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Reg. No.

THE OPEN UNIVERSITY OF SRI LANKA
B.Sc. Degree Programme and
Stand Alone Courses in Science - 2015/2016
CMU2221/CME4221 - Organic Chemistry 1
CONTINUOUS ASSESSMENT TEST 1

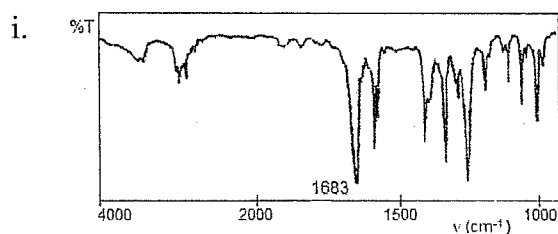
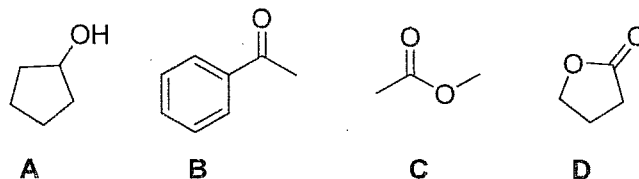
| Ques No. | Max. | Marks |
|--------------|------|-------|
| 1 | 30 | |
| 2 | 20 | |
| 3 | 20 | |
| 4 | 30 | |
| Total | 100 | |

Date: Saturday, 9th April 2016

Time: 14.30 - 15.30 hrs.

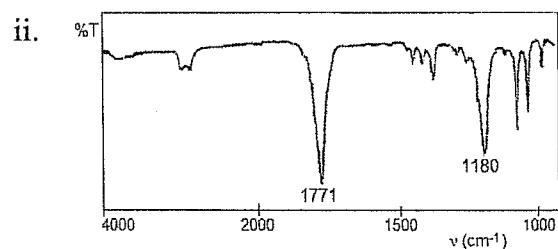
Write answers only in the space provided.

1. Giving reasons assign the correct compound out of the following (A-D) to its IR spectrum.



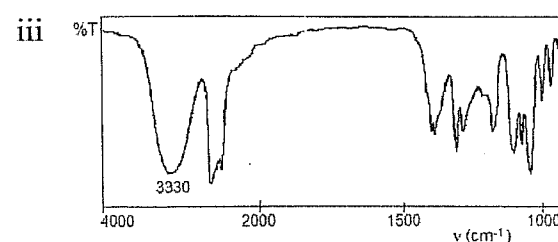
Compound

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Compound

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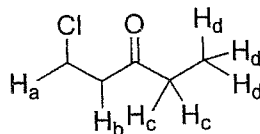
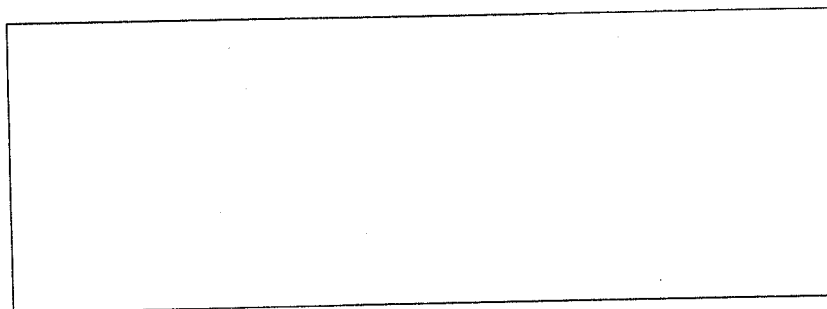
Compound

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(30 marks)

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2. Draw the ^1H NMR spectrum you would expect for the compound **E** showing multiplicities and relative positions of the signals from TMS (δ values not necessary). Label all the peaks.

Compound **E**

chemical shift

TMS

(20 marks)

3. CH_3CHO shows a UV absorption at 190 nm ($\log \epsilon = 2.0$) and at 290 nm ($\log \epsilon = 1.0$).

i. Indicate the electronic transitions responsible for the above absorptions.

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ii. Explain the observed differences in wave lengths of these two absorptions.

