

NAN-003-001607 Seat No. \_\_\_\_\_

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

March / April - 2017

C-602 : Organic Chemistry & Spectroscopy (New Course)

Faculty Code : 003 Subject Code : 001607

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

[Total Marks: 70

**Instructions**: (1) Total three questions, all are compulsory.

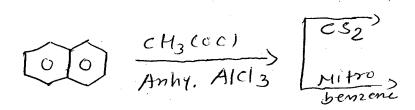
- (2) Q. 1 carries 20 marks, should be written in the answer book.
- 1 Answer the following questions:

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- (1) Write the structure of Isoprene.
- (2) Give full name of RDX and TNT.
- (3) Define Amino acids.
- (4) Write Ninhydrine Test.
- (5) Draw chair form of Cyclohexane.
- (6) Complete the reaction:

$$C_{10}H_{16}O \xrightarrow{\phantom{-}(i)\,KMNO_4\phantom{}}\phantom{-}(ii)\,CrO_4\phantom{}$$

- (7) Give an example and structure of basic amino acid.
- (8) Give structure and uses of PARATHION.
- (9) Complete the reaction (Friedal Craft acylation)



- (10) Give an example of Ullmann reaction.
- (11) What is chemical shift in NMR?
- (12) Define equivalent and non-equivalent protons.
- (13) How many types of Hydrogen are present in Methyl ethyl ether?
- (14) Give structure formula of compound which give rise only one NMR signal
  - (i)  $C_5H_8Cl_4$
- (15) Give structure formula which give only two NMR signal
  - (i)  $C_3H_8O_2$
- (16) >C = O group identify by which spectra?
- (17) Define Base Peak.
- (18) How many minimum carbon required for Mc-Lafferty rearrangement?
- (19) Which information we get from IR and NMR spectra?
- (20) Give no. of signal and splitting of each signal of  $\mathrm{CH_3-CH=CH_2}$
- 2 (a) Answer any three from the following:
  - (1) Give synthesis of Phenyl alanine from Benzaldehyde.
  - (2) Give synthesis of Geranic acid.
  - (3) Give synthesis of Glycylalanine by Bergman method.
  - (4) Complete the Anthrace reaction with H<sub>2</sub>/Ni
  - (5) Give synthesis and uses of P.E.T.N.
  - (6) Give synthesis and use of Musk Ketone.
  - (b) Answer any three from the following:
    - (1) Give synthesis and uses of Baygon and Biphenyl.
    - (2) Give synthesis and uses of RDX.
    - (3) Write a short note on classification of Proteins.

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- (4) Give conversion of Terebic acid from  $\alpha$ -terpeniol.
- (5) Give synthesis of Naphthalene by Haworth Reaction.
- (6) Give synthesis of  $\alpha$ -Terpineol.
- (c) Answer any two from the following:

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- (1) Explain constitution of Citral.
- (2) Explain chemical properties of Antracene.
- (3) What is Polypeptide? Give synthesis of Thyroxine.
- (4) Explain conformational analysis of cyclohexane.
- (5) Give synthesis of:
  - (i) Baygon
  - (ii) Biphenyl.
- 3 (a) Answer any three from the following:

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- (1) Why TMS is used as internal standard?
- (2) Give structure formula of following compound which give rise only two NMR signal.
  - (i)  $C_4H_6O_4$
  - (ii)  $C_2H_3F_3O$ .
- (3) Which Nuclei give NMR spectra?
- (4) A compound give peak at 1250 Hz. If NMR spectra taken in 100 MHz. Calculate chemical shift in  $\delta_{ppm}$  and  $\tau_{ppm}$ .
- (5) Explain the principle of Mass spectroscopy.
- (6) How many signals would you expect in NMR spectrum of Trans-1,2-Dimethyl cyclopropane?
- (b) Answer any three from the following:

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- (1) Distinguish chemical shift  $\delta$  and coupling constant "J".
- (2) Give possible isomer of dibromo propane and give no. of NMR signals.
- (3) Explain Deuterium labeling.
- (4) What is isotropic and anisotropic proton? Explain in short with example.
- (5) Distinguish the following compound by NMR spectra: (i) C<sub>6</sub>H<sub>5</sub>COCH<sub>3</sub> and (ii) C<sub>6</sub>H<sub>5</sub>CH<sub>2</sub>CHO

