



AAH-003-001607 Seat No. _____

B. Sc. (Sem. VI) (CBCS) Examination

April/May - 2016

C-602 : Organic Chemistry & Spectroscopy

(New Course)

Faculty Code : 003

Subject Code : 001607

Time : $2\frac{1}{2}$ Hours]

[Total Marks : 70

- Instructions :** (1) All questions are compulsory.
(2) Answer of Q. 1 (MCQ) should be written in the answer book.

1 Answer the following : 20

- (1) Citral is extracted from
(A) lemon grass oil (B) jasmine oil
(C) camphor oil (D) peppermint oil
- (2) Biuret test is utilised in the identification of
(A) terpenoid (B) protein
(C) alkaloid (D) dyes
- (3) Citral on heating with KHSO_4 gives
(A) p-Cymene (B) Geraniol
(C) Geranic acid (D) Laevulic acid
- (4) Chlorination of diphenylmethane in presence of sunlight gives
(A) 4,4' -Dichloro diphenylmethane
(B) Diphenyldichbro methane
(C) Chlorobiphenyl methane
(D) None of these

- (5) Staggered conformer of ethane is having energy
- Lower than eclipsed conformer
 - Higher than eclipsed conformer
 - Equal to eclipsed conformer
 - Both (A) and (B)
- (6) Citral is oxidised with silver oxide to obtain
- Laevullinic acid
 - Geraniol
 - Geranic acid
 - p-Cymene
- (7) Baygon is used as
- Perfume
 - Explosive
 - Insecticide
 - Protien
- (8) Diphenylmethane is oxidised with chromic acid to yield
- Benzhydrol
 - Dibenzyl
 - Benzophenone
 - Fluorene
- (9) α -Terpeniol reacts with Tilden's reagent to give compound molecular formula is
- $C_{10}H_{17}O_2NCl$
 - $C_9H_{17}O_2NCl$
 - $C_{10}H_{18}O_2NCl$
 - $C_9H_{18}O_2NCl$
- (10) In n-Butane conformation analysis after the rotation of 180° to obtain the form of
- Eclipsed
 - Gauche - staggered
 - Antistaggered
 - None of these
- (11) How many signals in NMR spectra will be obtained from the following compounds respectively?
- $CH_3CH_2CHBr_2$ - 1, 1 - Dibromopropane
 - $CH_3CHBrCH_2Br$ - 1, 2 - Dibromopropane
 - $CH_3CBr_2CH_3$ - 2, 2 - Dibromopropane
 - $CH_2BrCH_2CH_2Br$ - 1, 3 - Dibromopropane
- (3), (4), (1), (2)
 - (3), (4), (2), (1)
 - (3), (2), (1), (4)
 - (4), (3), (1), (2)

- (12) In mass spectroscopy, the stability of removed cation is
- (A) $3^\circ > 2^\circ > 1^\circ$ (B) $1^\circ > 2^\circ > 3^\circ$
 (C) $1^\circ = 2^\circ = 3^\circ$ (D) $1^\circ = 2^\circ > 3^\circ$
- (13) Analysis of an organic compound shows it to be 64.3% carbon. It displays a molecular ion at $m/z = 112$ amu in the mass spectrum. Which of the following is plausible molecular formula for this compound ? Which factor can affect chemical shift ?
- (A) C_8H_{16} (B) $C_7H_{12}O$
 (C) $C_6H_8O_2$ (D) $C_5H_4O_3$
- (14) Nitrile group peak is identified from
- (A) NMR spectra (B) IR spectra
 (C) UV spectra (D) TLC
- (15) How Many types of hydrogen are present in the 1-Bromopropane ?
- (A) Two (B) Three
 (C) Four (D) One
- (16) Mc – Lafferty rearrangement requires
- (A) Minimum three carbon in chain
 (B) Minimum two carbon in chain
 (C) Minimum one carbon in chain
 (D) None of these
- (17) In mass spectroscopy, all fragment possess
- (A) Same m/e value (B) Different M/e value
 (C) Equal mass value (D) Equal δ value
- (18) The peak having highest intensity is known as
- (A) Metastable ion peak (B) Parent peak
 (C) Molecular ion peak (D) Base peak

- (19) The NMR spectrum of diethyl ether shows
- (A) Two Peaks, both quartet
 - (B) Two Peaks, one a triplet, the other a quartet
 - (C) Two Peaks, both triplet
 - (D) Two Peaks, one a triplet, the other a doublet
- (20) Carbonyl group peak is identified from
- (A) TLC
 - (B) NMR spectra
 - (C) UV spectra
 - (D) IR spectra

2 (a) Answer the following : (any **three**) **6**

- (i) Give the synthesis of Naphthalene by Haworth method.
- (ii) Define :
 - (1) Proteins
 - (2) Polynuclear hydrocarbons
- (iii) Give the synthesis of Citral from Geranic acid.
- (iv) How will you determine the presence of $> C = O$ group in terpenoids ?
- (v) Give synthesis and use of P.E.T.N.
- (vi) Give synthesis and use of Musk ketone.

(b) Answer the following : (any **three**) **9**

- (i) Write a short note on classification of proteins.
- (ii) Give synthesis methods of Biphenyl.
- (iii) Prove that : Citral is an unsaturated aldehyde.
- (iv) Prove that : α – terpineol contains –OH alcoholic group at the 8th position.
- (v) Discuss the isoelectric point of an amino acid.
- (vi) Give synthesis of α – terpineol.

(c) Answer the following : (any two) 10

- (i) Give synthesis and uses of :
 - (1) RDX
 - (2) Carbendazin.
- (ii) Explain chemical properties of Anthracene.
- (iii) Discuss conformation of n-butane with energy diagram.
- (iv) Write any three methods for preparation of polypeptides.
- (v) Explain relative energies and the stability of the various conformations of cyclohexane.