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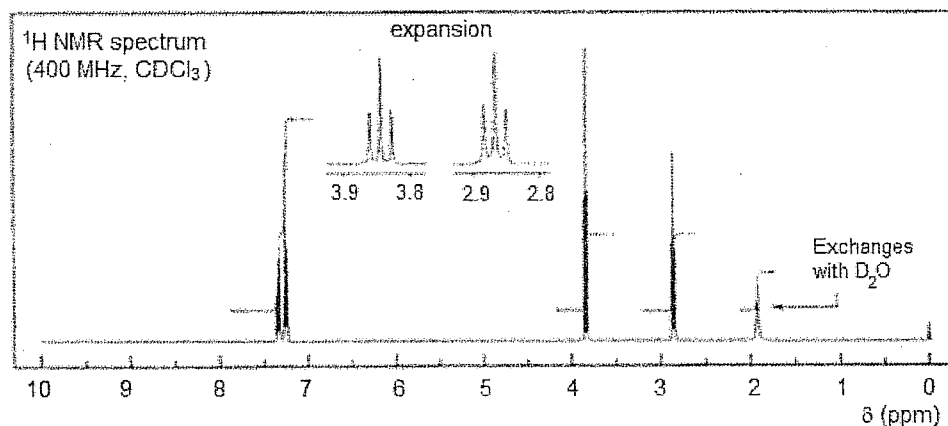
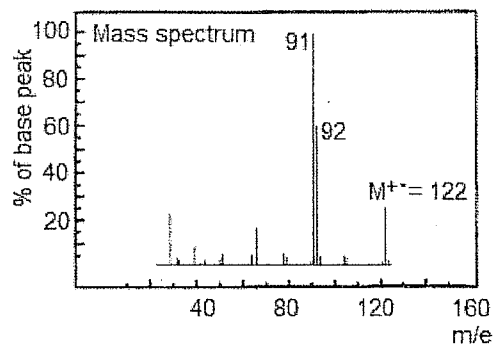
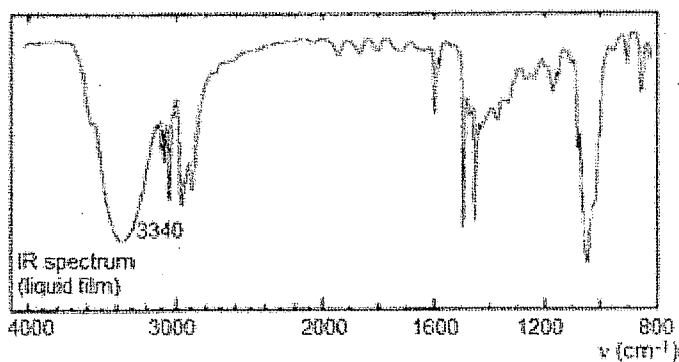
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(20 marks)

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4. **F** is an organic compound containing eight carbon atoms. IR, Mass and ^1H NMR spectra of **F** are given below.



i. What is/are the functional group/s present in **F**?

.....

ii. What are the different spin systems present in **F**?

iii. Give the structure of **F**.

iv. Draw the fragmentation pattern for the formation of the molecular fragment responsible for the peak at m/e 91.

v. Give reason why it forms the base peak in the mass spectrum.

(30 marks)

Reg. No.

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Name :.....

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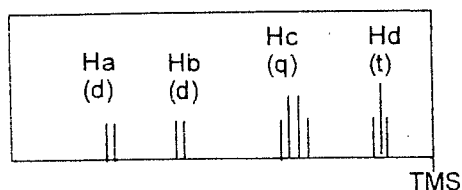
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B.Sc. Degree Programme and Stand Alone Courses in Science – 2015/2016
 CMU2221/CME4221 – Organic Chemistry I
 CAT 2- Answer Guide

1. i. Compound B : Sharp absorption band at 1683 cm^{-1} is due to $\text{C}=\text{O}$ stretching. Lower carbonyl frequency than normal value indicates a conjugated carbonyl group.
 ii. Compound D : Band at 1771 cm^{-1} due to $\text{C}=\text{O}$ str. and the band at 1180 cm^{-1} for $\text{C}-\text{O}$ str. suggests an ester group. Ester carbonyl appears at a higher frequency than for a normal acyclic one when they are in a strained ring. \therefore The compound cannot be C, but D.
 iii. Compound A : Broad band at 3330 cm^{-1} indicates a presence of a $-\text{OH}$ group. Only cpd. having $-\text{OH}$ is A.

2.



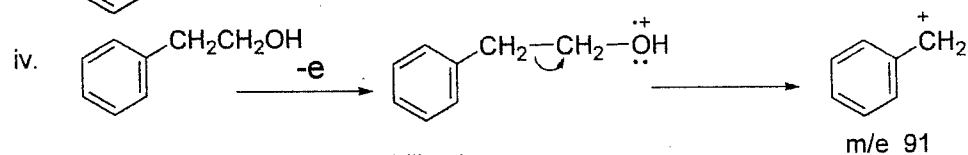
3. i. $n \longrightarrow \pi^*$ (290 nm), $\pi \longrightarrow \pi^*$ (190 nm)
 ii. $n \longrightarrow \pi^*$ electronic transitions require less energy than $\pi \longrightarrow \pi^*$ electronic transitions. When energy (E) is lesser λ will be higher as $E \propto 1/\lambda$.

4.

i. $-\text{OH}$

ii. $-\text{OH}$, $-\text{CH}_2\text{CH}_2-$,

iii.



v. Benzyl cation is resonance stabilized.