(6) Derive the structure and compound from given spectral information:

Mole. Formula : 
$$C_9H_{10}$$

NMR : a Quintet 2.04 
$$\delta_{ppm}$$
 2H

b Triplet 2.91 
$$\delta_{ppm}$$
 4H

c Singlet 7.17 
$$\delta_{ppm}$$
 4H

(c) Answer any two from the following:

- **10**
- (1) Explain Mass instrumentation with diagram.
- (2) Explain factors Effecting on chemical shift.
- (3) Assign the structure from the following data : M.F. =  $C_6H_{10}O_3$ .

$$IR = 3400-3500 \text{ cm}^{-1}, 2990, 2810, 1725, 1025 \text{ cm}^{-1}.$$

NMR = a Triplet 
$$\delta_{ppm} = 1.0$$
 3H

b Singlet 
$$\delta_{ppm} = 2.1$$
 3H

c Quartet 
$$\delta_{ppm} = 2.5$$
 2H

d Singlet 
$$\delta_{ppm} = 2.4$$
 1H

e Singlet 
$$\delta_{ppm} = 3.2$$
 1H

- (4) Assign the structure from the following data with explanation.
  - (i) Molecular weight = 174 g/mole.

(ii) 
$$C = 55.1\%, H = 8.0\%$$

(iii) IR = 
$$2941 \text{ cm}^{-1}$$
,  $2850$ ,  $1740$ ,  $1450$ ,  $1020 \text{ cm}^{-1}$ 

Signal	Splitting	$\delta_{ppm}$	Intensity
a	Quartet	2.61	8.2 square
b	Singlet	4.14	8.1 square
c	Triplet	1.27	12.4 square

- (5) Assign the structure formula from the spectral data.
  - (i) Vapour density : 51 gram/mole,M.W. = 102 gram/mole
  - (ii) C = 70.5%, H = 13.72%

IR = 3500 (s), 2950(s), 2900(m), 2850(m), 1452(sh), 1385(sh), 1080(s) cm<sup>-1</sup>.

NMR =	Signal	Splitting	$\tau_{ppm}$	Intensity
	a	Triplet	8.8	6H
	b	Quintet	8.5	$4\mathrm{H}$
	c	Multiplet	8.3	1H
	d	Doublet	6.1	2H
	e	Singlet	5.5	1H

Spectral Ba	ta	
Infra - Red Data		
Alkene (streteching)	-C-H	2850-2960(v)
Alkene	-0-11 =C-H	3100-9200(m)
	=C-H	320 <b>0-3</b> 300(s)
Alkyene Aromatic	ArC-H	3010-3100(m)
Aromatic ring	C=C	1500-1600(y)
Aromauc img	0-0	(two to thre
Alkene	>C=C<	1610-1680(v)
the contract of the contract o	_C=C <sup>2</sup> .	
Alkyene		2100-2260(s)
Alkene (Bending)	-C-H	1340(w)
*	$-C(C_2H_3)_3$	1430-1470(m)
Contract to the second of the	0.40**	1380-1385(s)
	$-C(CH_2)_3$	1365 (8)
Aldehyde		2820-2000(w)&2650 2760
Adehyde	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)&1740-17
Alecohols, 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	0.	1070-1150(s)
Alkene bending H		-690(s)
disulstituted Cis.	<b>:Q</b> )	333(3)
Alkene bending disulstituted Cis.  H C=C		
disulstituted Trans.	?_ <sub>н</sub>	960-970(s)
Aromatic substitution:		
Type C-H out of plane bending	•	
No of adiabant II at		
Tio. or adjacent 11 awn.	orthosubs	range cm
<u> </u>		750(s) & 700
* * * * * * * * * * * * * * * * * * *		750 790
		780
2		830

NMR Data: Cl		
ypes of proton	C	hemical shift in $\delta_{ppm}$
Primary	R-CH <sub>3.</sub>	0.9
Secondary	$R_2$ - $CH_3$	1.3
Tertiary	R <sub>3</sub> -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
Florides	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
Adehyde	R-CHO	9-10
Hydroxylic	R-OH	<b>.</b> 1-5.5
Phenolic	Ar-OH	4-12
Carboxylic	R-COOH	10.5-12
Amino	$R-NH_2$	1.5