3 (a) Answer the following : (any **three**) 6

- (i) Define Equivalent and Non-equivalent protons.
- (ii) Define Base peak and Molecular ion peak.
- (iii) Give the structure for the compound $C_4H_6O_3$ giving only one NMR signal.
- (iv) How many signals would you expect in NMR spectrum of Trans-1,2-Dimethyl cyclopropane ?
- (v) Give the structure for the compound $C_8H_{18}O$ giving only one NMR signal.
- (vi) How will you distinguish the phenyl acetic acid and methyl benzoate with the help of NMR spectrum?

(b) Answer the following : (any **three**) 9

- (i) Discuss importance of TMS in NMR spectroscopy.
- (ii) Explain shortly : "Principle of mass spectroscopy".
- (iii) How many kinds of hydrogen are present in the following compound ?
 - (A) $C_6H_5-CH_3$
 - (B) $CH_2=CH_2$
 - (C) $CH_3-CH=CH_2$
- (iv) Sketch the NMR spectrum of Ethyl benzoate.
- (v) Explain Mc-Lafferty rearrangement.

- (vi) Assign the structure to a compound having following characteristics.

M.F. : $C_5H_6O_3$

IR : 2965, 1825, 1780, 1280 cm^{-1}

NMR : (a) Quintet 2H δ ppm = 1.8

(B) Triplet 4H δ ppm = 2.4

- (c) Answer the following : (any two) : 10

- (i) Explain the factors affecting chemical shift in NMR Spectrum.
- (ii) Explain characteristics of Mass spectra of alkane.
- (iii) Explain with suitable example "Enantiomeric proton and Diastereomeric proton".
- (iv) Assign the structure to a compound from the following spectral results with explanation.

M.F. : $C_8H_{14}O_4$

IR : 2990, 2885, 1730, 1035 cm^{-1}

NMR : (a) Triplet 6H δ ppm=1.85

(B) Singlet 4H δ ppm=2.0

(C) Quartet 4H δ ppm=4.3

- (v) Assign the structure to a compound from the following spectral results with explanation.

Molecular weight : 199 gm/mol [C=54.27%, H=5.53%, Br=40.2% (At. Wt. of Br=80 gm/mol)]

IR : 3025, 2910, 1600, 1580, 1505, 1440, 750, 710 cm^{-1}

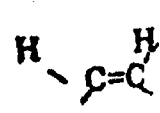
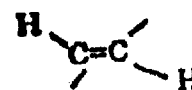
NMR : (a) Singlet 5H δ ppm=7.22

(B) Triplet 2H δ ppm=2.74

(C) Triplet 2H δ ppm=3.38

(D) Quintet 2H δ ppm=1.85

Spectral Data

Infra - Red Data		
Alkene (stretching)	-C-H	2850-2960(v)
Alkene	=C-H	3100-3200(m)
Alkyene	=C-H	3200-3300(s)
Aromatic	ArC-H	3010-3100(m)
Aromatic ring	C=C	1500-1600(v) (two to three)
Alkene	>C=C<	1610-1680(v)
Alkyene	-C=C ² .	2100-2260(s)
Alkene (Bending)	-C-H	1340(w)
	-C(C ₂ H ₃) ₃	1430-1470(m) & 1380-1385(s)
	-C(CH ₂) ₃	1365 (s)
Aldehyde	-C-H	2820-2000(w)&2850 2760(s)
Aldehyde	C=O	1740-1720(s)
Ketone	C=O	1725-1710(s)
Carboxylic acid	C=O	1725-1705(s)
Ester	C=O	1750-1730(s)
Amide	C=O	1670-1640(s)
Anhydride	C=O	1810-1860(s)&1740-1790
Alcohols, Ethers, esters		
Carboxylic acids, Anhydride	C-O	1300-1000(s)
Alcohols, phenols :		
Free	O-H	3650-3600(sh)
bonded	O-H	3500-3200(b)
Carboxylic acids free		
Free	O-H	3500-3650(m)
H-bonded	O-H	2500-3200(b)
amines (stretch)	N-H	3330-3500(m)
Bnding	-N-H	1640-1550(m)
Nitrile	-C=N	2210-2280(s)
Ether	-O-	1070-1150(s)
Alkene bending		-690(s)
disubstituted Cis.		
disubstituted Trans.		960-970(s)
Aromatic substitution :		
Type C-H out of plane bending		
No. of adjacent H atom.		range cm
5		750(s) & 700(s)
4		750
3		780
2		830
1		850

NMR Data : Chemical Shift

Types of proton		Chemical shift in δ_{ppm}
Primary	$R-CH_3$	0.9
Secondary	R_2-CH_2	1.3
Tertiary	R_3-CH	1.5
Vinylic	$C=C-H$	4.6-5.9
Acetylinic	$Cr-C-H$	2.3
Aromatic	$Ar-H$	6-8.5
Benzylic	$Ar-C-H$	2.2-3
Allylic	$C=C-CH_3$	1.7
Fluorides	$H-C-F$	4-4.5
Chlorides	$HC-Cl$	3.4
Bromides	$HC-Br$	2.5-4
Iodides	$HC-I$	2.4
Alcohols	$HC-OH$	3.4-4
Ethers	$HC-OR$	3.3-4
Esters	$R-COO-CH$	3.7-4.1
Acids	$HC-COOH$	2-2.6
Carbonyl comp.	$HC-C=O$	2-2.7
Aldehyde	$R-CHO$	9-10
Hydroxylic	$R-OH$	1-5.5
Phenolic	$Ar-OH$	4-12
Carboxylic	$R-COOH$	10.5-12
Amino	$R-NH_2$	1.5