2R,3R)-tartaric acid is $11.682^{\circ} \text{ g}^{-1} \text{ cm}^3 \text{ dm}^{-1}$

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

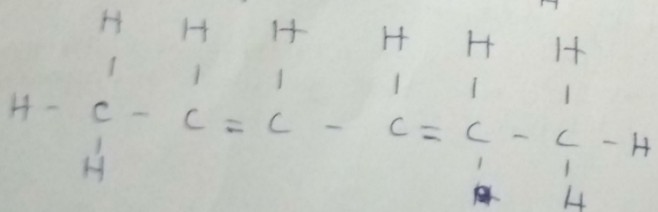
- i) Hexa-2,4-diene
- ii) 2,3-Dimethylbut-2-ene

POSSIBLE ISOMERS

Hexa-2,4-diene $[\text{CH}_3(\text{CH}=\text{CH}=\text{CHCH}_3)]$

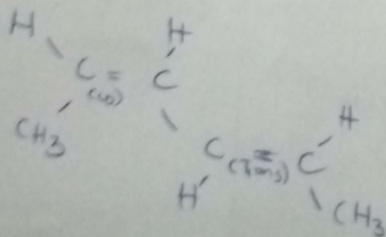


i)



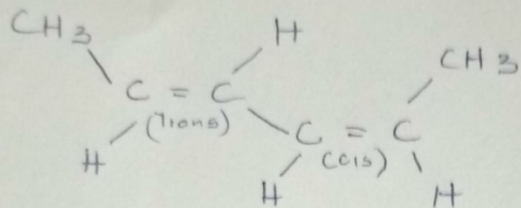
Hexa-2,4-diene

ii)

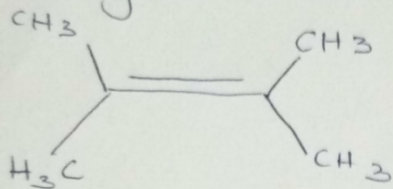


(cis-), trans-2,3-dimethylbut-2-ene

b) Trans-1, cis-4-dimethylbut-2-ene



a) 2,3-Dimethylbut-2-ene



There are no isomers for the compound.

